

Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\  
 Data File : BF089489.D  
 Acq On : 1 Aug 2016 2:03  
 Operator : UM/SJ  
 Sample : H4307-01  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 1057-SOUTH

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-BF072716.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         | 1.667       | 5             | 7           | 46           | rVB      | 880262         | 2006172       | 15.74%          | 1.253%        |
| 2         | 3.827       | 192           | 196         | 201          | rBV      | 88918          | 172390        | 1.35%           | 0.108%        |
| 3         | 4.376       | 242           | 244         | 246          | rBV      | 129051         | 174674        | 1.37%           | 0.109%        |
| 4         | 4.479       | 250           | 253         | 255          | rVB      | 244670         | 381006        | 2.99%           | 0.238%        |
| 5         | 4.536       | 255           | 258         | 260          | rBV      | 210505         | 288492        | 2.26%           | 0.180%        |
| 6         | 4.570       | 260           | 261         | 264          | rVB      | 141903         | 234966        | 1.84%           | 0.147%        |
| 7         | 4.787       | 277           | 280         | 284          | rBV2     | 274646         | 627984        | 4.93%           | 0.392%        |
| 8         | 4.901       | 288           | 290         | 291          | rBV      | 354210         | 428612        | 3.36%           | 0.268%        |
| 9         | 4.924       | 291           | 292         | 295          | rVB      | 319341         | 402625        | 3.16%           | 0.251%        |
| 10        | 4.981       | 295           | 297         | 298          | rBV2     | 117424         | 151548        | 1.19%           | 0.095%        |
| 11        | 5.016       | 298           | 300         | 302          | rVB      | 684965         | 765400        | 6.01%           | 0.478%        |
| 12        | 5.061       | 302           | 304         | 306          | rBV      | 934403         | 1117674       | 8.77%           | 0.698%        |
| 13        | 5.096       | 306           | 307         | 310          | rVB      | 197911         | 203401        | 1.60%           | 0.127%        |
| 14        | 5.187       | 313           | 315         | 317          | rBV      | 354034         | 556263        | 4.36%           | 0.347%        |
| 15        | 5.244       | 317           | 320         | 321          | rVB      | 420093         | 534641        | 4.20%           | 0.334%        |
| 16        | 5.290       | 321           | 324         | 328          | rVB2     | 378457         | 781962        | 6.14%           | 0.488%        |
| 17        | 5.370       | 328           | 331         | 335          | rBV      | 653750         | 862454        | 6.77%           | 0.538%        |
| 18        | 5.462       | 335           | 339         | 342          | rBV2     | 535968         | 1074294       | 8.43%           | 0.671%        |
| 19        | 5.542       | 344           | 346         | 350          | rVB3     | 199585         | 456247        | 3.58%           | 0.285%        |
| 20        | 5.610       | 350           | 352         | 356          | rVB2     | 649421         | 1446960       | 11.35%          | 0.903%        |
| 21        | 5.690       | 356           | 359         | 360          | rBV      | 418901         | 715621        | 5.62%           | 0.447%        |
| 22        | 5.736       | 360           | 363         | 366          | rVB3     | 1625225        | 3221927       | 25.28%          | 2.012%        |
| 23        | 5.793       | 366           | 368         | 376          | rVB4     | 992835         | 2517766       | 19.76%          | 1.572%        |
| 24        | 5.942       | 376           | 381         | 383          | rBV3     | 1061501        | 2003525       | 15.72%          | 1.251%        |
| 25        | 6.010       | 383           | 387         | 389          | rBV2     | 1929873        | 3497381       | 27.44%          | 2.184%        |
| 26        | 6.044       | 389           | 390         | 393          | rVB2     | 1178556        | 1529611       | 12.00%          | 0.955%        |
| 27        | 6.102       | 393           | 395         | 397          | rBV      | 1350057        | 1753454       | 13.76%          | 1.095%        |
| 28        | 6.159       | 397           | 400         | 405          | rVB3     | 1221555        | 3143252       | 24.66%          | 1.962%        |
| 29        | 6.273       | 405           | 410         | 412          | rBV3     | 1585595        | 4263314       | 33.45%          | 2.662%        |
| 30        | 6.342       | 412           | 416         | 423          | rVB2     | 2440295        | 6109792       | 47.94%          | 3.815%        |
| 31        | 6.467       | 423           | 427         | 428          | rBV4     | 810032         | 1617174       | 12.69%          | 1.010%        |
| 32        | 6.547       | 432           | 434         | 442          | rVB3     | 1501796        | 5171523       | 40.58%          | 3.229%        |
| 33        | 6.685       | 442           | 446         | 451          | rVB4     | 1798238        | 5328700       | 41.81%          | 3.327%        |
| 34        | 6.810       | 451           | 457         | 458          | rBV5     | 1106648        | 3696274       | 29.00%          | 2.308%        |

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 Sampling : 1  
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 Stop Thrs : 0  
 Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |          |         |        |
|----|--------|------|------|------|------|---------|----------|---------|--------|
| 35 | 6.845  | 458  | 460  | 461  | rVV  | 1304098 | 1527391  | 11.98%  | 0.954% |
| 36 | 6.913  | 464  | 466  | 471  | rVB3 | 910994  | 2802431  | 21.99%  | 1.750% |
| 37 | 7.016  | 471  | 475  | 477  | rBV  | 1003135 | 2412569  | 18.93%  | 1.506% |
| 38 | 7.085  | 477  | 481  | 486  | rVB3 | 1490280 | 4540985  | 35.63%  | 2.835% |
| 39 | 7.187  | 487  | 490  | 493  | rVV4 | 615316  | 1421874  | 11.16%  | 0.888% |
| 40 | 7.325  | 500  | 502  | 508  | rVB3 | 1797307 | 5672689  | 44.51%  | 3.542% |
| 41 | 7.473  | 509  | 515  | 519  | rBV4 | 1595750 | 7651952  | 60.04%  | 4.777% |
| 42 | 7.576  | 519  | 524  | 529  | rVV7 | 1661515 | 8255706  | 64.78%  | 5.154% |
| 43 | 7.656  | 529  | 531  | 534  | rVB2 | 1277821 | 2656908  | 20.85%  | 1.659% |
| 44 | 7.827  | 540  | 546  | 552  | rBV5 | 2196920 | 12744419 | 100.00% | 7.957% |
| 45 | 8.033  | 560  | 564  | 567  | rBV3 | 1130109 | 3543897  | 27.81%  | 2.213% |
| 46 | 8.296  | 583  | 587  | 589  | rBV  | 1642867 | 4099229  | 32.16%  | 2.559% |
| 47 | 8.662  | 614  | 619  | 623  | rBV3 | 1962977 | 6205250  | 48.69%  | 3.874% |
| 48 | 8.890  | 634  | 639  | 642  | rVB3 | 1667097 | 4584464  | 35.97%  | 2.862% |
| 49 | 9.268  | 668  | 672  | 676  | rBV2 | 1559114 | 6274331  | 49.23%  | 3.917% |
| 50 | 9.576  | 697  | 699  | 701  | rBV3 | 984045  | 1666175  | 13.07%  | 1.040% |
| 51 | 9.839  | 720  | 722  | 726  | rVB3 | 942606  | 2319611  | 18.20%  | 1.448% |
| 52 | 9.919  | 726  | 729  | 732  | rBV4 | 898270  | 2316578  | 18.18%  | 1.446% |
| 53 | 10.262 | 756  | 759  | 764  | rVB3 | 1905787 | 5002108  | 39.25%  | 3.123% |
| 54 | 10.536 | 778  | 783  | 785  | rBV2 | 1916977 | 4492919  | 35.25%  | 2.805% |
| 55 | 10.982 | 820  | 822  | 825  | rVB2 | 1806000 | 3295768  | 25.86%  | 2.058% |
| 56 | 11.085 | 829  | 831  | 834  | rVV2 | 806344  | 1235817  | 9.70%   | 0.772% |
| 57 | 11.554 | 870  | 872  | 873  | rBV  | 605908  | 827689   | 6.49%   | 0.517% |
| 58 | 12.045 | 912  | 915  | 917  | rVB2 | 836103  | 1584693  | 12.43%  | 0.989% |
| 59 | 12.102 | 917  | 920  | 921  | rBV  | 1079617 | 1742180  | 13.67%  | 1.088% |
| 60 | 12.537 | 956  | 958  | 959  | rBV  | 481753  | 577168   | 4.53%   | 0.360% |
| 61 | 12.617 | 963  | 965  | 969  | rVB2 | 1054289 | 1642615  | 12.89%  | 1.026% |
| 62 | 12.891 | 987  | 989  | 994  | rVB4 | 439568  | 1032307  | 8.10%   | 0.645% |
| 63 | 12.971 | 994  | 996  | 998  | rVB3 | 208329  | 259818   | 2.04%   | 0.162% |
| 64 | 13.634 | 1052 | 1054 | 1057 | rVB  | 246746  | 342192   | 2.69%   | 0.214% |
| 65 | 14.948 | 1167 | 1169 | 1177 | rVB  | 178258  | 202111   | 1.59%   | 0.126% |

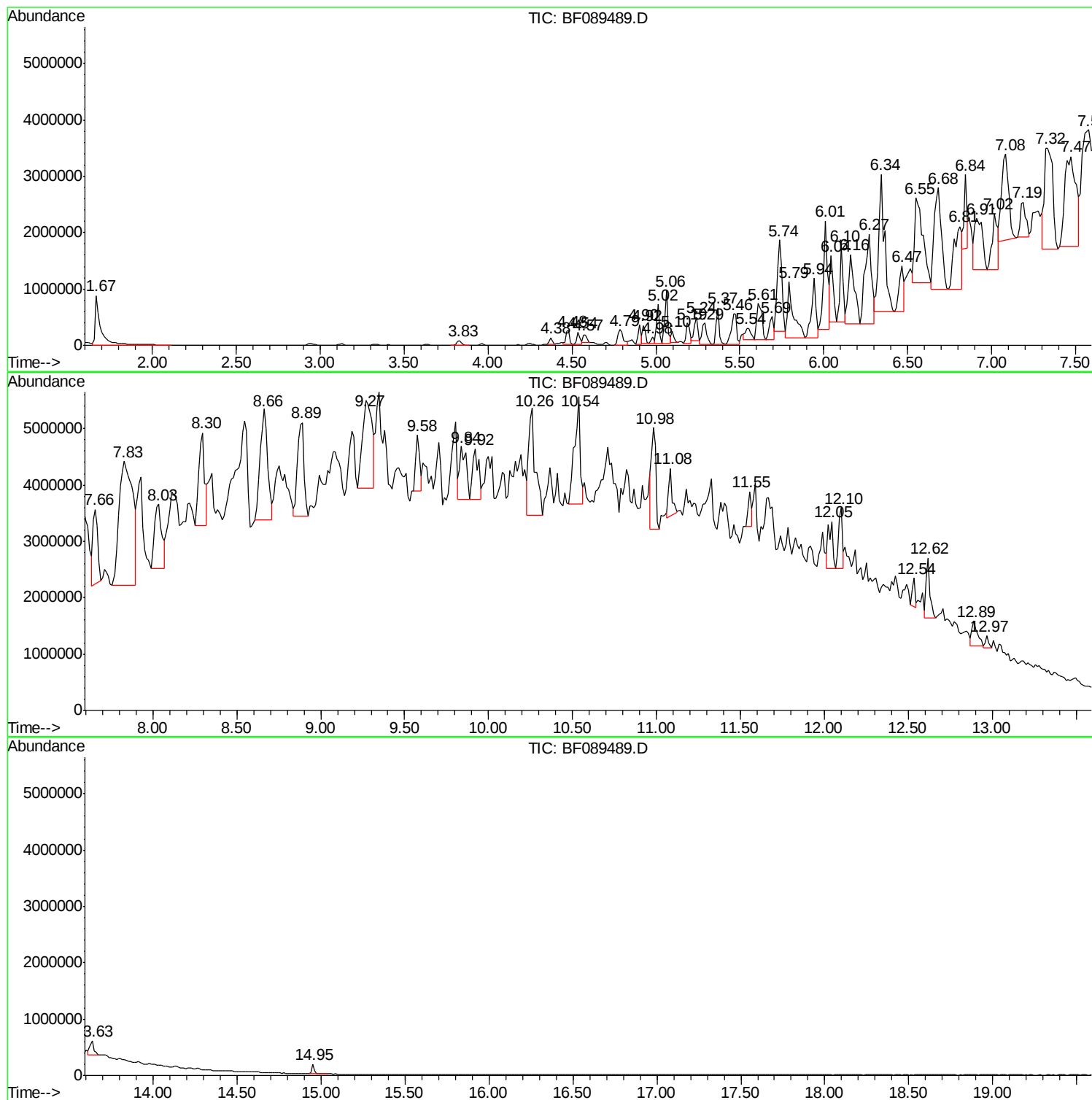
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Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

