

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020616\  
 Data File : BF084908.D  
 Acq On : 6 Feb 2016 13:35  
 Operator : SJ/IZ  
 Sample : H1310-02  
 Misc :  
 ALS Vial : 50 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SUP-B008(25-27)

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-BF020116.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         | 2.373       | 23            | 25          | 28           | rBV      | 215595         | 278045        | 4.20%           | 0.372%        |
| 2         | 2.498       | 30            | 36          | 45           | rVB3     | 14913          | 70024         | 1.06%           | 0.094%        |
| 3         | 3.447       | 115           | 119         | 122          | rBV      | 44789          | 86605         | 1.31%           | 0.116%        |
| 4         | 4.121       | 176           | 178         | 184          | rVB2     | 93218          | 150886        | 2.28%           | 0.202%        |
| 5         | 4.830       | 237           | 240         | 249          | rVB      | 1363392        | 2262443       | 34.15%          | 3.027%        |
| 6         | 5.081       | 260           | 262         | 265          | rBV      | 1547309        | 1879040       | 28.37%          | 2.514%        |
| 7         | 5.207       | 270           | 273         | 276          | rBV      | 4127567        | 6624378       | 100.00%         | 8.864%        |
| 8         | 5.264       | 276           | 278         | 281          | rBV      | 4041037        | 4329523       | 65.36%          | 5.793%        |
| 9         | 5.470       | 294           | 296         | 299          | rBV      | 920705         | 1237018       | 18.67%          | 1.655%        |
| 10        | 5.664       | 311           | 313         | 316          | rVB      | 69149          | 69520         | 1.05%           | 0.093%        |
| 11        | 5.813       | 324           | 326         | 328          | rVB      | 204954         | 169975        | 2.57%           | 0.227%        |
| 12        | 6.110       | 350           | 352         | 355          | rVB      | 563012         | 473356        | 7.15%           | 0.633%        |
| 13        | 6.179       | 356           | 358         | 360          | rBV      | 1611636        | 1484501       | 22.41%          | 1.986%        |
| 14        | 6.213       | 360           | 361         | 363          | rVB      | 1124269        | 808216        | 12.20%          | 1.081%        |
| 15        | 6.259       | 363           | 365         | 368          | rVB      | 1506095        | 1198762       | 18.10%          | 1.604%        |
| 16        | 6.316       | 368           | 370         | 374          | rBV      | 3254458        | 4901790       | 74.00%          | 6.559%        |
| 17        | 6.430       | 378           | 380         | 383          | rBV      | 3894449        | 4816065       | 72.70%          | 6.444%        |
| 18        | 6.487       | 383           | 385         | 388          | rVB      | 3997979        | 3526127       | 53.23%          | 4.718%        |
| 19        | 6.659       | 398           | 400         | 404          | rBV      | 1265727        | 1140450       | 17.22%          | 1.526%        |
| 20        | 6.716       | 404           | 405         | 408          | rVB      | 1272246        | 1359264       | 20.52%          | 1.819%        |
| 21        | 6.819       | 412           | 414         | 415          | rBV      | 4486864        | 3973224       | 59.98%          | 5.316%        |
| 22        | 6.922       | 420           | 423         | 424          | rBV      | 211274         | 251072        | 3.79%           | 0.336%        |
| 23        | 6.944       | 424           | 425         | 427          | rVB      | 245062         | 175840        | 2.65%           | 0.235%        |
| 24        | 6.990       | 427           | 429         | 431          | rBV      | 694334         | 634205        | 9.57%           | 0.849%        |
| 25        | 7.059       | 433           | 435         | 438          | rBV      | 141501         | 135047        | 2.04%           | 0.181%        |
| 26        | 7.139       | 440           | 442         | 445          | rVB      | 278329         | 533894        | 8.06%           | 0.714%        |
| 27        | 7.230       | 445           | 450         | 453          | rBV2     | 2898302        | 4312711       | 65.10%          | 5.771%        |
| 28        | 7.356       | 459           | 461         | 463          | rBV      | 158008         | 171840        | 2.59%           | 0.230%        |
| 29        | 7.436       | 466           | 468         | 469          | rBV      | 341014         | 333487        | 5.03%           | 0.446%        |
| 30        | 7.470       | 469           | 471         | 472          | rVB      | 401767         | 443488        | 6.69%           | 0.593%        |
| 31        | 7.619       | 482           | 484         | 487          | rVB      | 531652         | 547594        | 8.27%           | 0.733%        |
| 32        | 7.687       | 487           | 490         | 492          | rBV      | 1032200        | 1030452       | 15.56%          | 1.379%        |
| 33        | 7.733       | 492           | 494         | 496          | rVV2     | 109727         | 205577        | 3.10%           | 0.275%        |
| 34        | 7.790       | 496           | 499         | 500          | rVB3     | 106280         | 148400        | 2.24%           | 0.199%        |

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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-BF020116.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 7.973  | 510  | 515  | 516  | rBV2 | 4224885 | 5555803 | 83.87% | 7.434% |
| 36 | 8.019  | 516  | 519  | 520  | rVB2 | 227908  | 250655  | 3.78%  | 0.335% |
| 37 | 8.225  | 533  | 537  | 541  | rBV3 | 67604   | 109050  | 1.65%  | 0.146% |
| 38 | 8.339  | 545  | 547  | 549  | rVV  | 76368   | 73661   | 1.11%  | 0.099% |
| 39 | 8.419  | 549  | 554  | 555  | rVV  | 80964   | 100564  | 1.52%  | 0.135% |
| 40 | 8.659  | 573  | 575  | 577  | rVV  | 1008198 | 814793  | 12.30% | 1.090% |
| 41 | 8.762  | 581  | 584  | 586  | rVB  | 262409  | 377381  | 5.70%  | 0.505% |
| 42 | 9.025  | 604  | 607  | 610  | rBV  | 4761460 | 5322542 | 80.35% | 7.122% |
| 43 | 9.699  | 663  | 666  | 667  | rBV  | 1239273 | 1447007 | 21.84% | 1.936% |
| 44 | 9.722  | 667  | 668  | 671  | rVB  | 115384  | 106204  | 1.60%  | 0.142% |
| 45 | 10.488 | 732  | 735  | 737  | rBV  | 3037439 | 3377011 | 50.98% | 4.519% |
| 46 | 11.173 | 792  | 795  | 797  | rBV  | 1564836 | 1374137 | 20.74% | 1.839% |
| 47 | 12.602 | 917  | 920  | 922  | rBV  | 98533   | 111301  | 1.68%  | 0.149% |
| 48 | 12.762 | 931  | 934  | 937  | rBV  | 4255831 | 4054231 | 61.20% | 5.425% |
| 49 | 13.299 | 979  | 981  | 983  | rBV  | 78841   | 71902   | 1.09%  | 0.096% |
| 50 | 13.802 | 1023 | 1025 | 1027 | rBV  | 1040417 | 930725  | 14.05% | 1.245% |
| 51 | 15.174 | 1142 | 1145 | 1147 | rBV  | 774453  | 896439  | 13.53% | 1.199% |

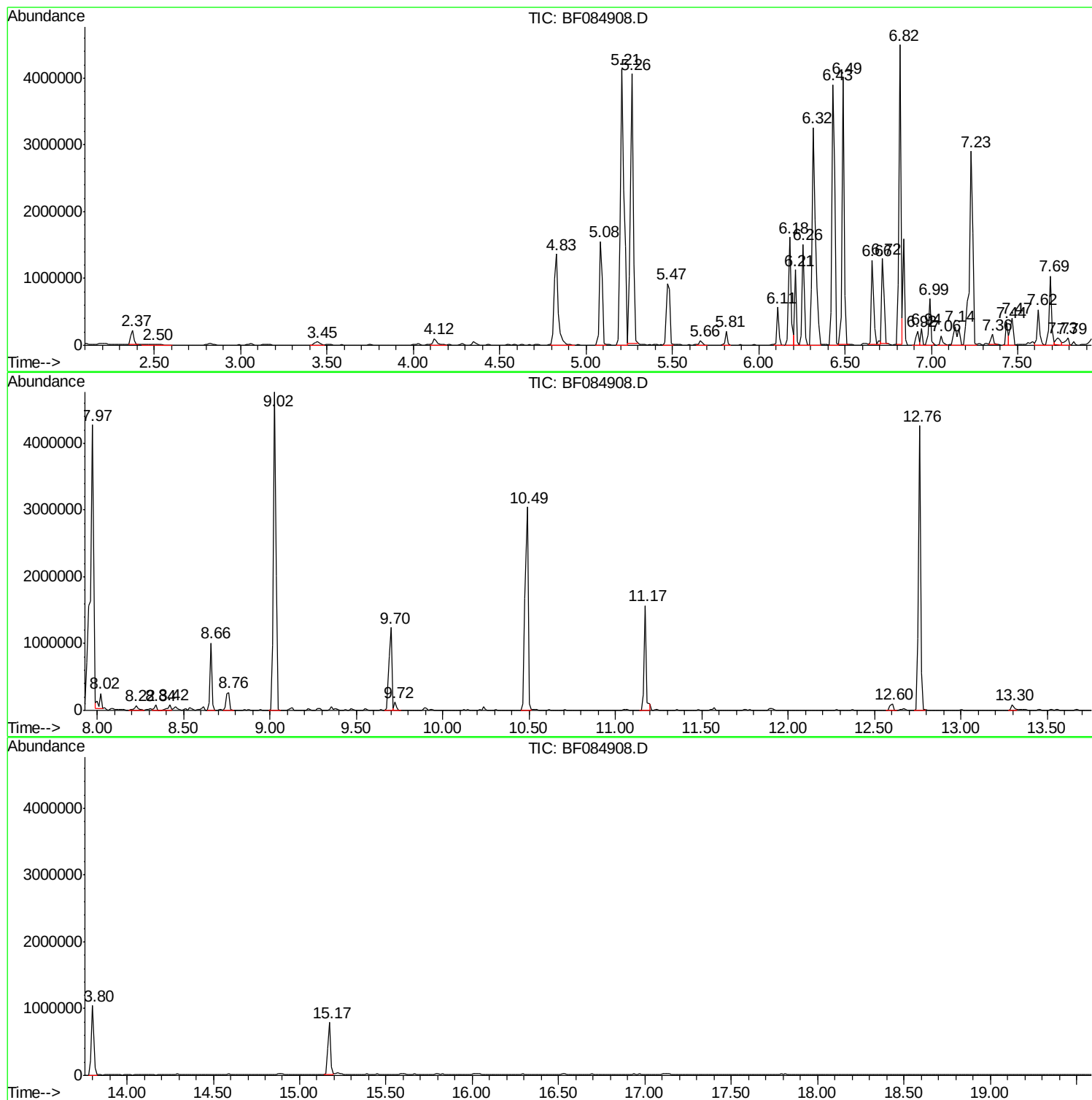
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ClientSampled :  
SUP-B008(25-27)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.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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Instrument :  
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ClientSampleId :  
SUP-B008(25-27)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