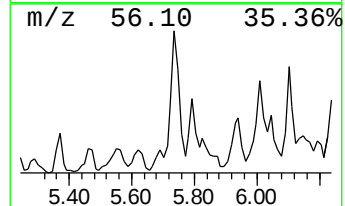
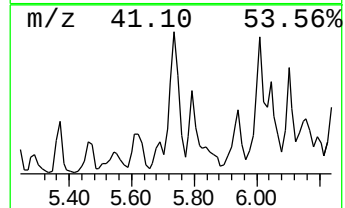
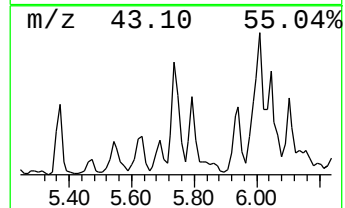
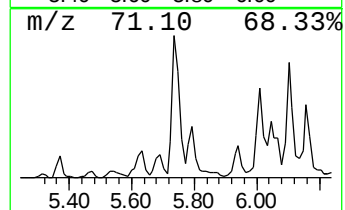
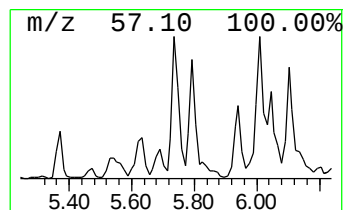
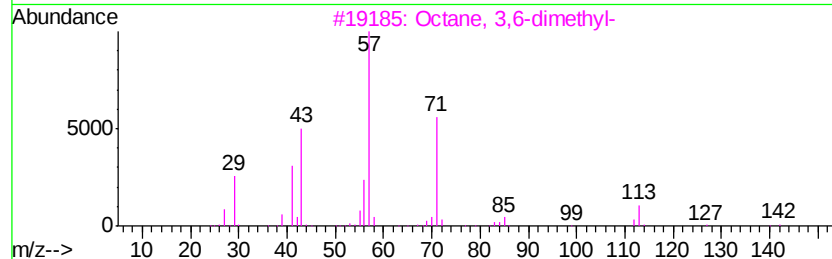
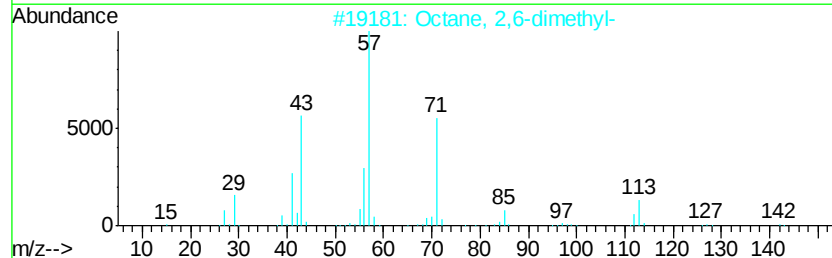
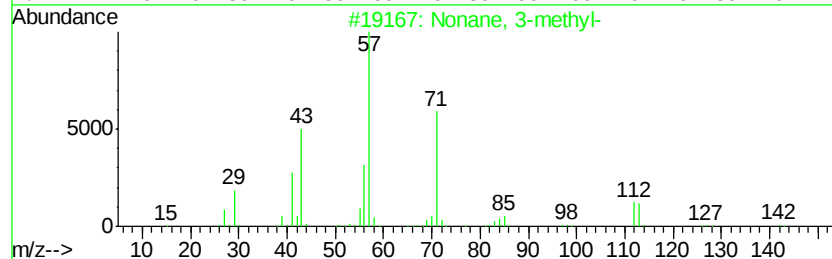
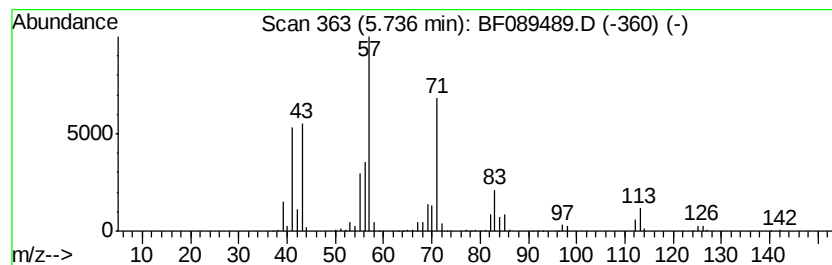
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Nonane, 3-methyl- Concentration Rank 11

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.74 | 39.85 ng | 3221930 | 1,4-Dichlorobenzene-d4 | 6.50 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Nonane, 3-methyl-                   | 142 | C10H22    | 005911-04-6  | 81   |
| 2    |    | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 72   |
| 3    |    | Octane, 3,6-dimethyl-               | 142 | C10H22    | 015869-94-0  | 64   |
| 4    |    | Sulfurous acid, 2-ethylhexyl pen... | 264 | C13H28O3S | 1000309-18-9 | 53   |
| 5    |    | 3-Bromooctane                       | 192 | C8H17Br   | 000999-64-4  | 52   |



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ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

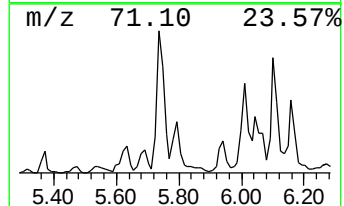
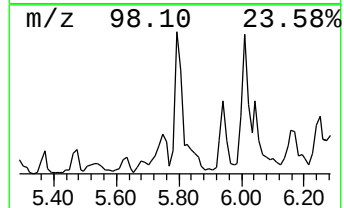
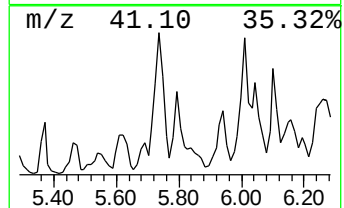
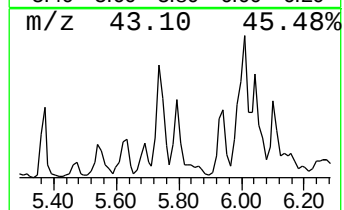
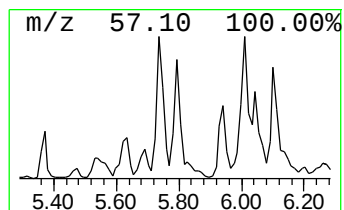
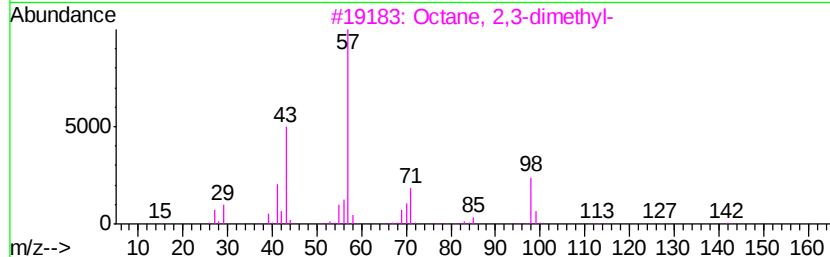
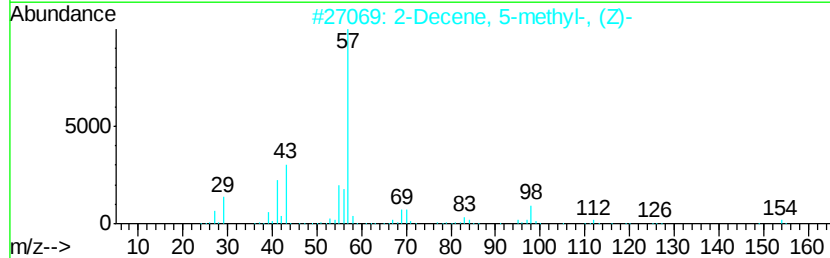
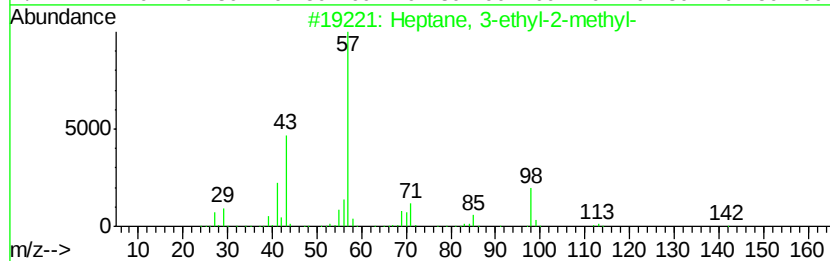
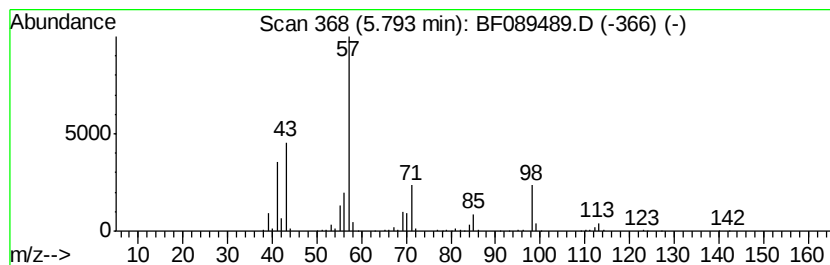
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Heptane, 3-ethyl-2-methyl- Concentration Rank 14