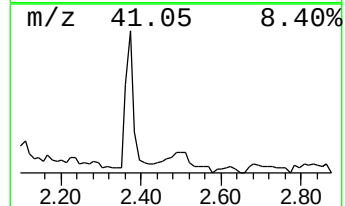
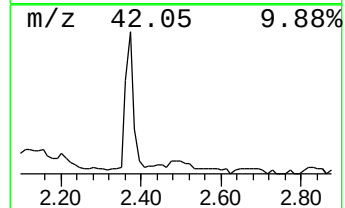
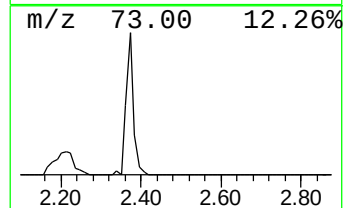
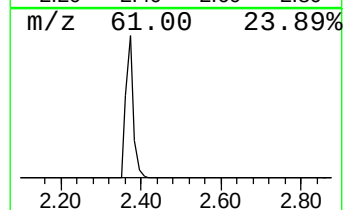
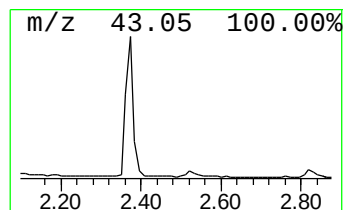
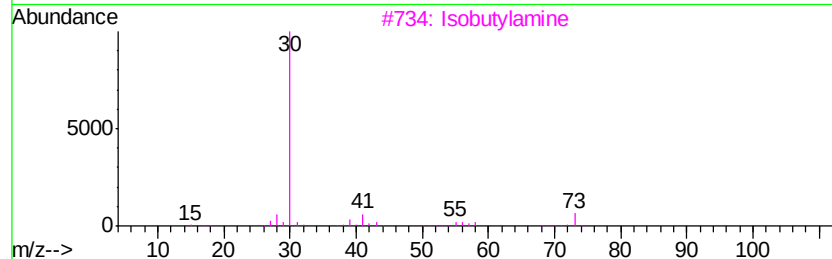
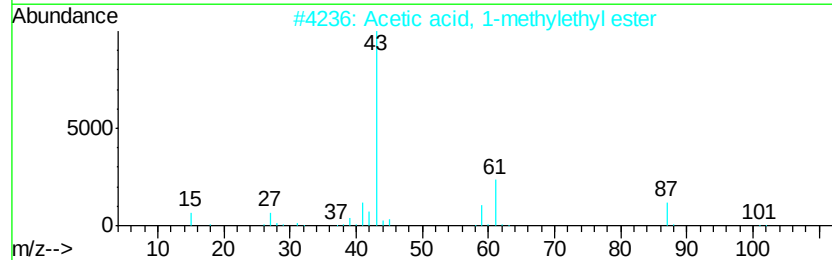
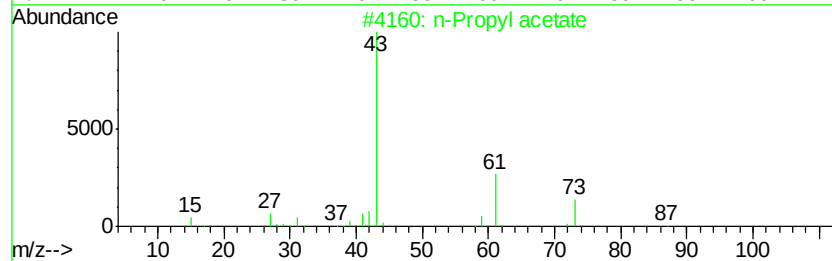
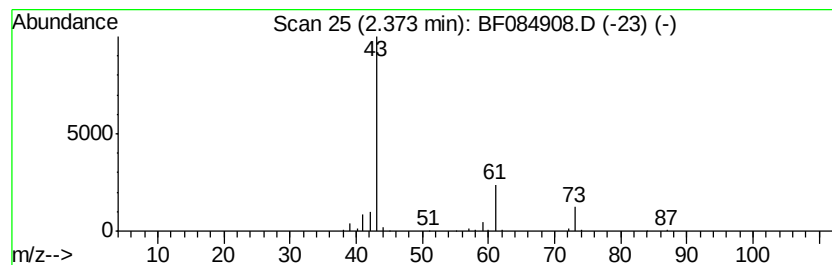
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 n-Propyl acetate Concentration Rank 15

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 2.37 | 4.88 ng | 278045 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | n-Propyl acetate                 | 102 | C5H10O2 | 000109-60-4 | 83   |
| 2    |    | Acetic acid, 1-methylethyl ester | 102 | C5H10O2 | 000108-21-4 | 36   |
| 3    |    | Isobutylamine                    | 73  | C4H11N  | 000078-81-9 | 5    |
| 4    |    | 1-Butanamine                     | 73  | C4H11N  | 000109-73-9 | 5    |
| 5    |    | Pentane                          | 72  | C5H12   | 000109-66-0 | 4    |



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ClientSampleID :  
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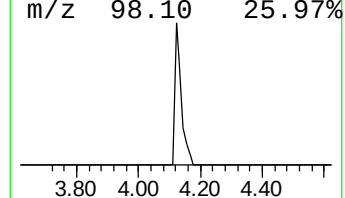
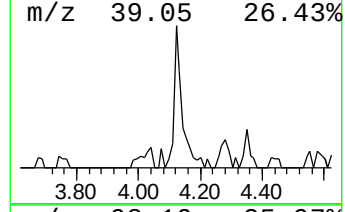
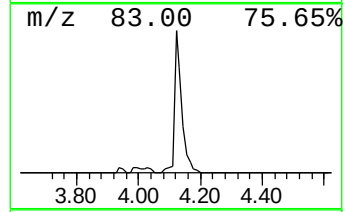
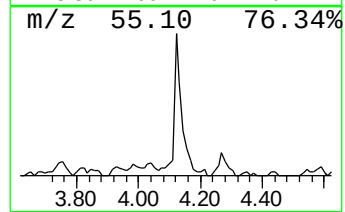
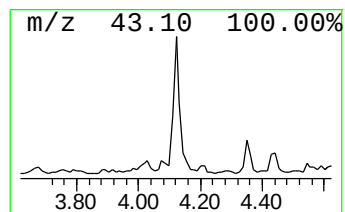
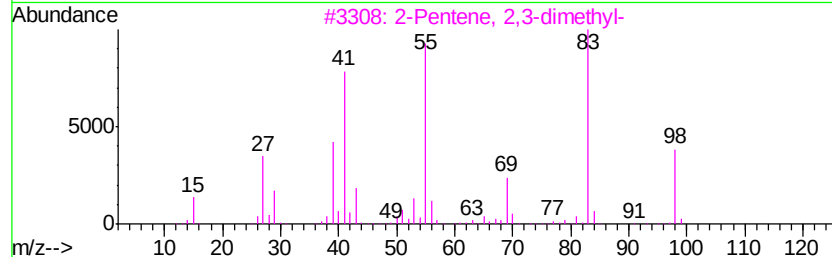
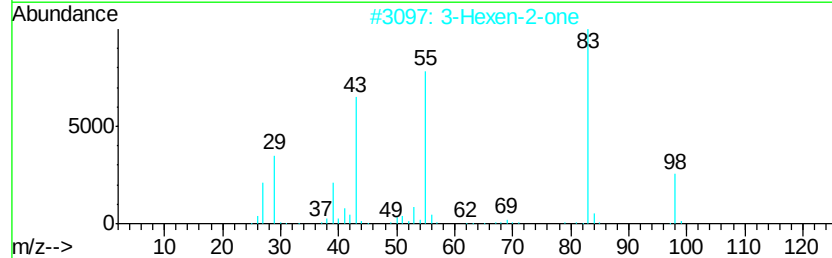
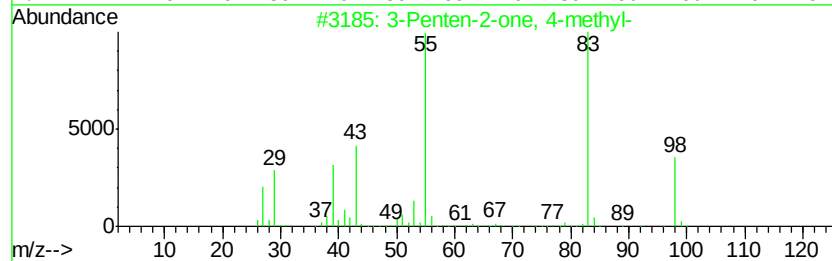
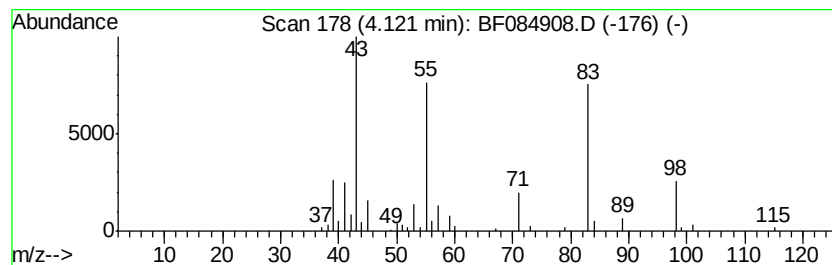
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 20

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.12 | 2.65 ng | 150886 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | 5 | Tentative ID                   | MW | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|----|---------|-------------|------|
| 1    |    |   | 3-Penten-2-one, 4-methyl-      | 98 | C6H10O  | 000141-79-7 | 76   |
| 2    |    |   | 3-Hexen-2-one                  | 98 | C6H10O  | 000763-93-9 | 68   |
| 3    |    |   | 2-Pentene, 2,3-dimethyl-       | 98 | C7H14   | 010574-37-5 | 64   |
| 4    |    |   | 2-Pentene, 3,4-dimethyl-, (E)- | 98 | C7H14   | 004914-92-5 | 58   |
| 5    |    |   | 2-Pentene, 3,4-dimethyl-, (Z)- | 98 | C7H14   | 004914-91-4 | 58   |



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ALS Vial : 50 Sample Multiplier: 1

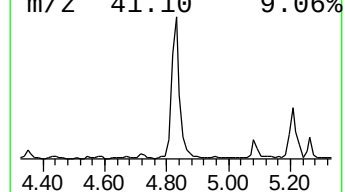
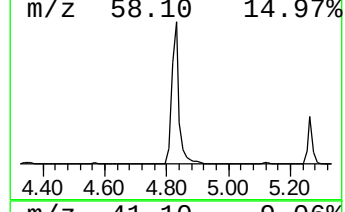
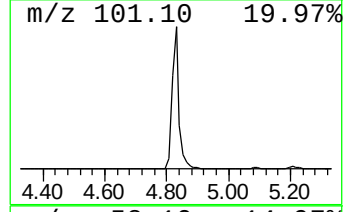
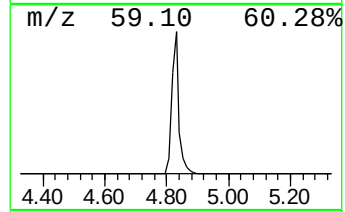
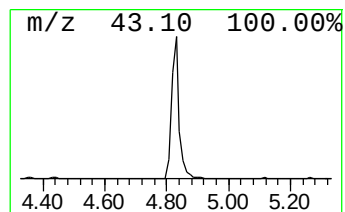
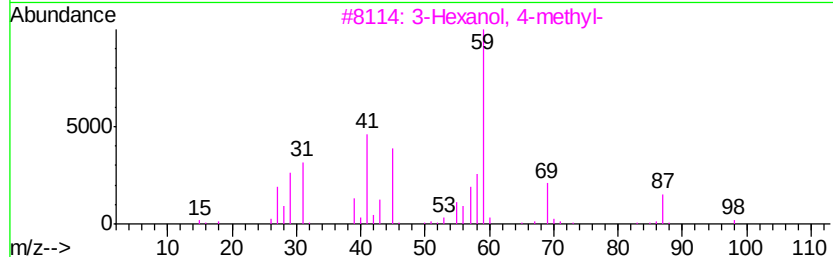
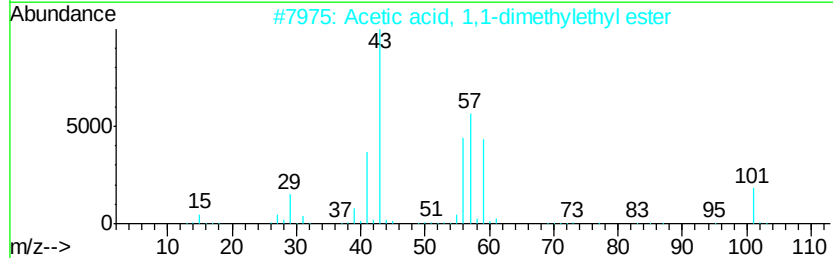
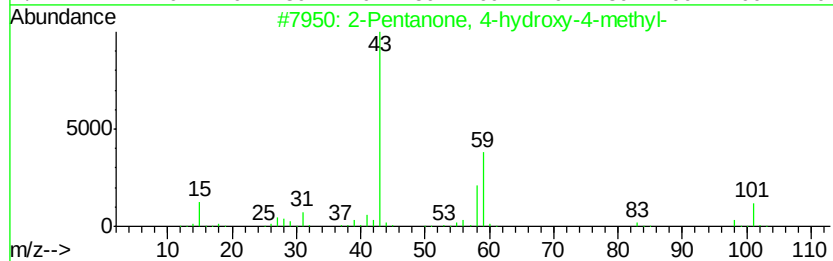
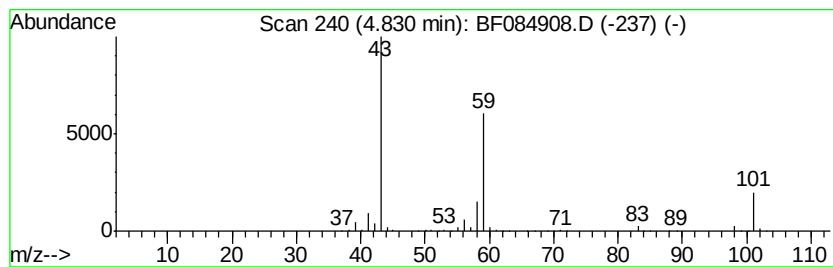
Instrument :  
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ClientSampleId :  
SUP-B008(25-27)

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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 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

| R.T.    | EstConc                             | Area         | Relative to ISTD       | R.T.        |      |      |
|---------|-------------------------------------|--------------|------------------------|-------------|------|------|
| 4.83    | 39.68 ng                            | 2262440      | 1,4-Dichlorobenzene-d4 | 6.66        |      |      |
| Hit# of | 5                                   | Tentative ID | MW                     | MolForm     | CAS# | Qual |
| 1       | 2-Pentanone, 4-hydroxy-4-methyl-    | 116          | C6H12O2                | 000123-42-2 | 56   |      |
| 2       | Acetic acid, 1,1-dimethylethyl e... | 116          | C6H12O2                | 000540-88-5 | 39   |      |
| 3       | 3-Hexanol, 4-methyl-                | 116          | C7H16O                 | 000615-29-2 | 28   |      |
| 4       | Acetic acid, cyano-, 1,1-dimethy... | 141          | C7H11NO2               | 001116-98-9 | 17   |      |
| 5       | (+)-4-Amino-4,5-dihydro-2(3H)-f...  | 101          | C4H7NO2                | 016504-58-8 | 9    |      |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

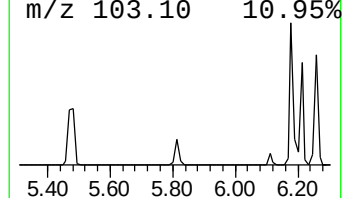
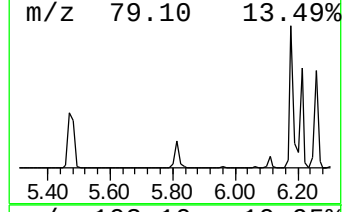
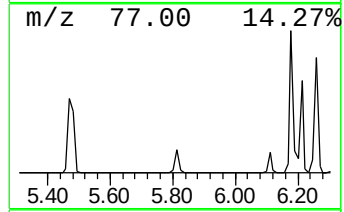
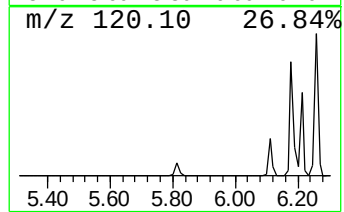
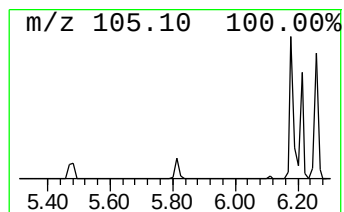
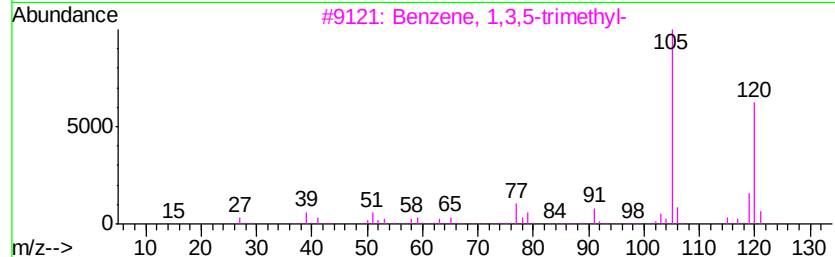
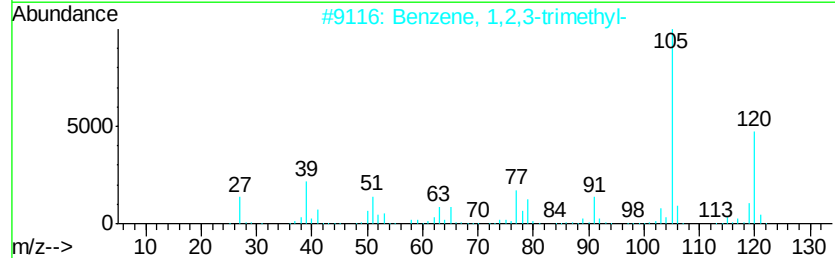
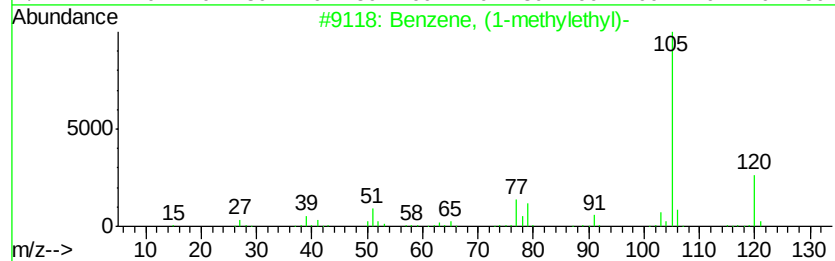
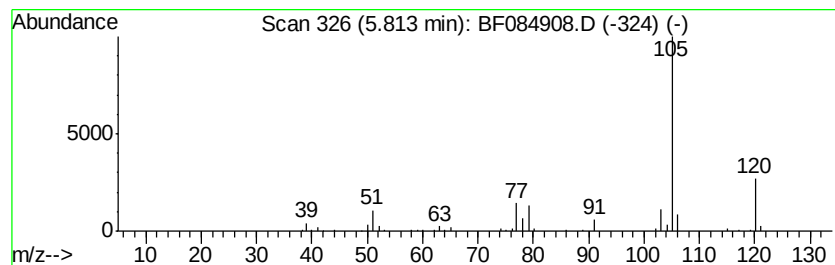
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Benzene, (1-methylethyl)- Concentration Rank 19