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.79 | 31.14 ng | 2517770 | 1,4-Dichlorobenzene-d4 | 6.50 |

| Hit# | of | Tentative ID                      | MW  | MolForm  | CAS#         | Qual |
|------|----|-----------------------------------|-----|----------|--------------|------|
| 1    | 5  | Heptane, 3-ethyl-2-methyl-        | 142 | C10H22   | 014676-29-0  | 72   |
| 2    |    | 2-Decene, 5-methyl-, (Z)-         | 154 | C11H22   | 074645-86-6  | 59   |
| 3    |    | Octane, 2,3-dimethyl-             | 142 | C10H22   | 007146-60-3  | 52   |
| 4    |    | Oxalic acid, isobutyl nonyl ester | 272 | C15H28O4 | 1000309-37-4 | 50   |
| 5    |    | Hexane, 3-ethyl-2,5-dimethyl-     | 142 | C10H22   | 052897-04-8  | 50   |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

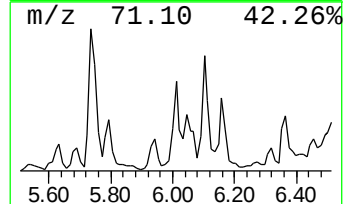
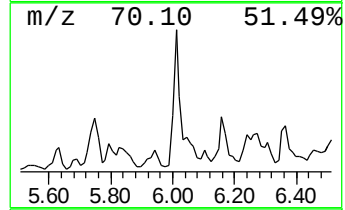
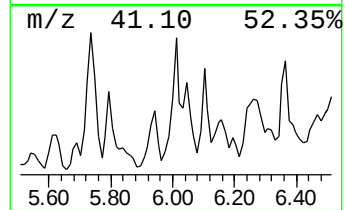
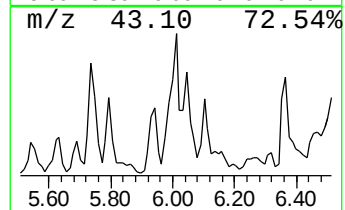
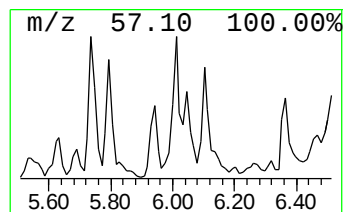
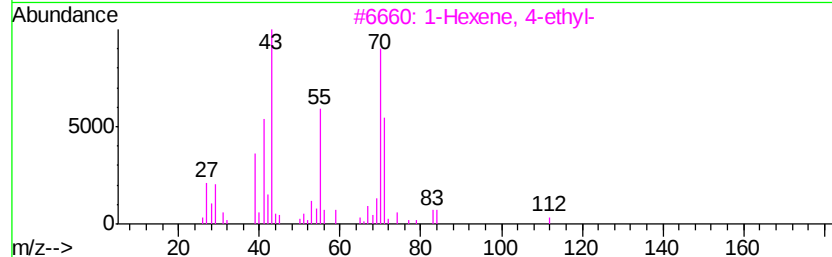
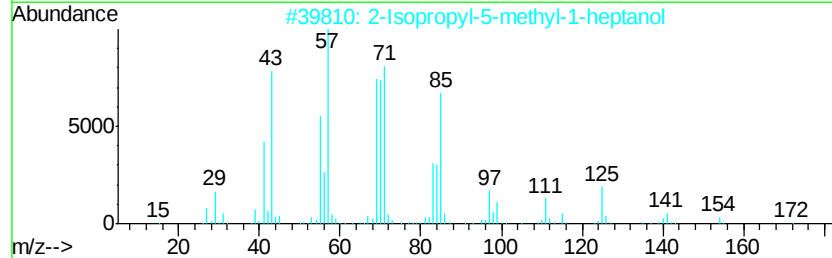
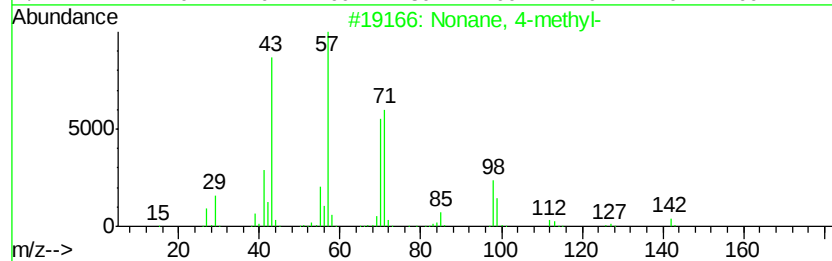
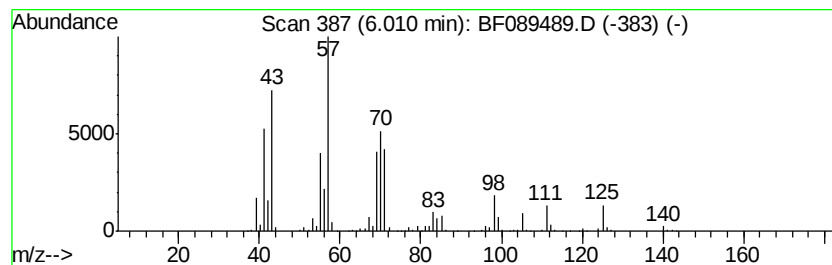
TIC Library : C:\Database\NIST11.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 4 Nonane, 4-methyl- Concentration Rank 10

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.01 | 43.25 ng | 3497380 | 1,4-Dichlorobenzene-d4 | 6.50 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 4-methyl-               | 142 | C10H22  | 017301-94-9 | 53   |
| 2    |    | 2-Isopropyl-5-methyl-1-heptanol | 172 | C11H24O | 091337-07-4 | 50   |
| 3    |    | 1-Hexene, 4-ethyl-              | 112 | C8H16   | 016746-85-3 | 50   |
| 4    |    | 1-Hexene, 3,5,5-trimethyl-      | 126 | C9H18   | 004316-65-8 | 43   |
| 5    |    | Nonane, 2-methyl-               | 142 | C10H22  | 000871-83-0 | 43   |



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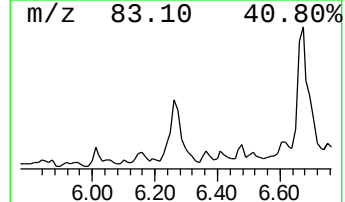
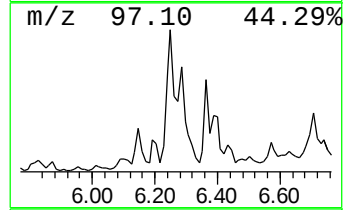
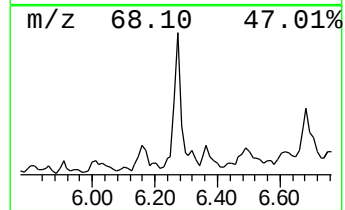
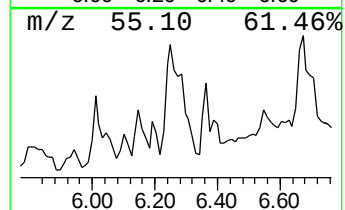
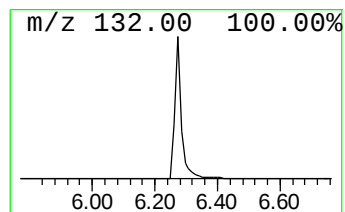
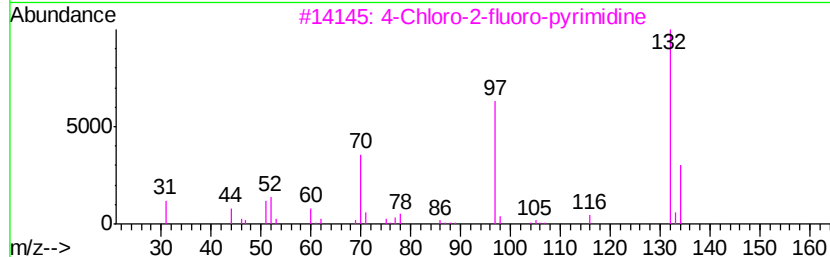
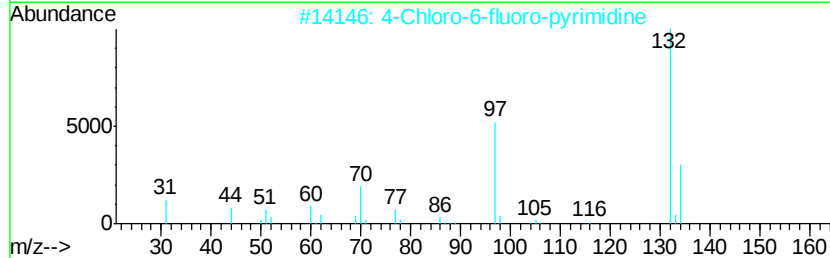
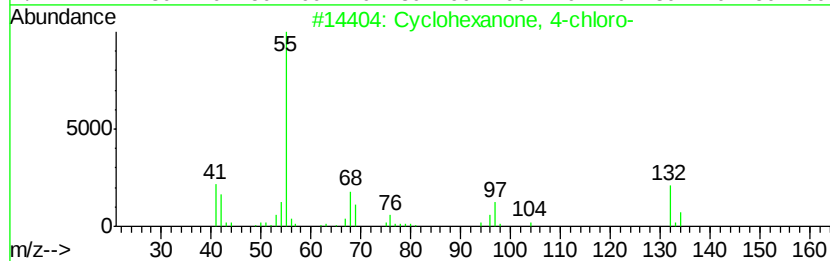
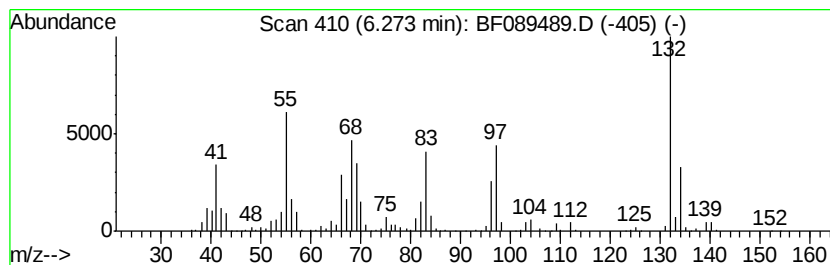
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 unknown6.27 Concentration Rank 8

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.27 | 52.73 ng | 4263310 | 1,4-Dichlorobenzene-d4 | 6.50 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Cyclohexanone, 4-chloro-            | 132 | C6H9ClO   | 021299-26-3 | 38   |
| 2    |    | 4-Chloro-6-fluoro-pyrimidine        | 132 | C4H2ClFN2 | 051422-01-6 | 27   |
| 3    |    | 4-Chloro-2-fluoro-pyrimidine        | 132 | C4H2ClFN2 | 051422-00-5 | 27   |
| 4    |    | Cyclopentane, 2-isopropyl-1,3-di... | 140 | C10H20    | 032281-85-9 | 14   |
| 5    |    | Cyclopentane, 1-methyl-3-(2-meth... | 140 | C10H20    | 029053-04-1 | 14   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