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.81 | 2.98 ng | 169975 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (1-methylethyl)-  | 120 | C9H12   | 000098-82-8 | 94   |
| 2    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 3    |    | Benzene, 1,3,5-trimethyl-  | 120 | C9H12   | 000108-67-8 | 78   |
| 4    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 72   |
| 5    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 72   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020616\  
Data File : BF084908.D  
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Operator : SJ/IZ  
Sample : H1310-02  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

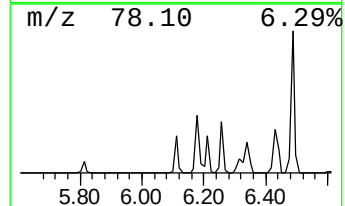
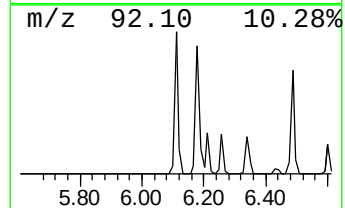
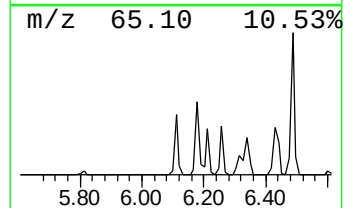
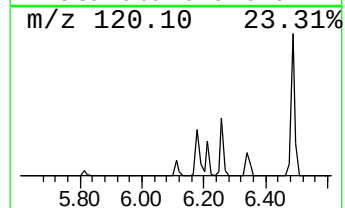
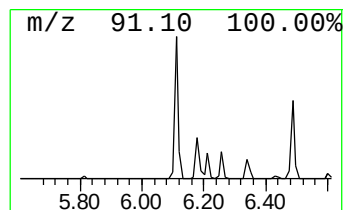
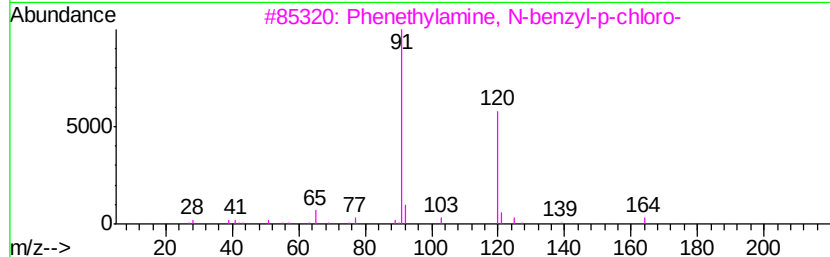
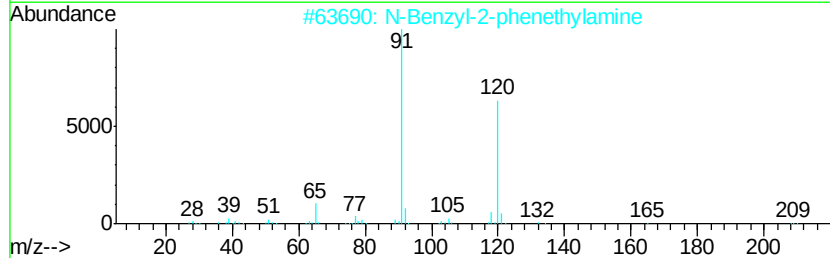
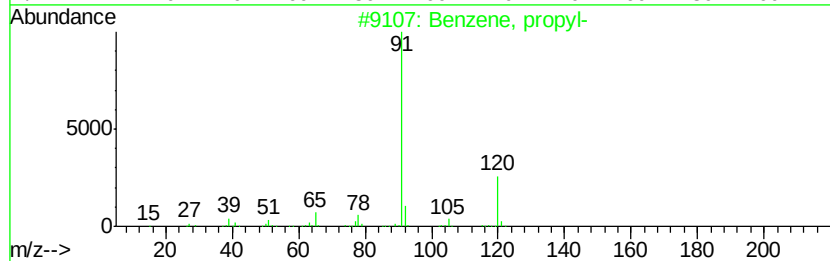
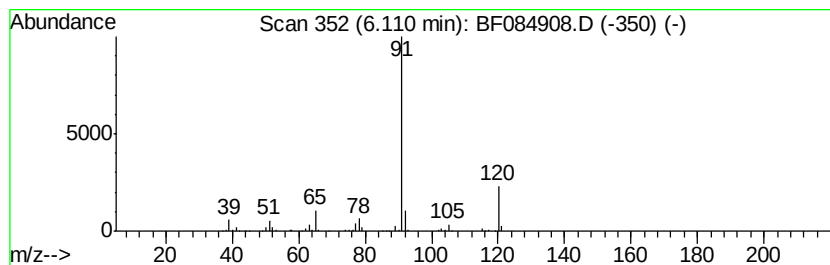
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SUP-B008(25-27)

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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 8 Benzene, propyl- Concentration Rank 14

| R.T.      | EstConc                              | Area   | Relative to ISTD       | R.T.        |      |
|-----------|--------------------------------------|--------|------------------------|-------------|------|
| 6.11      | 8.30 ng                              | 473356 | 1,4-Dichlorobenzene-d4 | 6.66        |      |
| Hit# of 5 | Tentative ID                         | MW     | MolForm                | CAS#        | Qual |
| 1         | Benzene, propyl-                     | 120    | C9H12                  | 000103-65-1 | 91   |
| 2         | N-Benzyl-2-phenethylamine            | 211    | C15H17N                | 003647-71-0 | 83   |
| 3         | Phenethylamine, N-benzyl-p-chloro-   | 245    | C15H16ClN              | 013622-43-0 | 72   |
| 4         | Propanenitrile, 3-[(phenylmethyl)... | 160    | C10H12N2               | 000706-03-6 | 72   |
| 5         | 1,2-Ethanediamine, N,N'-bis(phen...  | 240    | C16H20N2               | 000140-28-3 | 64   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020616\  
Data File : BF084908.D  
Acq On : 6 Feb 2016 13:35  
Operator : SJ/IZ  
Sample : H1310-02  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SUP-B008(25-27)

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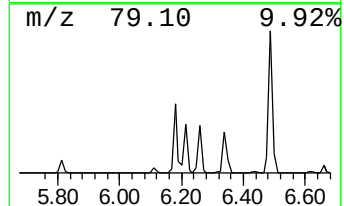
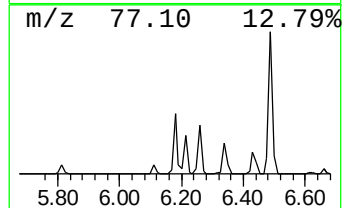
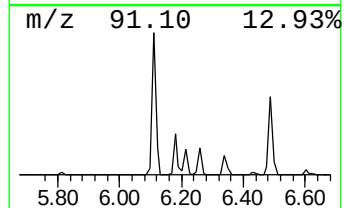
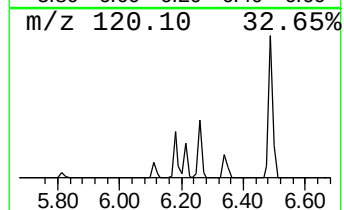
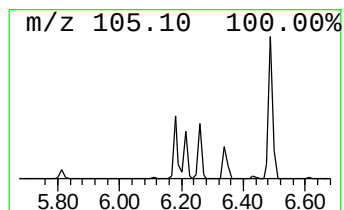
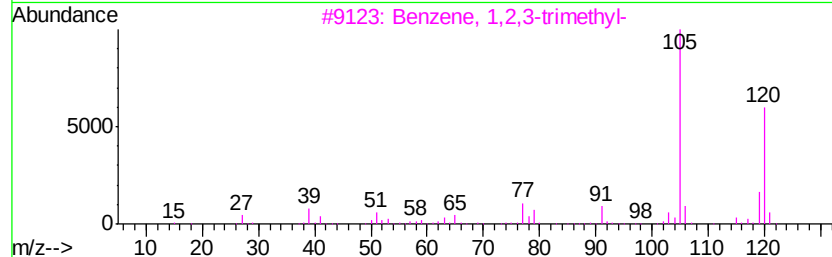
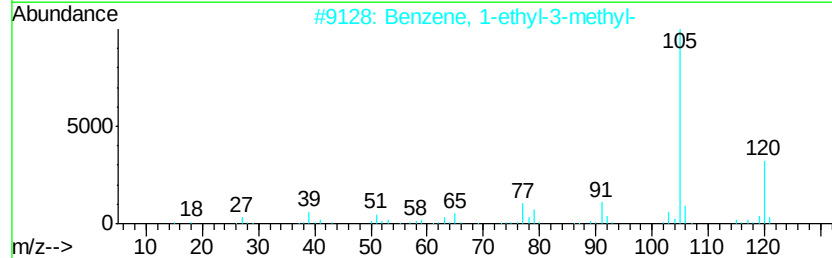
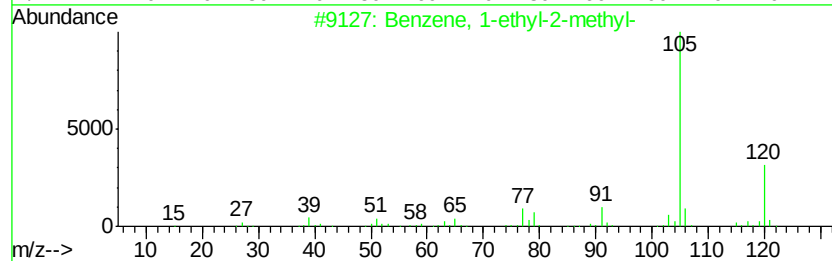
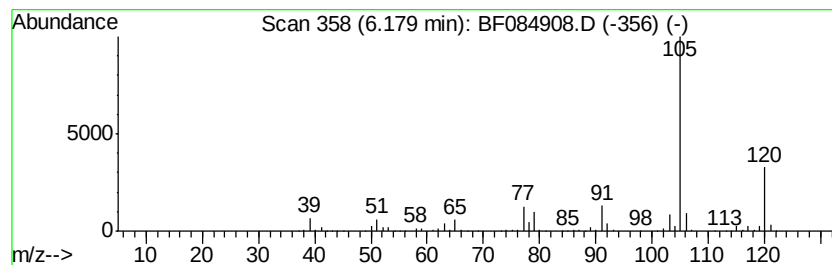
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.18 | 26.03 ng | 1484500 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 2    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 95   |
| 3    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 4    |    | Benzene, 1,3,5-trimethyl-  | 120 | C9H12   | 000108-67-8 | 91   |
| 5    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020616\  
Data File : BF084908.D  
Acq On : 6 Feb 2016 13:35  
Operator : SJ/IZ  
Sample : H1310-02  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SUP-B008(25-27)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