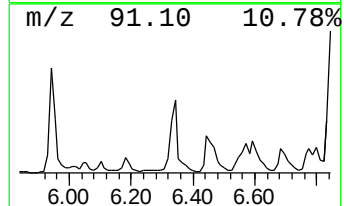
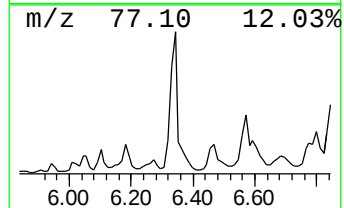
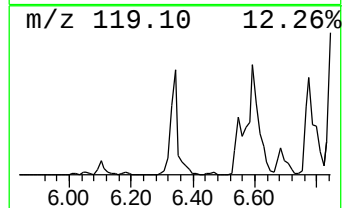
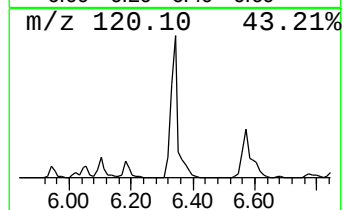
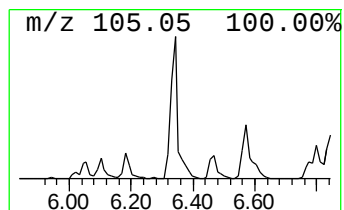
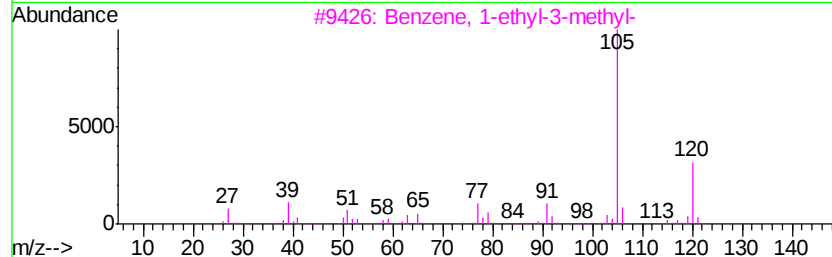
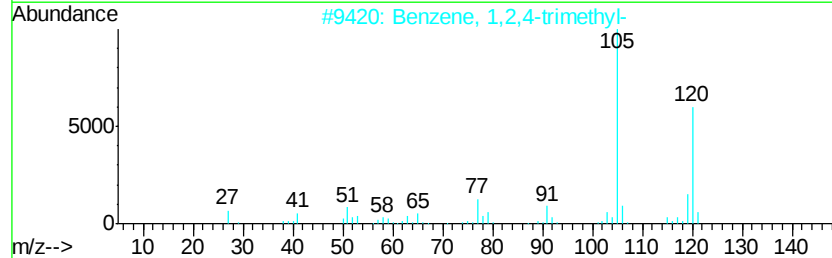
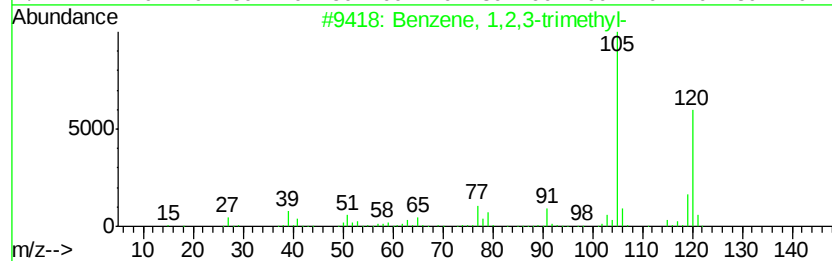
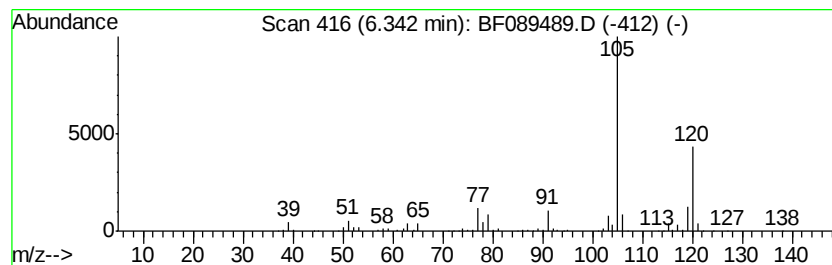
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.34 | 75.56 ng | 6109790 | 1,4-Dichlorobenzene-d4 | 6.50 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 95   |
| 3    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 4    |    | Mesitylene                 | 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\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

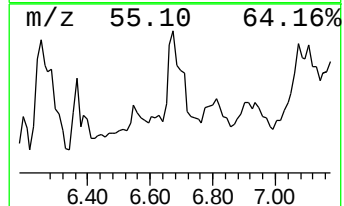
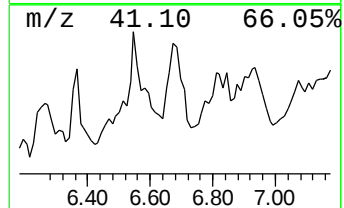
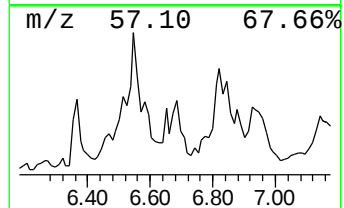
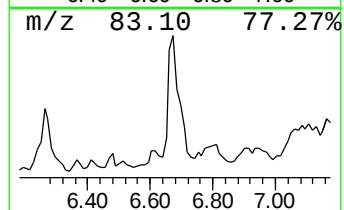
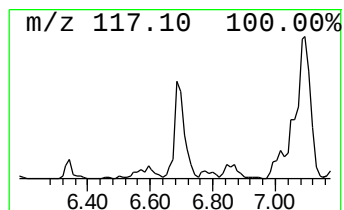
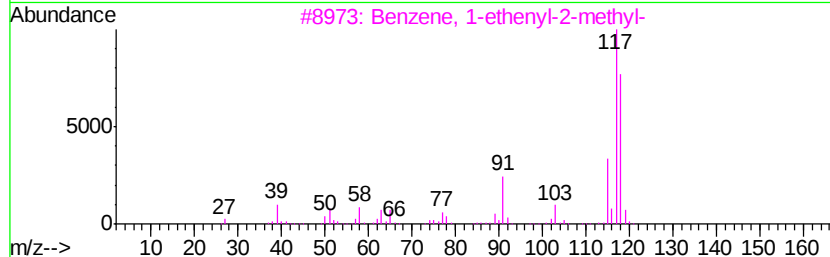
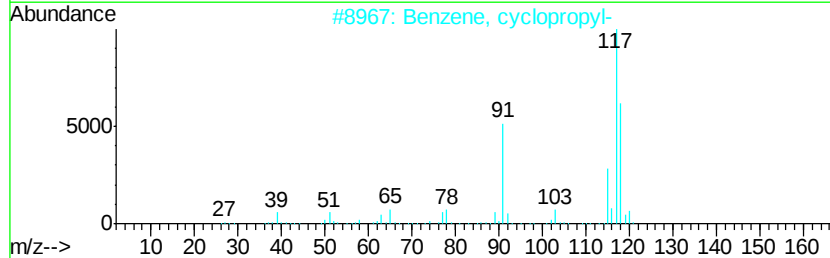
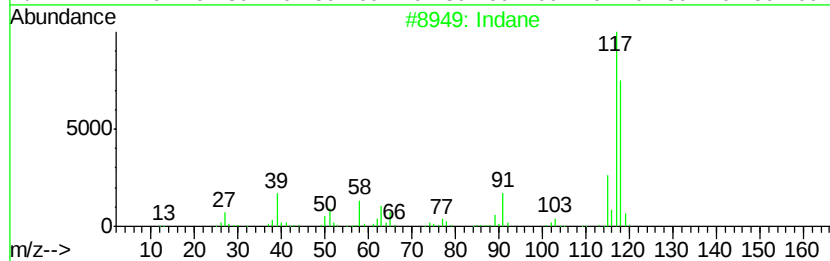
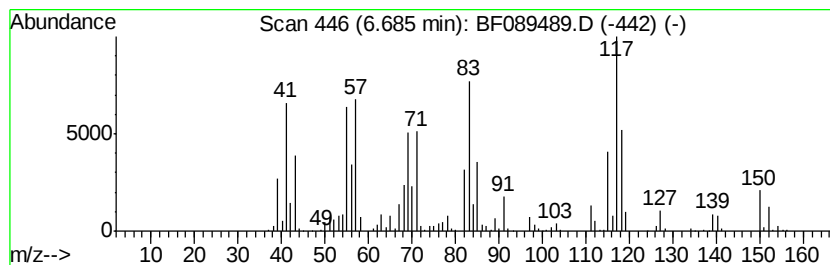
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 unknown6.68 Concentration Rank 4

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.68 | 65.90 ng | 5328700 | 1,4-Dichlorobenzene-d4 | 6.50 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#         | Qual |
|------|----|----------------------------------|-----|---------|--------------|------|
| 1    | 5  | Indane                           | 118 | C9H10   | 000496-11-7  | 38   |
| 2    |    | Benzene, cyclopropyl-            | 118 | C9H10   | 000873-49-4  | 30   |
| 3    |    | Benzene, 1-ethenyl-2-methyl-     | 118 | C9H10   | 000611-15-4  | 30   |
| 4    |    | Tetracyclo[3.3.1.0(2,8).0(4,6)]- | 118 | C9H10   | 1000191-13-7 | 30   |
| 5    |    | Deltacyclene                     | 118 | C9H10   | 007785-10-6  | 30   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

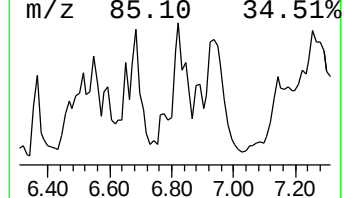
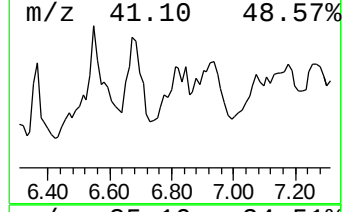
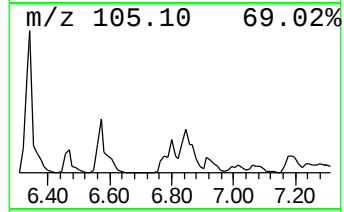
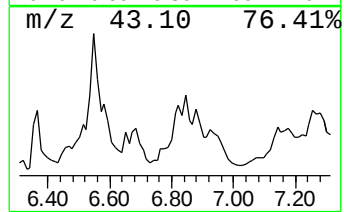
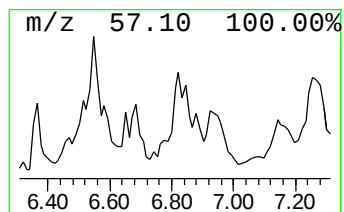
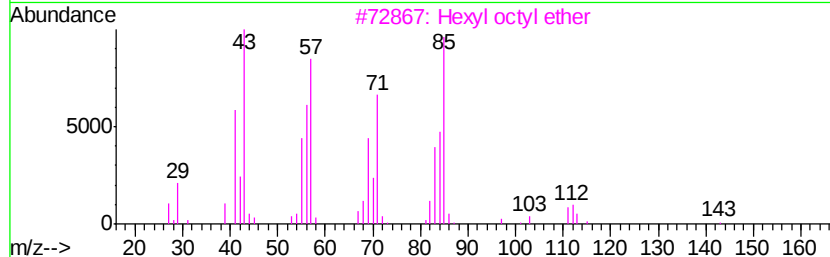
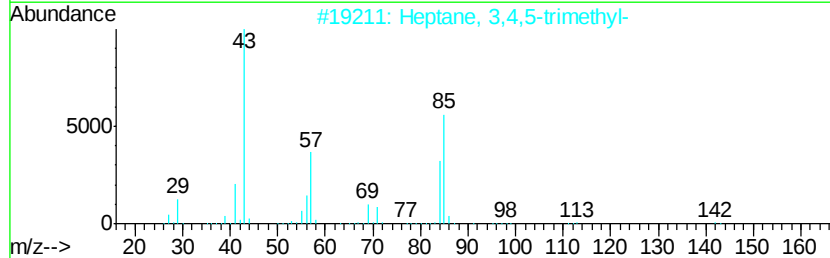
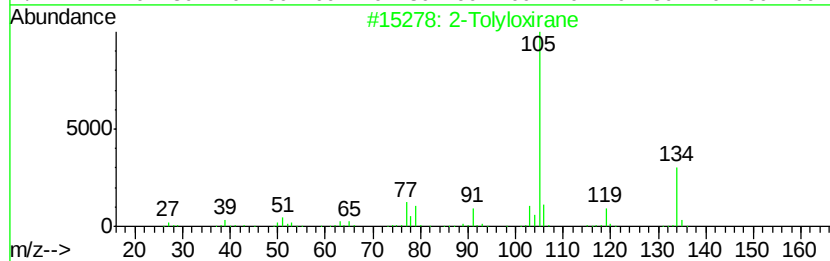
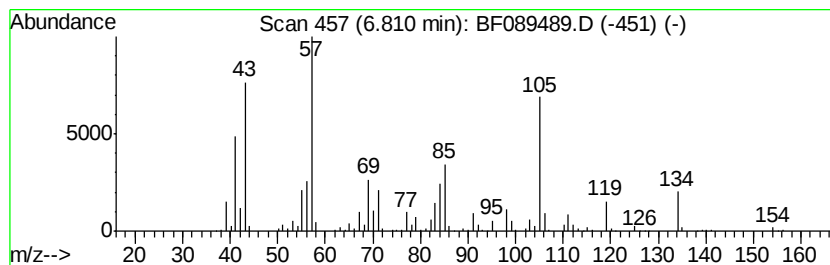
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 2-Tolyloxirane Concentration Rank 9

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.81 | 45.71 ng | 3696270 | 1,4-Dichlorobenzene-d4 | 6.50 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Tolyloxirane            | 134 | C9H10O  | 002783-26-8 | 53   |
| 2    |    | Heptane, 3,4,5-trimethyl- | 142 | C10H22  | 020278-89-1 | 52   |
| 3    |    | Hexyl octyl ether         | 214 | C14H30O | 017071-54-4 | 43   |
| 4    |    | Heptane, 4-ethyl-         | 128 | C9H20   | 002216-32-2 | 43   |
| 5    |    | Hexane, 2,4-dimethyl-     | 114 | C8H18   | 000589-43-5 | 35   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

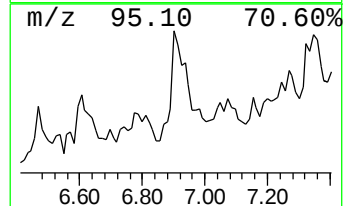
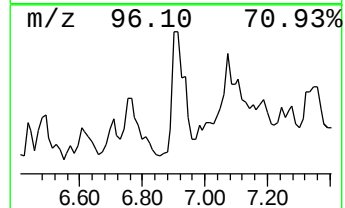
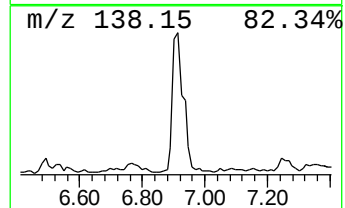
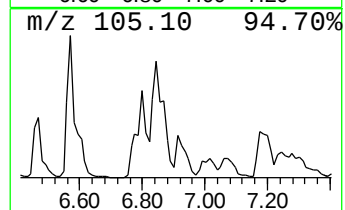
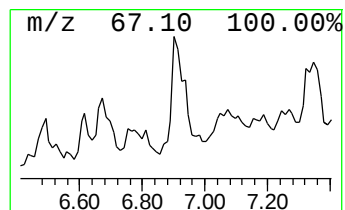
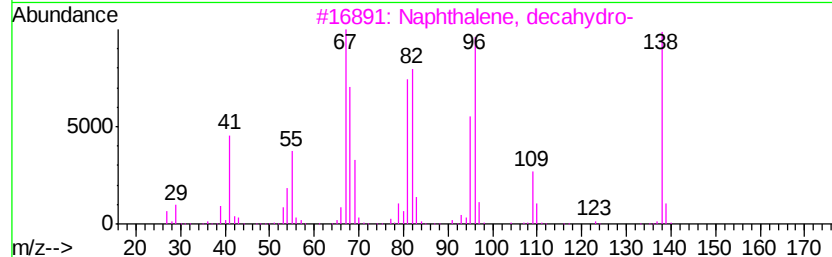
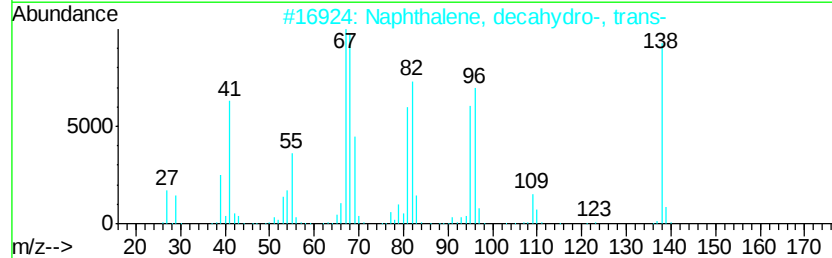
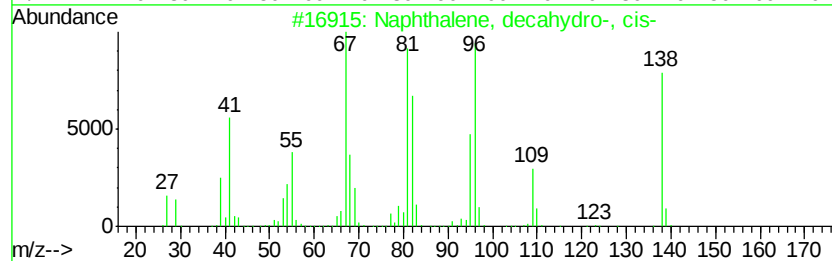
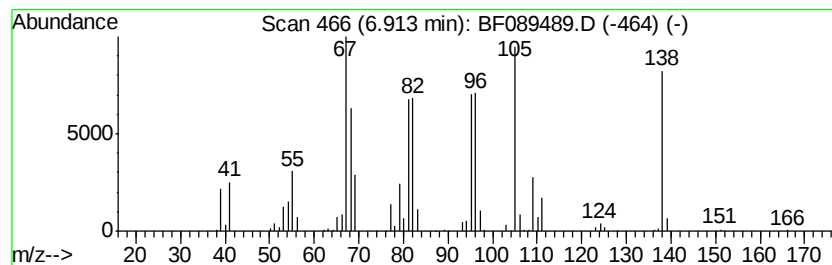
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Naphthalene, decahydro-, cis- Concentration Rank 12

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.91 | 34.66 ng | 2802430 | 1,4-Dichlorobenzene-d4 | 6.50 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, decahydro-, cis-   | 138 | C10H18  | 000493-01-6 | 96   |
| 2    |    | Naphthalene, decahydro-, trans- | 138 | C10H18  | 000493-02-7 | 95   |
| 3    |    | Naphthalene, decahydro-         | 138 | C10H18  | 000091-17-8 | 93   |
| 4    |    | Spiro[4.5]decane                | 138 | C10H18  | 000176-63-6 | 87   |
| 5    |    | Cyclohexene, 1-butyl-           | 138 | C10H18  | 003282-53-9 | 76   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

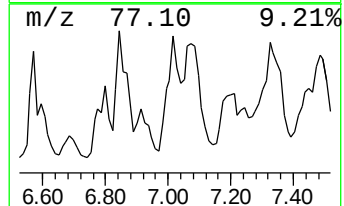
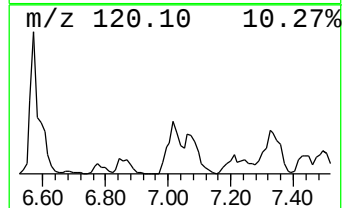
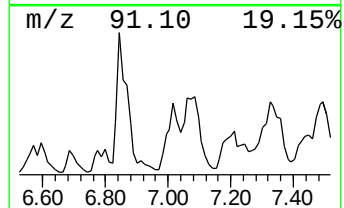
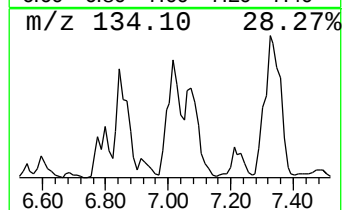
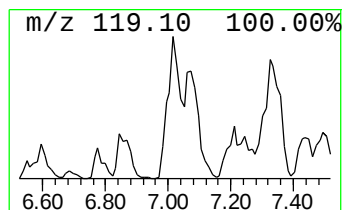
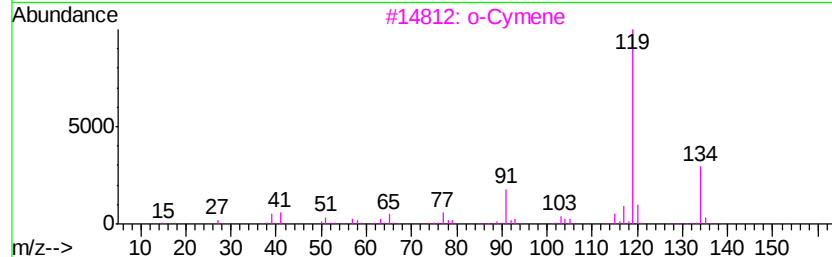
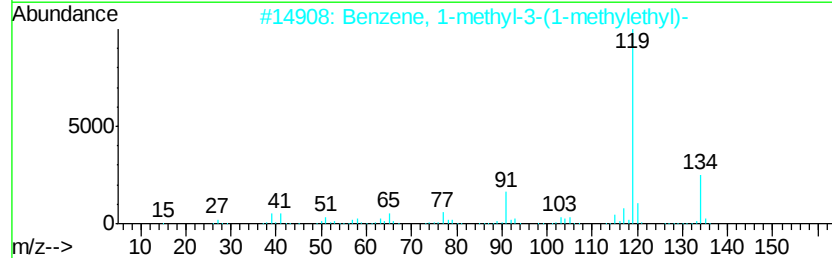
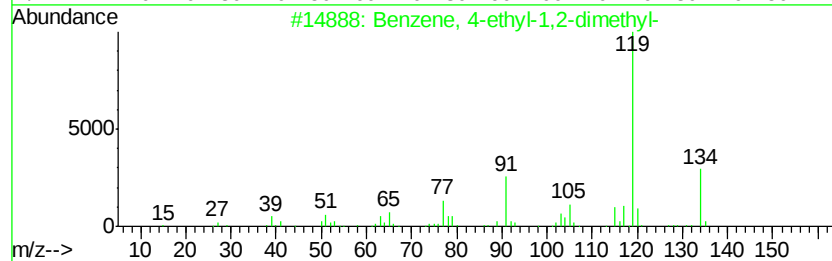
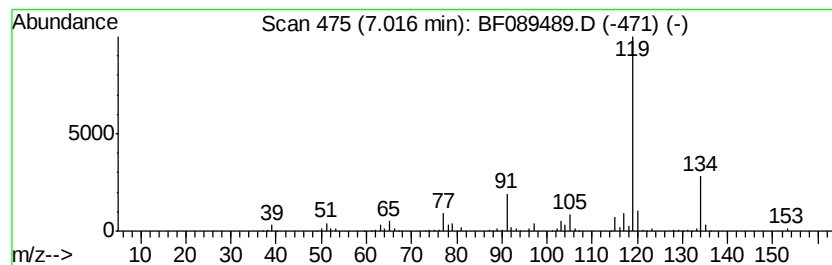
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 15