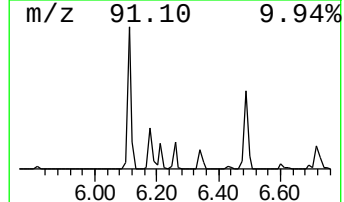
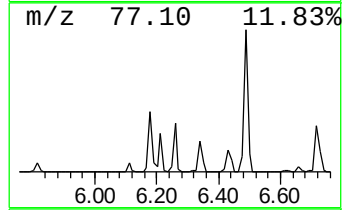
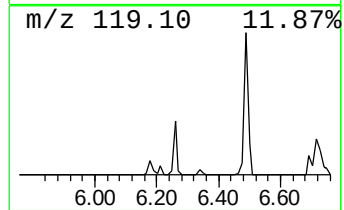
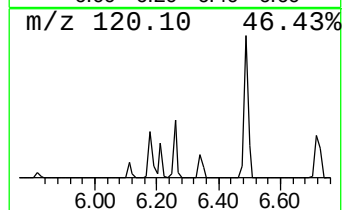
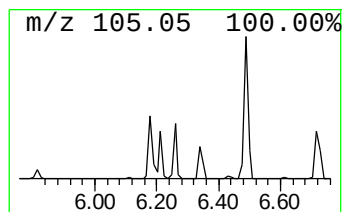
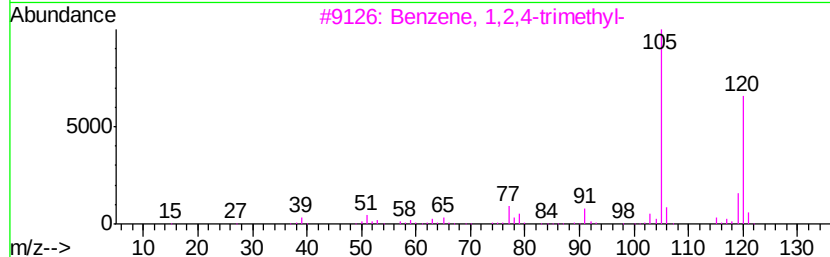
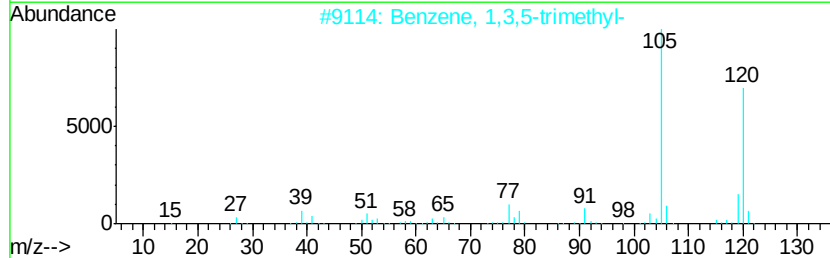
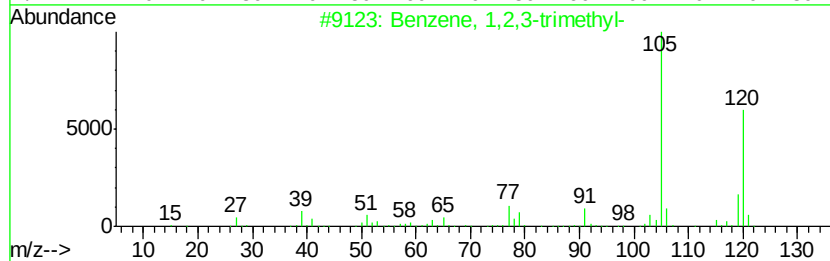
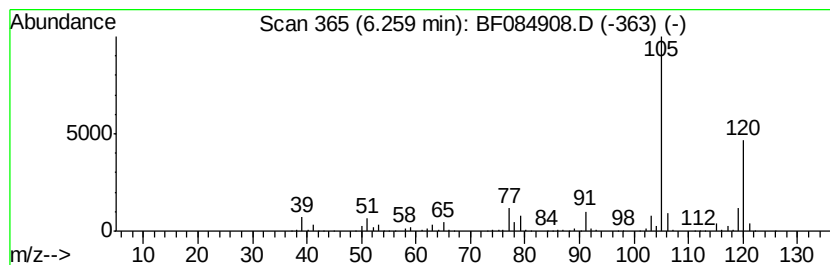
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 10

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.26 | 21.02 ng | 1198760 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    | Benzene, 1,3,5-trimethyl-  | 120 | C9H12   | 000108-67-8 | 94   |
| 3    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 94   |
| 4    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 93   |
| 5    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020616\  
Data File : BF084908.D  
Acq On : 6 Feb 2016 13:35  
Operator : SJ/IZ  
Sample : H1310-02  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SUP-B008(25-27)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

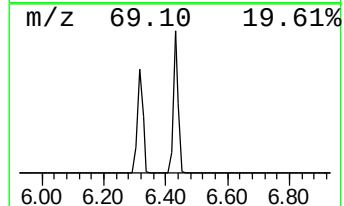
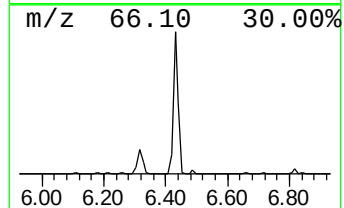
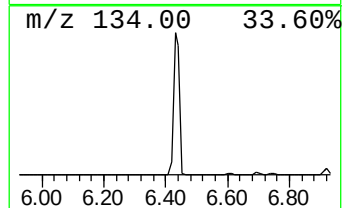
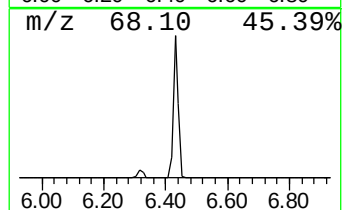
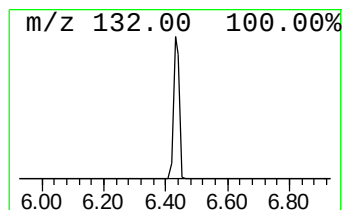
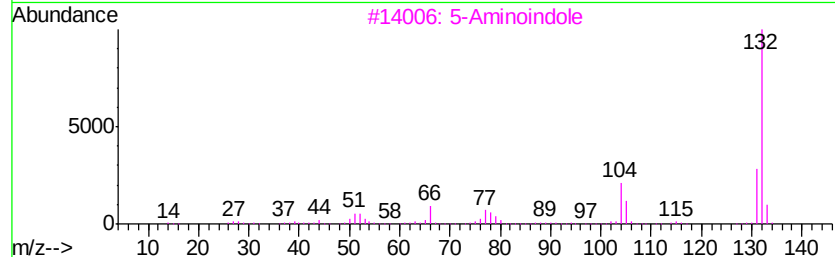
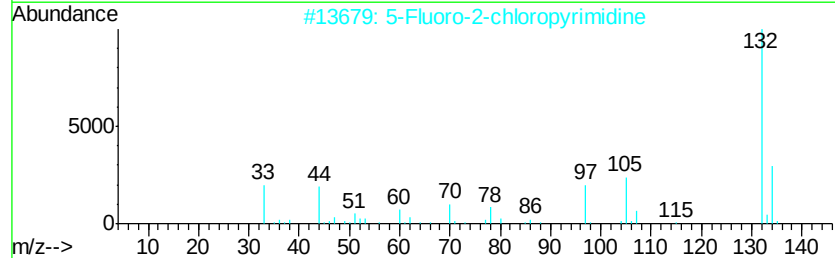
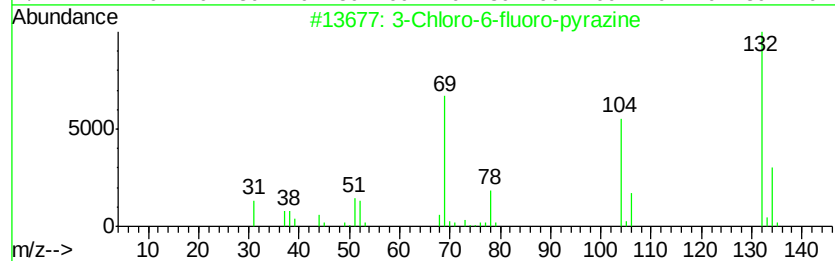
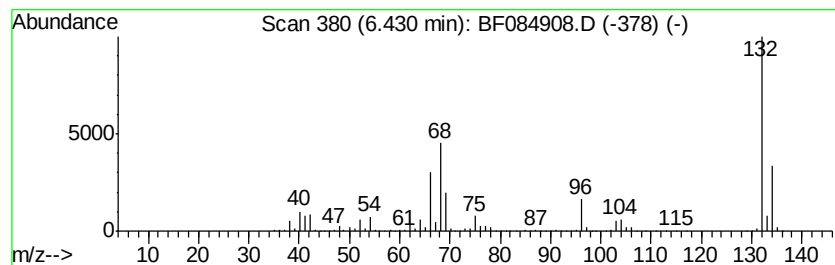
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 unknown6.43 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.43 | 84.46 ng | 4816070 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 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    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2    | 022291-85-6  | 9    |



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Data File : BF084908.D  
Acq On : 6 Feb 2016 13:35  
Operator : SJ/IZ  
Sample : H1310-02  
Misc :  
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Instrument :  
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ClientSampleId :  
SUP-B008(25-27)