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.02 | 29.84 ng | 2412570 | 1,4-Dichlorobenzene-d4 | 6.50 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-ethyl-1,2-dimethyl-       | 134 | C10H14  | 000934-80-5 | 96   |
| 2    |    | Benzene, 1-methyl-3-(1-methylethyl-) | 134 | C10H14  | 000535-77-3 | 95   |
| 3    |    | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 95   |
| 4    |    | p-Cymene                             | 134 | C10H14  | 000099-87-6 | 94   |
| 5    |    | Benzene, 1-ethyl-2,3-dimethyl-       | 134 | C10H14  | 000933-98-2 | 93   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

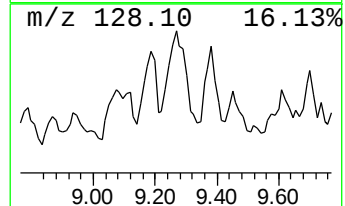
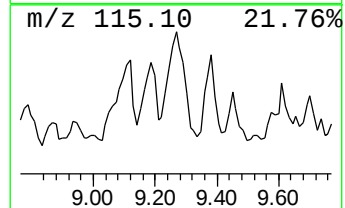
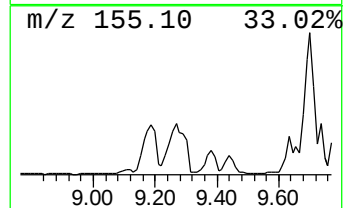
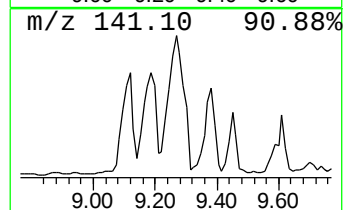
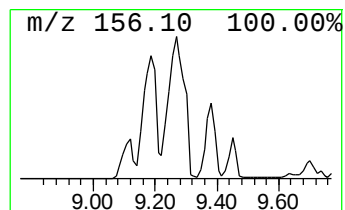
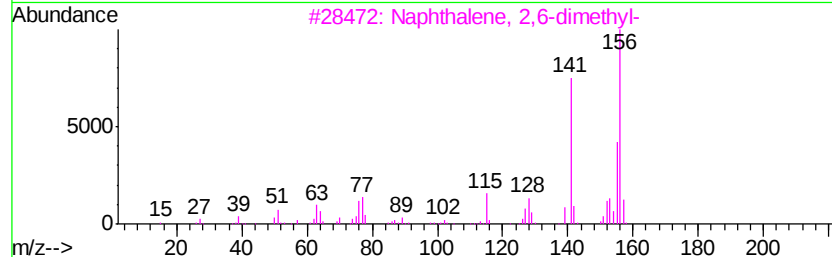
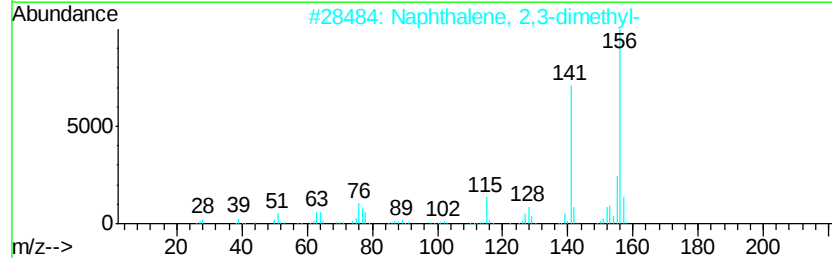
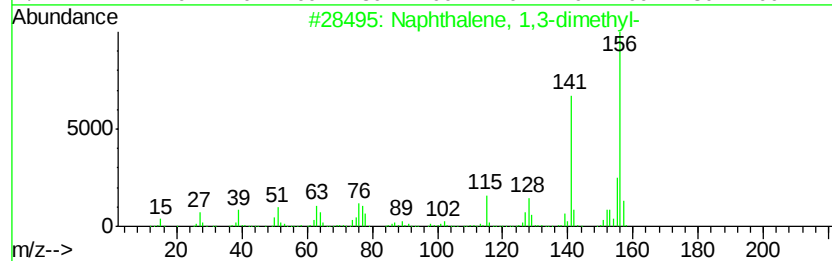
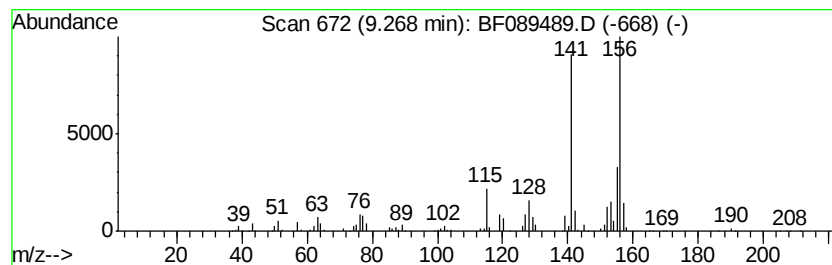
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BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Naphthalene, 1,3-dimethyl- Concentration Rank 2

| R.T.      | EstConc                    | Area    | Relative to ISTD |             | R.T. |
|-----------|----------------------------|---------|------------------|-------------|------|
| 9.27      | 75.31 ng                   | 6274330 | Acenaphthene-d10 |             | 9.59 |
| Hit# of 5 | Tentative ID               | MW      | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 1,3-dimethyl- | 156     | C12H12           | 000575-41-7 | 97   |
| 2         | Naphthalene, 2,3-dimethyl- | 156     | C12H12           | 000581-40-8 | 97   |
| 3         | Naphthalene, 2,6-dimethyl- | 156     | C12H12           | 000581-42-0 | 97   |
| 4         | Naphthalene, 1,7-dimethyl- | 156     | C12H12           | 000575-37-1 | 97   |
| 5         | Naphthalene, 1,4-dimethyl- | 156     | C12H12           | 000571-58-4 | 97   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

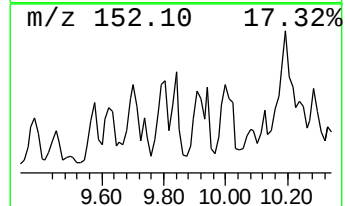
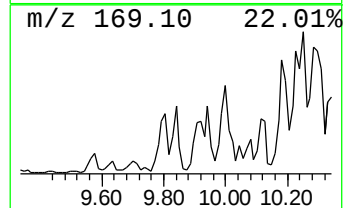
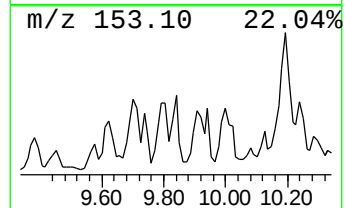
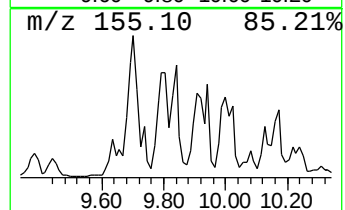
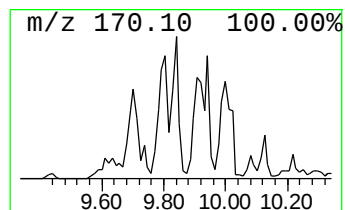
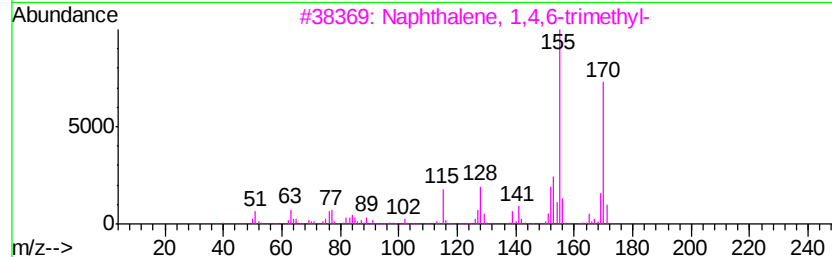
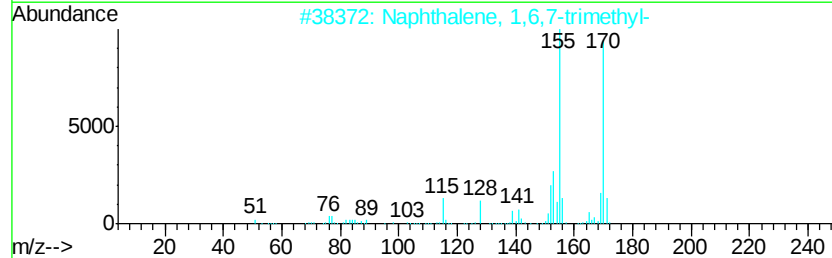
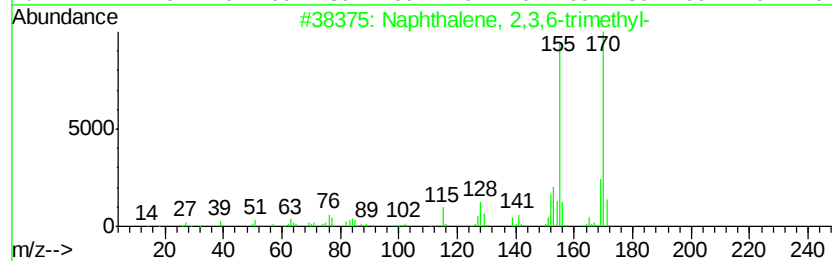
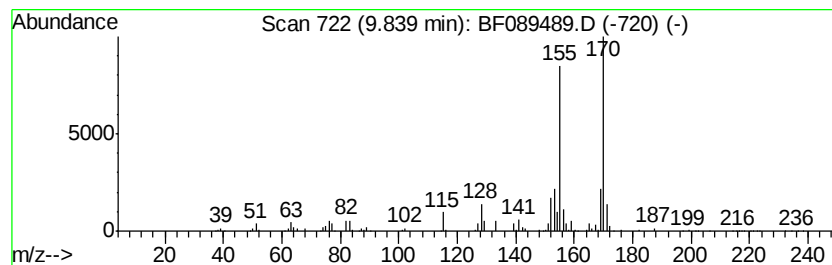
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Naphthalene, 2,3,6-trimethyl- Concentration Rank 17

| R.T. | EstConc  | Area    | Relative to ISTD | R.T. |
|------|----------|---------|------------------|------|
| 9.84 | 27.84 ng | 2319610 | Acenaphthene-d10 | 9.59 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 97   |
| 2    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 96   |
| 3    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 96   |
| 4    |    |   | 4,6,8-Trimethylazulene        | 170 | C13H14  | 000941-81-1 | 94   |
| 5    |    |   | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 93   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

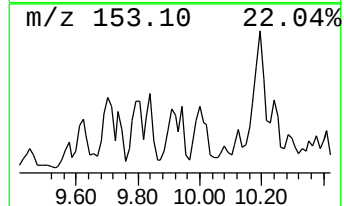
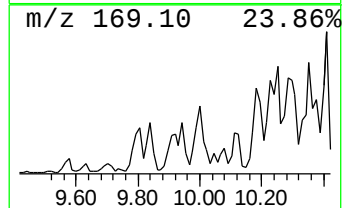
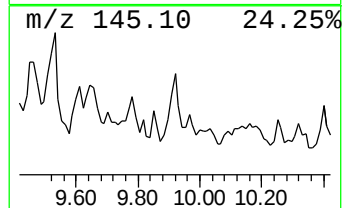
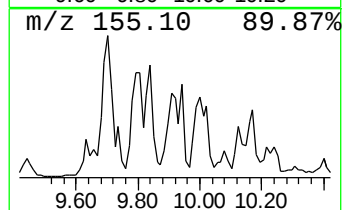
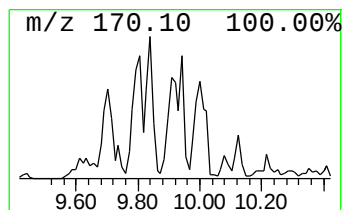
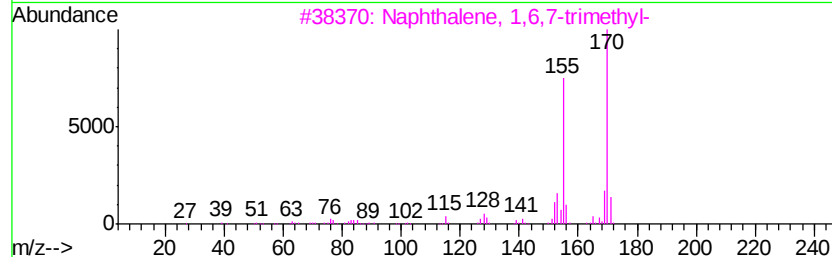
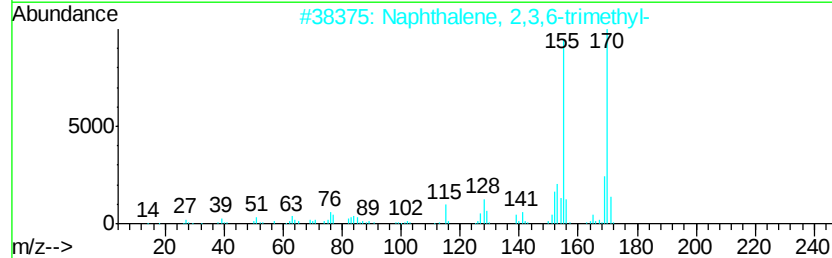
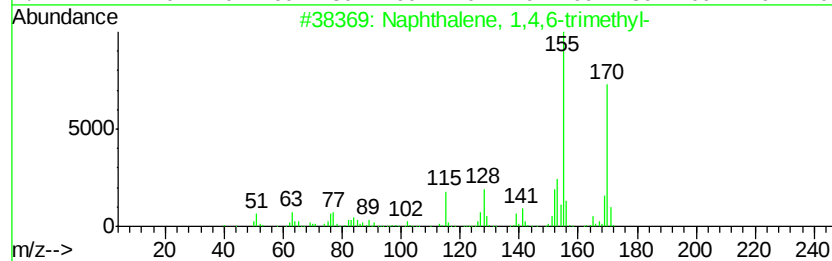
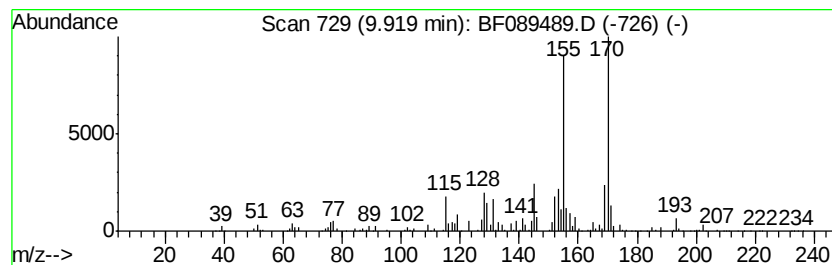
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Naphthalene, 1,4,6-trimethyl- Concentration Rank 18

| R.T. | EstConc  | Area    | Relative to ISTD | R.T. |
|------|----------|---------|------------------|------|
| 9.92 | 27.81 ng | 2316580 | Acenaphthene-d10 | 9.59 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#         | Qual |
|------|----|---|---------------------------------|-----|---------|--------------|------|
| 1    |    |   | Naphthalene, 1,4,6-trimethyl-   | 170 | C13H14  | 002131-42-2  | 97   |
| 2    |    |   | Naphthalene, 2,3,6-trimethyl-   | 170 | C13H14  | 000829-26-5  | 97   |
| 3    |    |   | Naphthalene, 1,6,7-trimethyl-   | 170 | C13H14  | 002245-38-7  | 96   |
| 4    |    |   | 4,6,8-Trimethylazulene          | 170 | C13H14  | 000941-81-1  | 94   |
| 5    |    |   | 3-(2-Methyl-propenyl)-1H-indene | 170 | C13H14  | 1000187-78-5 | 91   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

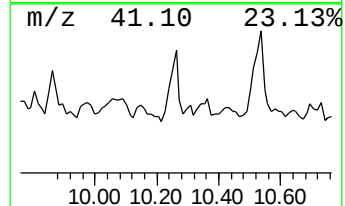
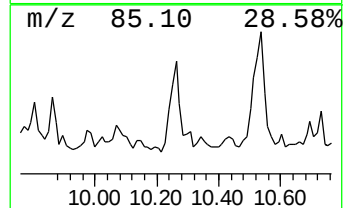
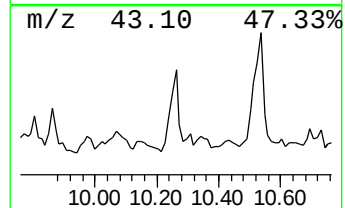
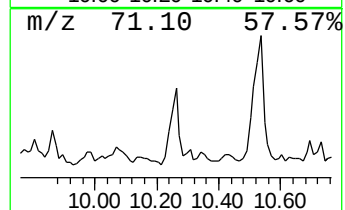
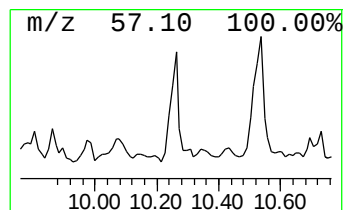
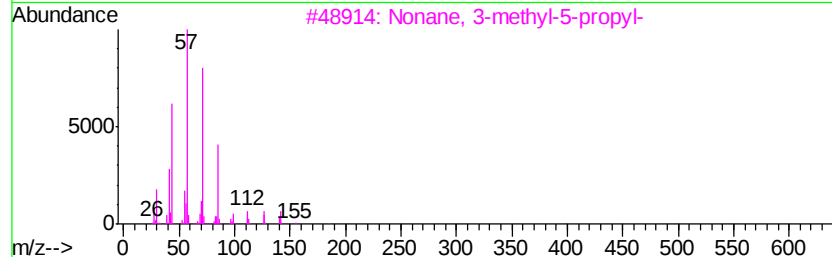
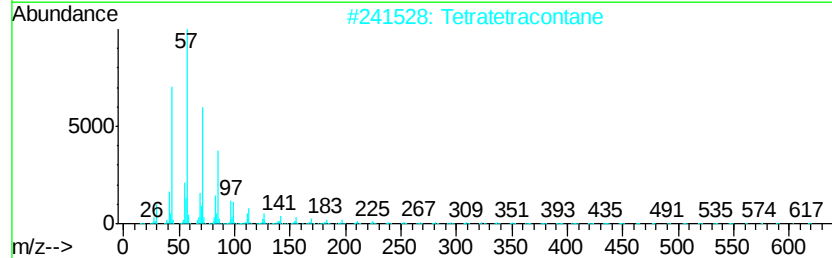
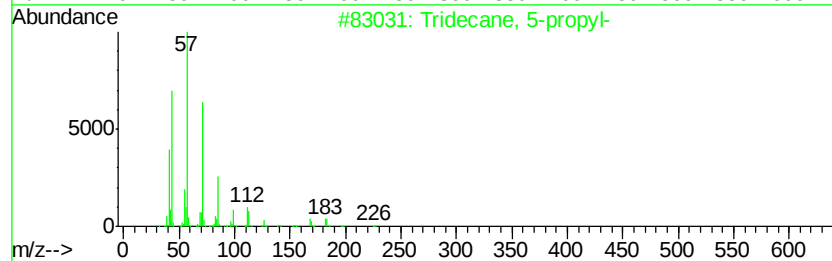
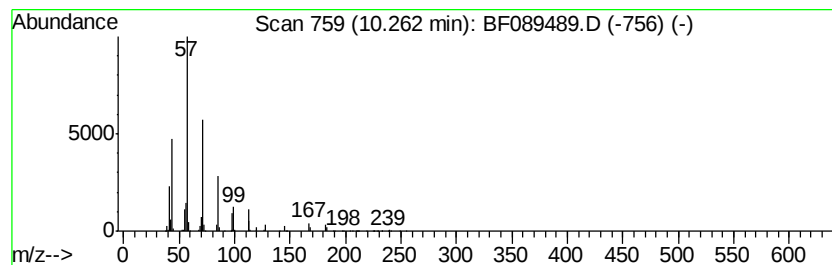
Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Tridecane, 5-propyl- Concentration Rank 6

| R.T.      | EstConc                    | Area    | Relative to ISTD |             | R.T. |
|-----------|----------------------------|---------|------------------|-------------|------|
| 10.26     | 60.04 ng                   | 5002110 | Acenaphthene-d10 |             | 9.59 |
| Hit# of 5 | Tentative ID               | MW      | MolForm          | CAS#        | Qual |
| 1         | Tridecane, 5-propyl-       | 226     | C16H34           | 055045-11-9 | 83   |
| 2         | Tetratetracontane          | 619     | C44H90           | 007098-22-8 | 64   |
| 3         | Nonane, 3-methyl-5-propyl- | 184     | C13H28           | 031081-18-2 | 59   |
| 4         | Nonane, 2-methyl-5-propyl- | 184     | C13H28           | 031081-17-1 | 59   |
| 5         | Dodecane, 4,9-dipropyl-    | 254     | C18H38           | 003054-63-5 | 58   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