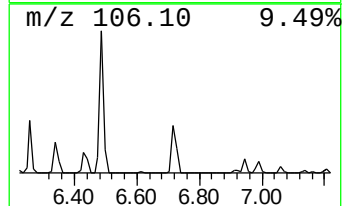
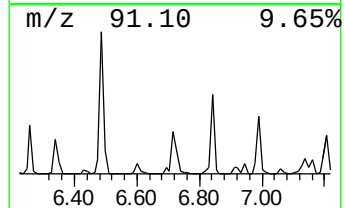
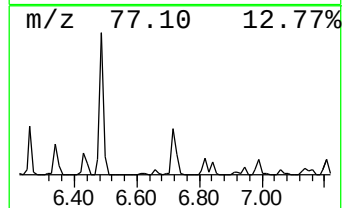
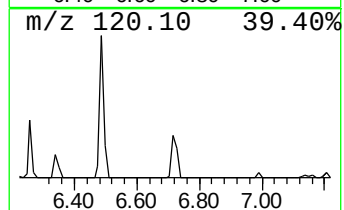
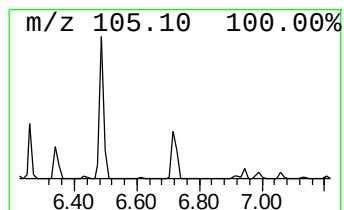
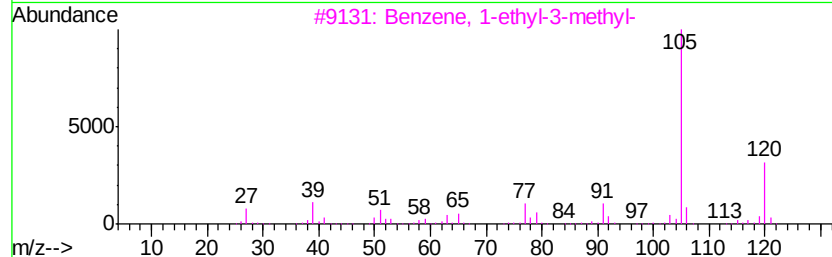
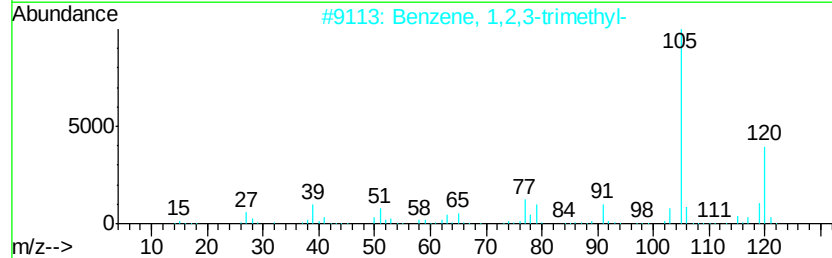
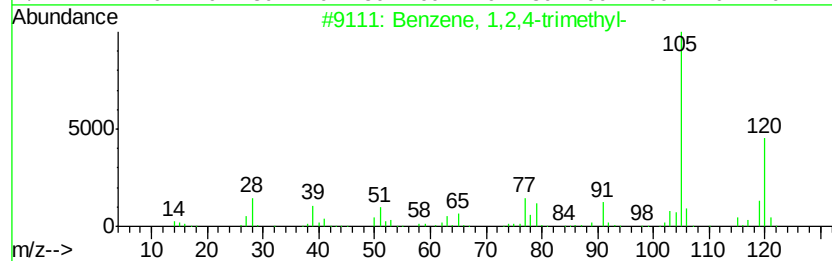
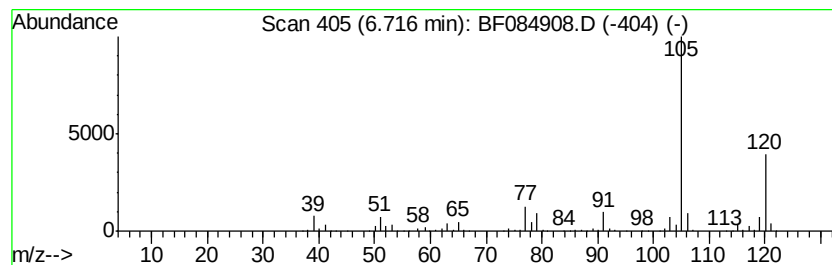
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 14 Benzene, 1,2,4-trimethyl- Concentration Rank 8

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.72 | 23.84 ng | 1359260 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 94   |
| 2    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 3    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 4    |    | Benzene, 1,3,5-trimethyl-  | 120 | C9H12   | 000108-67-8 | 91   |
| 5    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 91   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020616\  
Data File : BF084908.D  
Acq On : 6 Feb 2016 13:35  
Operator : SJ/IZ  
Sample : H1310-02  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SUP-B008(25-27)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

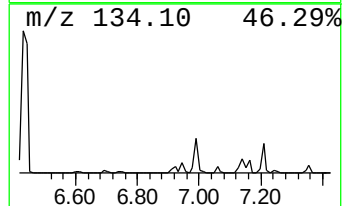
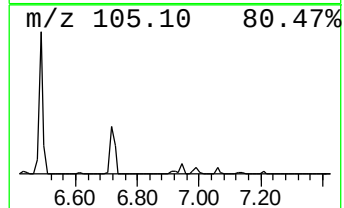
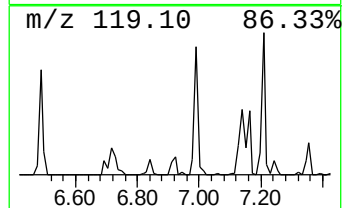
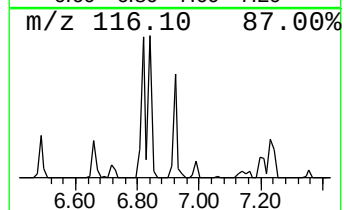
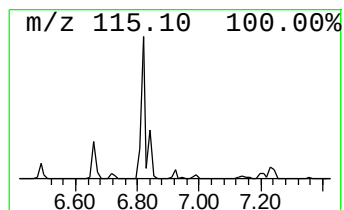
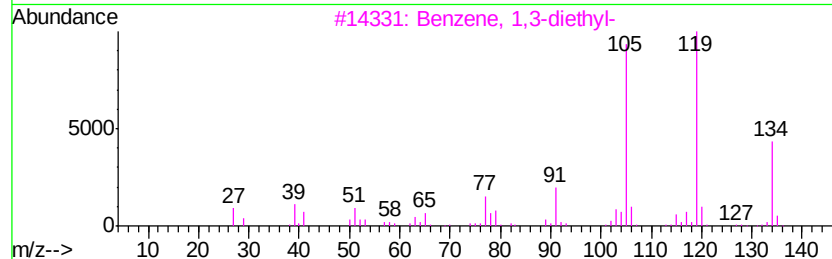
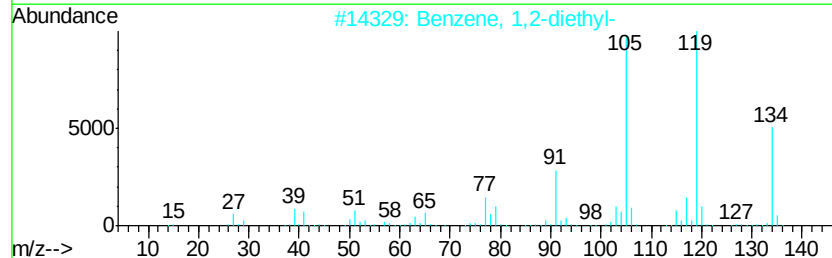
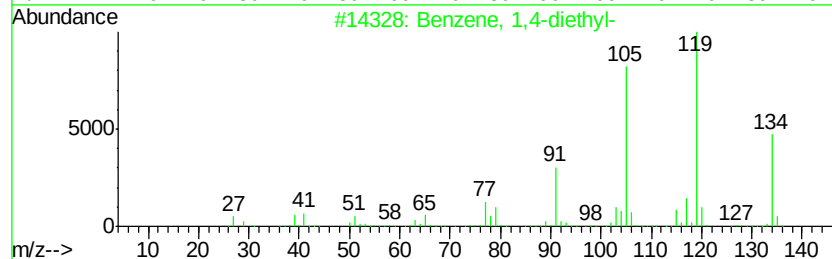
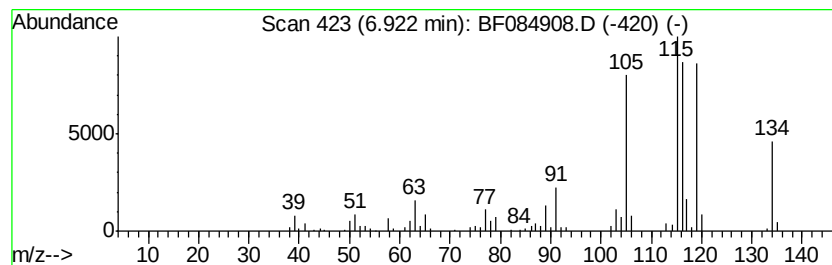
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzene, 1,4-diethyl- Concentration Rank 16

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.92 | 4.40 ng | 251072 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,4-diethyl- | 134 | C10H14  | 000105-05-5 | 90   |
| 2    |    |   | Benzene, 1,2-diethyl- | 134 | C10H14  | 000135-01-3 | 90   |
| 3    |    |   | Benzene, 1,3-diethyl- | 134 | C10H14  | 000141-93-5 | 86   |
| 4    |    |   | Indene                | 116 | C9H8    | 000095-13-6 | 86   |
| 5    |    |   | Benzene, 1-propynyl-  | 116 | C9H8    | 000673-32-5 | 50   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020616\  
Data File : BF084908.D  
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Sample : H1310-02  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SUP-B008(25-27)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

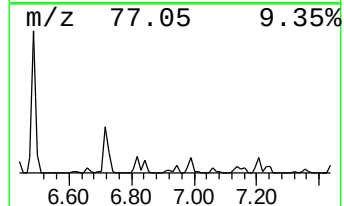
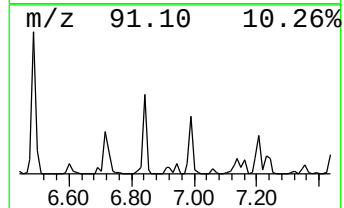
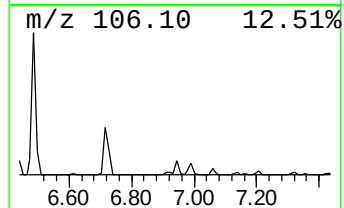
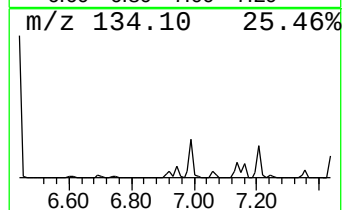
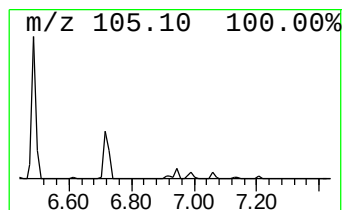
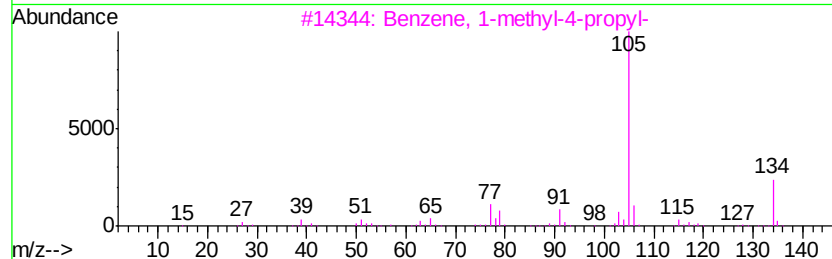
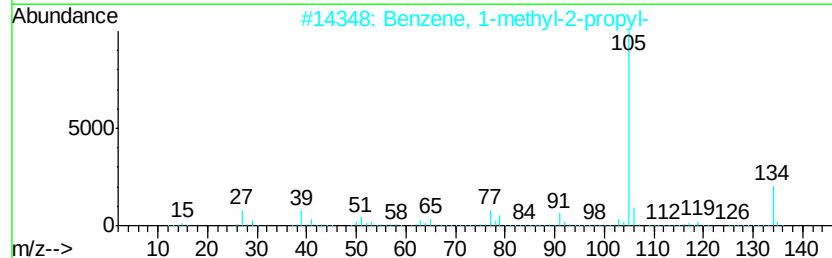
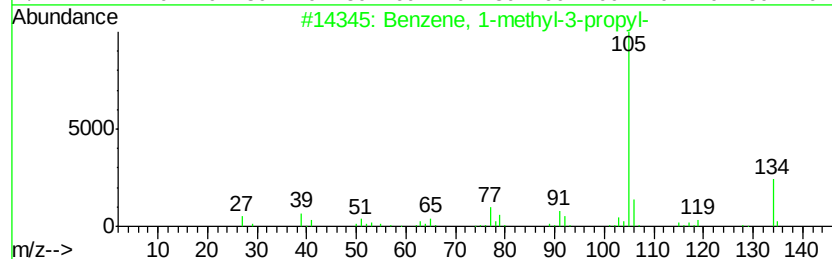
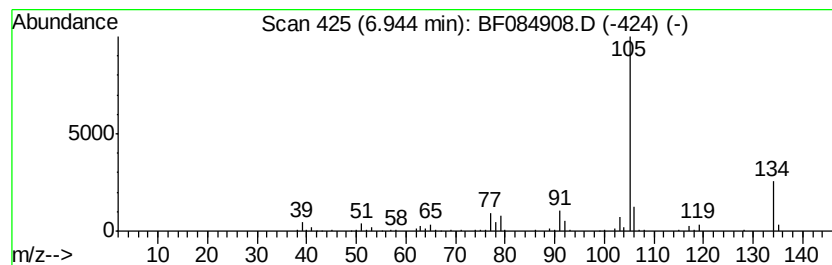
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Benzene, 1-methyl-3-propyl- Concentration Rank 18

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.94 | 3.08 ng | 175840 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-3-propyl-         | 134 | C10H14  | 001074-43-7 | 95   |
| 2    |    | Benzene, 1-methyl-2-propyl-         | 134 | C10H14  | 001074-17-5 | 91   |
| 3    |    | Benzene, 1-methyl-4-propyl-         | 134 | C10H14  | 001074-55-1 | 91   |
| 4    |    | 2-Tolyloxirane                      | 134 | C9H10O  | 002783-26-8 | 87   |
| 5    |    | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O  | 000093-53-8 | 81   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020616\  
Data File : BF084908.D  
Acq On : 6 Feb 2016 13:35  
Operator : SJ/IZ  
Sample : H1310-02  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
SUP-B008(25-27)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

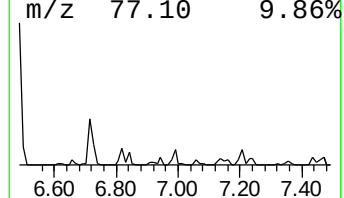
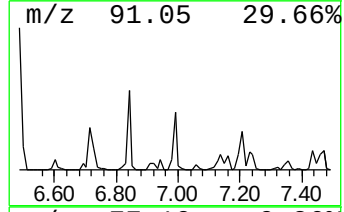
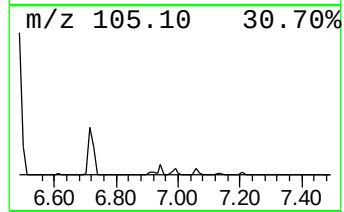
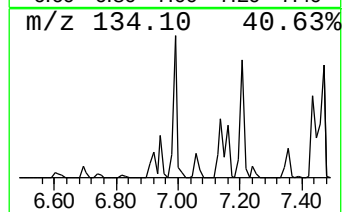
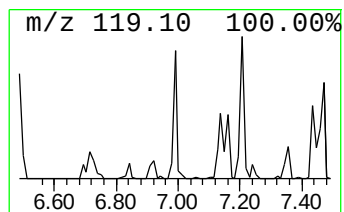
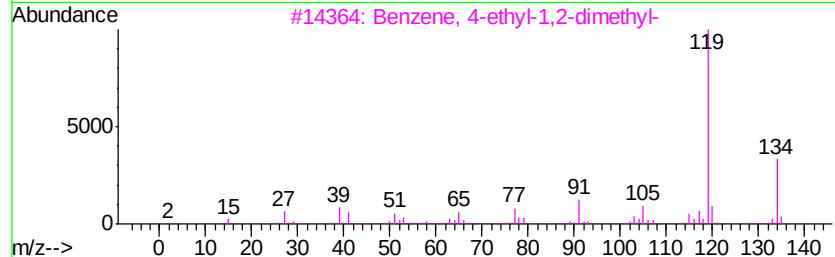
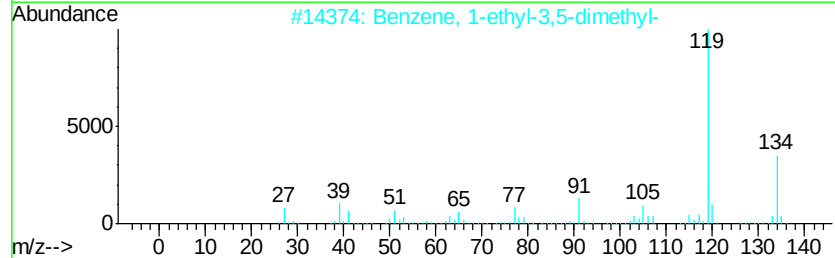
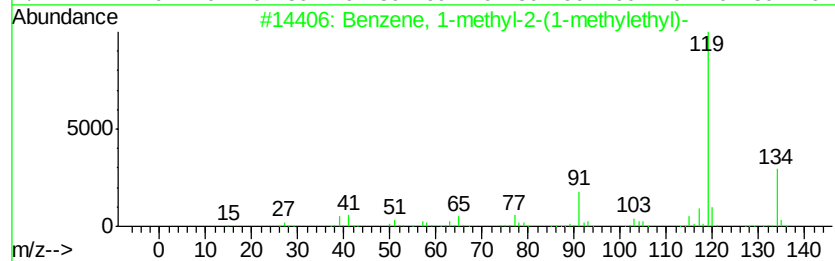
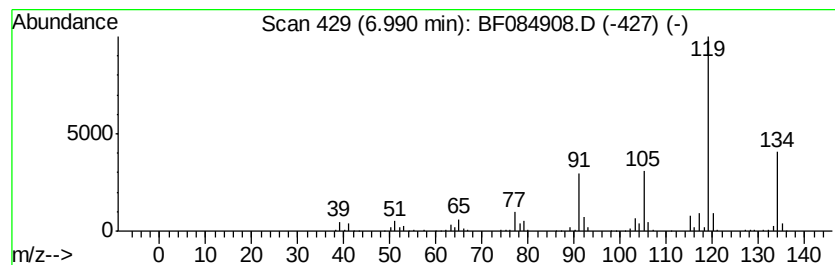
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Benzene, 1-methyl-2-(1-meth... Concentration Rank 12

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 6.99 | 11.12 ng | 634205 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 93   |
| 2    |    | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 90   |
| 3    |    | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 90   |
| 4    |    | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 90   |
| 5    |    | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 90   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020616\  
Data File : BF084908.D  
Acq On : 6 Feb 2016 13:35  
Operator : SJ/IZ  
Sample : H1310-02  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SUP-B008(25-27)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

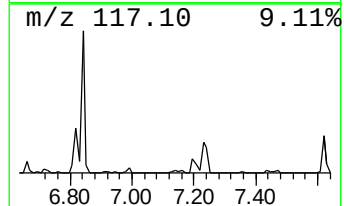
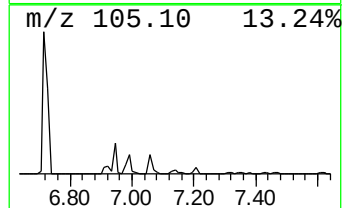
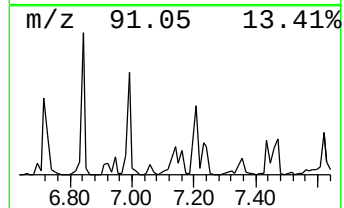
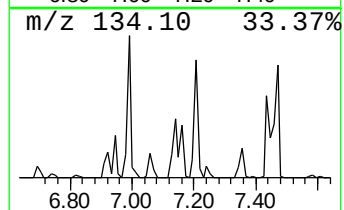
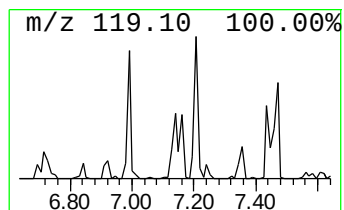
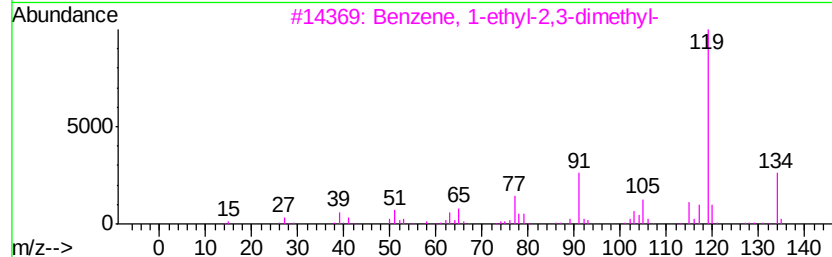
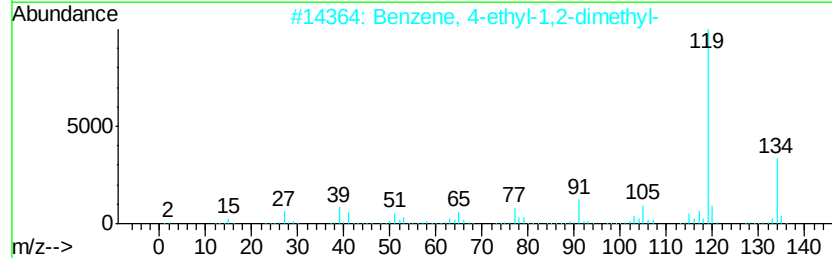
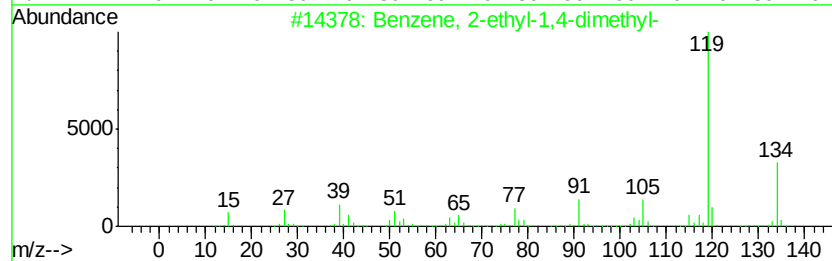
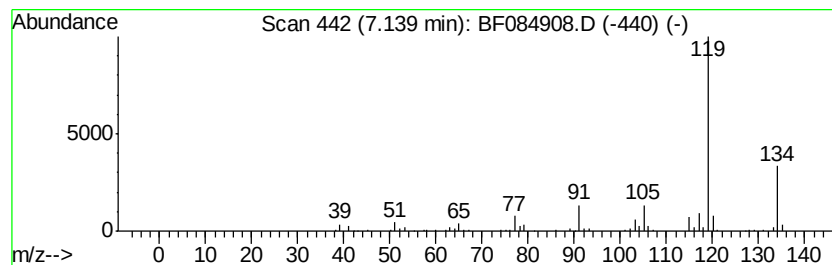
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 13

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 7.14 | 9.36 ng | 533894 | 1,4-Dichlorobenzene-d4 | 6.66 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 97   |
| 2    |    | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 95   |
| 3    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 95   |
| 4    |    | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 94   |
| 5    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 93   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020616\  
Data File : BF084908.D  
Acq On : 6 Feb 2016 13:35  
Operator : SJ/IZ  
Sample : H1310-02  
Misc :  
ALS Vial : 50 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SUP-B008(25-27)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

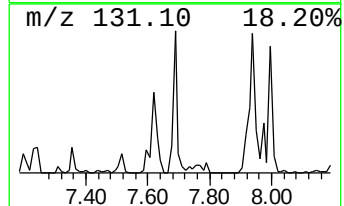
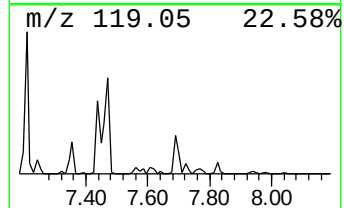
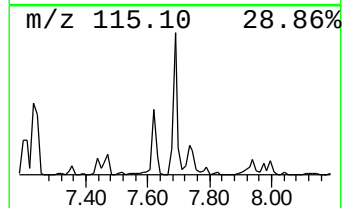
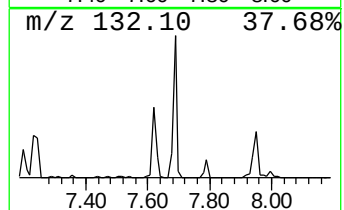
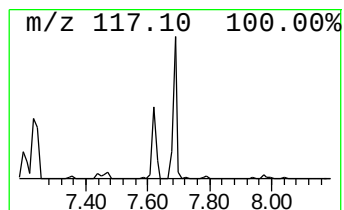
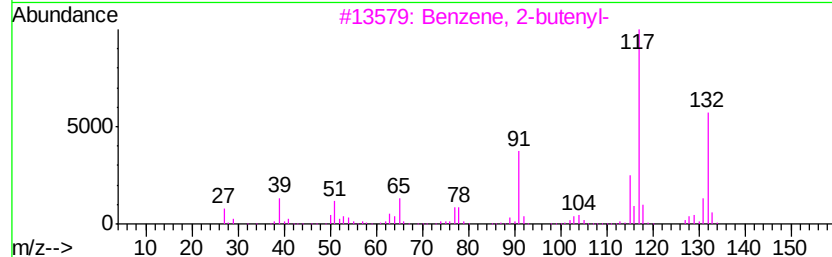
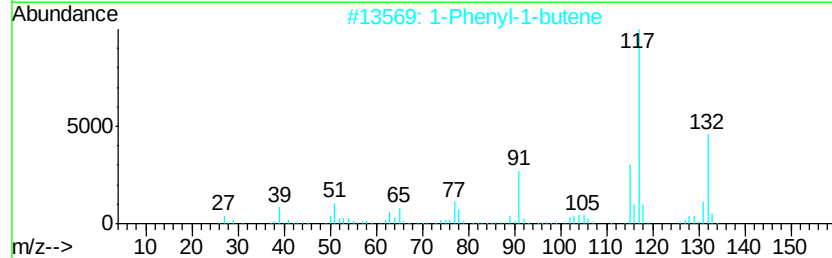
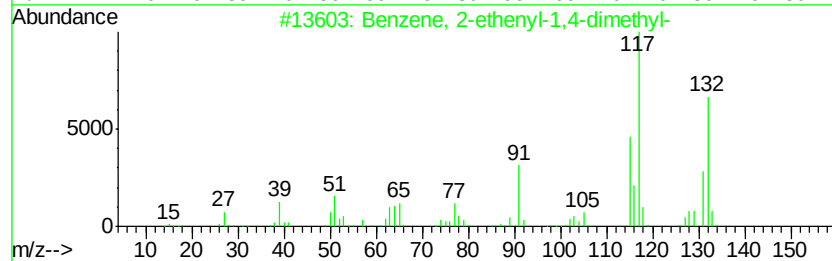
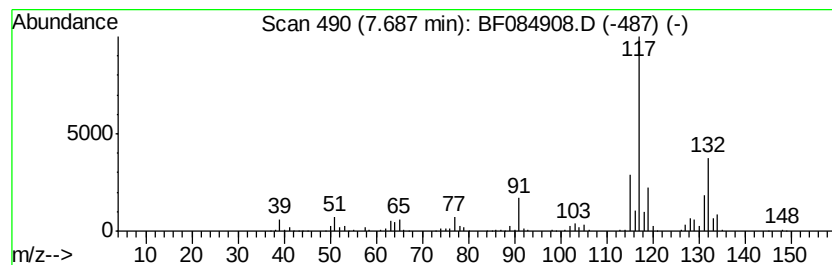
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 17

| R.T. | EstConc | Area    | Relative to ISTD | R.T. |
|------|---------|---------|------------------|------|
| 7.69 | 3.71 ng | 1030450 | Napthalene-d8    | 7.95 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethenyl-1,4-dimethyl-  | 132 | C10H12  | 002039-89-6 | 91   |
| 2    |    | 1-Phenyl-1-butene                 | 132 | C10H12  | 000824-90-8 | 87   |
| 3    |    | Benzene, 2-butenyl-               | 132 | C10H12  | 001560-06-1 | 87   |
| 4    |    | 1H-Indene, 2,3-dihydro-2-methyl-  | 132 | C10H12  | 000824-63-5 | 83   |
| 5    |    | Benzene, 1-methyl-2-(2-propenyl)- | 132 | C10H12  | 001587-04-8 | 83   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020616\  
 Data File : BF084908.D  
 Acq On : 6 Feb 2016 13:35  
 Operator : SJ/IZ  
 Sample : H1310-02  
 Misc :  
 ALS Vial : 50 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SUP-B008(25-27)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.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 |
| n-Propyl acetate     | 2.37 | 4.9     | ng    | 278045   | 1                     | 6.66 | 1140450 | 20.0 |
| 3-Penten-2-one, 4... | 4.12 | 2.6     | ng    | 150886   | 1                     | 6.66 | 1140450 | 20.0 |
| 2-Pentanone, 4-hy... | 4.83 | 39.7    | ng    | 2262440  | 1                     | 6.66 | 1140450 | 20.0 |
| Benzene, (1-methy... | 5.81 | 3.0     | ng    | 169975   | 1                     | 6.66 | 1140450 | 20.0 |
| Benzene, propyl-     | 6.11 | 8.3     | ng    | 473356   | 1                     | 6.66 | 1140450 | 20.0 |
| Benzene, 1-ethyl-... | 6.18 | 26.0    | ng    | 1484500  | 1                     | 6.66 | 1140450 | 20.0 |
| Benzene, 1,2,3-tr... | 6.26 | 21.0    | ng    | 1198760  | 1                     | 6.66 | 1140450 | 20.0 |
| unknown6.43          | 6.43 | 84.5    | ng    | 4816070  | 1                     | 6.66 | 1140450 | 20.0 |
| Benzene, 1,2,4-tr... | 6.72 | 23.8    | ng    | 1359260  | 1                     | 6.66 | 1140450 | 20.0 |
| Benzene, 1,4-diet... | 6.92 | 4.4     | ng    | 251072   | 1                     | 6.66 | 1140450 | 20.0 |
| Benzene, 1-methyl... | 6.94 | 3.1     | ng    | 175840   | 1                     | 6.66 | 1140450 | 20.0 |
| Benzene, 1-methyl... | 6.99 | 11.1    | ng    | 634205   | 1                     | 6.66 | 1140450 | 20.0 |
| Benzene, 2-ethyl-... | 7.14 | 9.4     | ng    | 533894   | 1                     | 6.66 | 1140450 | 20.0 |
| Benzene, 2-etheny... | 7.69 | 3.7     | ng    | 1030450  | 2                     | 7.95 | 5555800 | 20.0 |