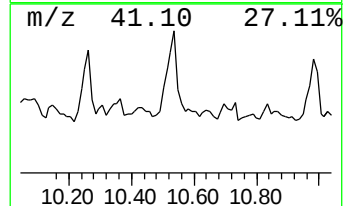
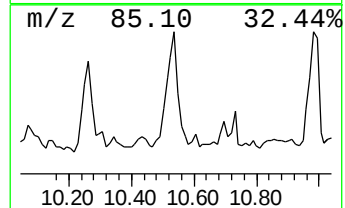
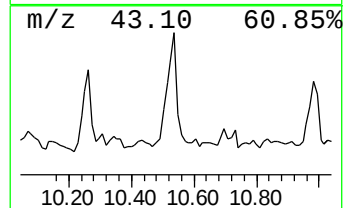
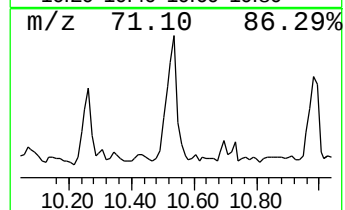
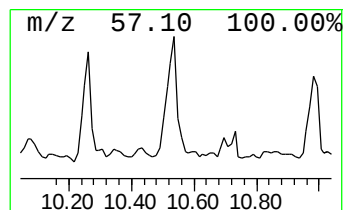
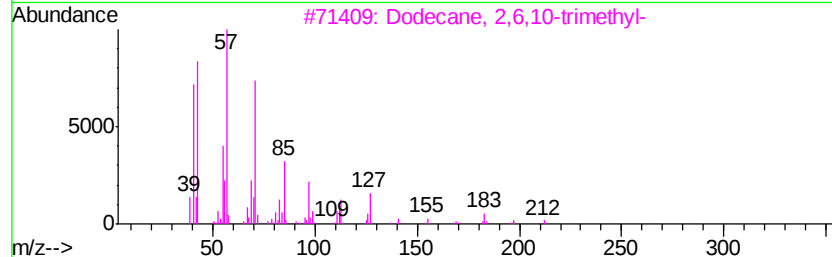
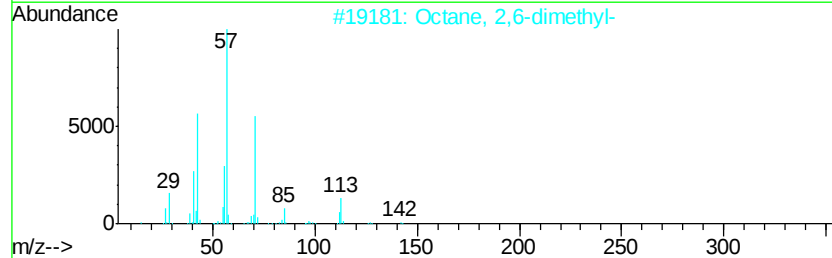
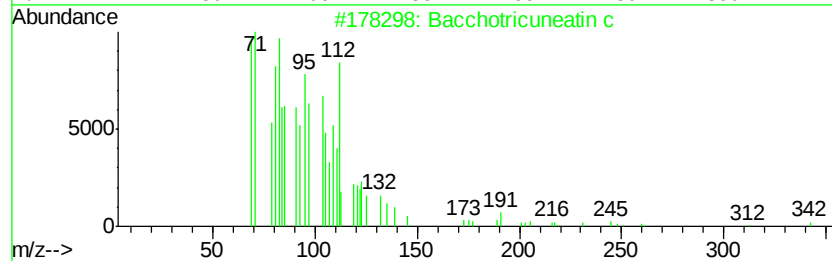
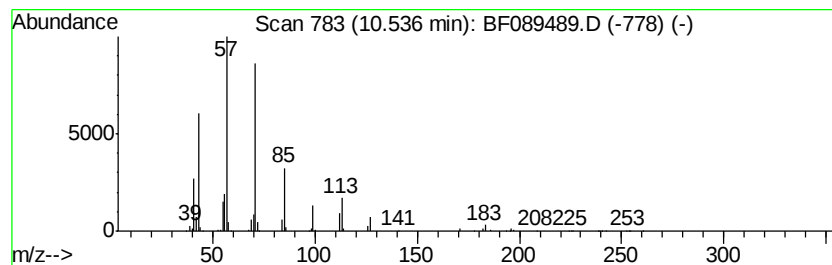
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Bacchotricuneatin c Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 10.54 | 72.71 ng | 4492920 | Phenanthrene-d10 | 11.05 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Bacchotricuneatin c                 | 342 | C20H22O5  | 066563-30-2  | 94   |
| 2    |    | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 80   |
| 3    |    | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32    | 003891-98-3  | 78   |
| 4    |    | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32    | 074645-98-0  | 72   |
| 5    |    | Sulfurous acid, 2-ethylhexyl tri... | 376 | C21H44O3S | 1000309-19-6 | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

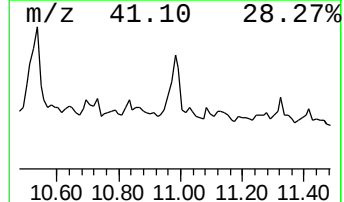
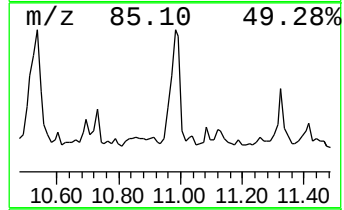
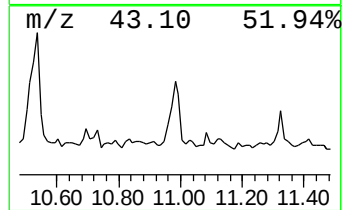
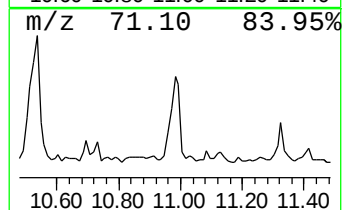
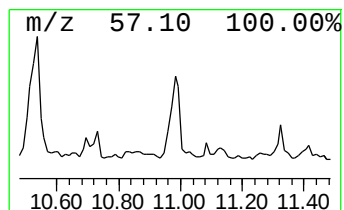
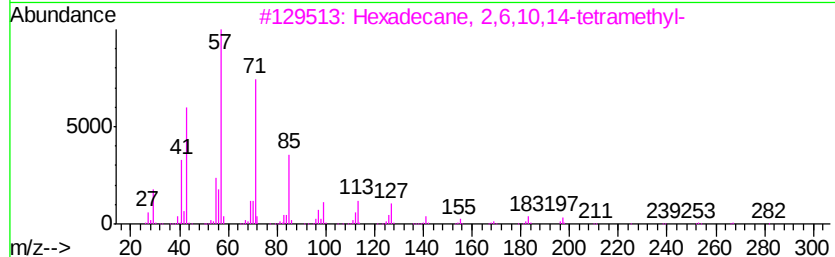
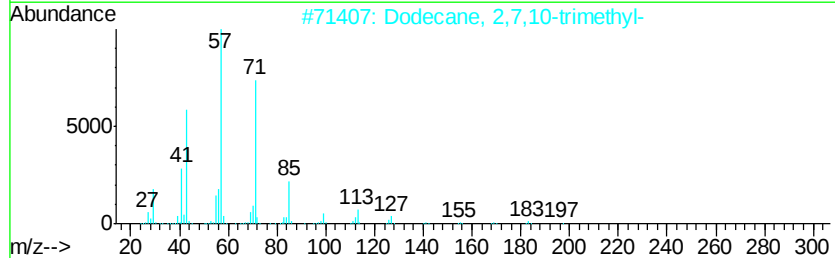
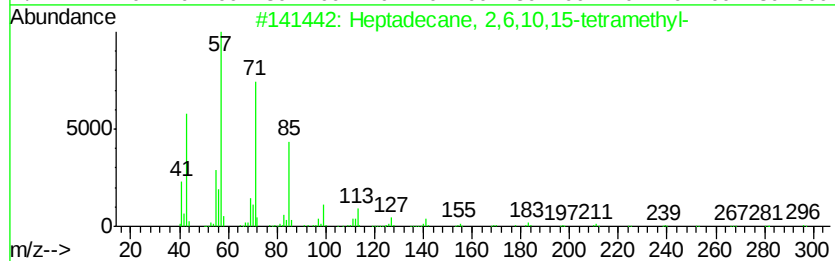
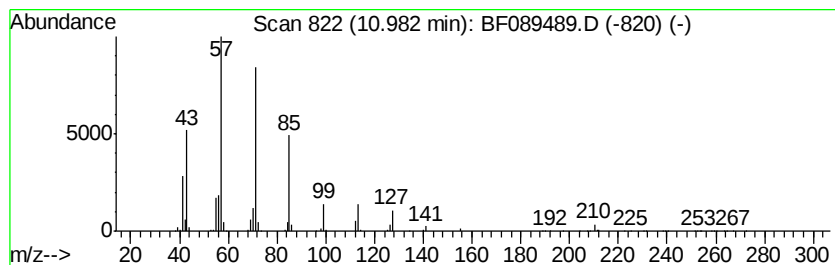
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 Heptadecane, 2,6,10,15-tetr... Concentration Rank 7

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 10.98 | 53.34 ng | 3295770 | Phenanthrene-d10 | 11.05 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 78   |
| 2    |    | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32  | 074645-98-0 | 72   |
| 3    |    | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 72   |
| 4    |    | Tetradecane, 4-methyl-              | 212 | C15H32  | 025117-24-2 | 64   |
| 5    |    | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32  | 003891-98-3 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

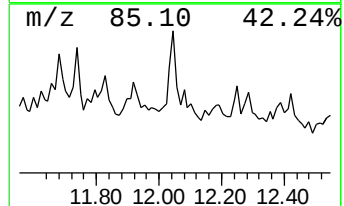
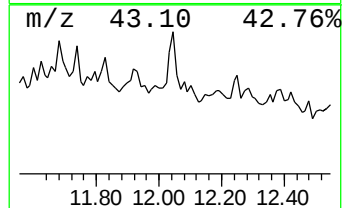
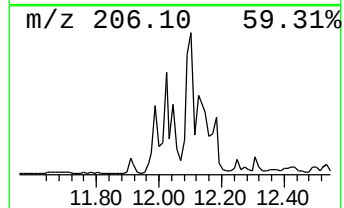
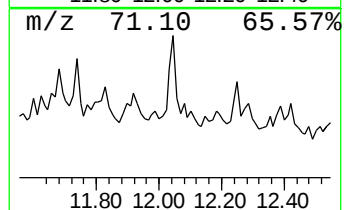
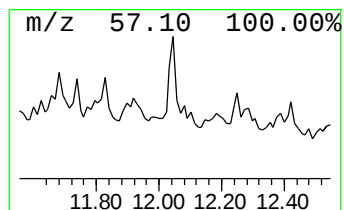
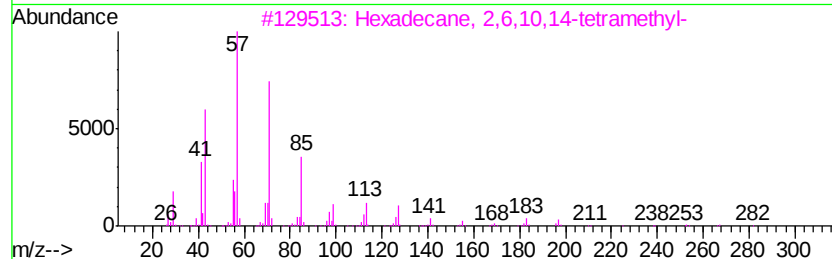
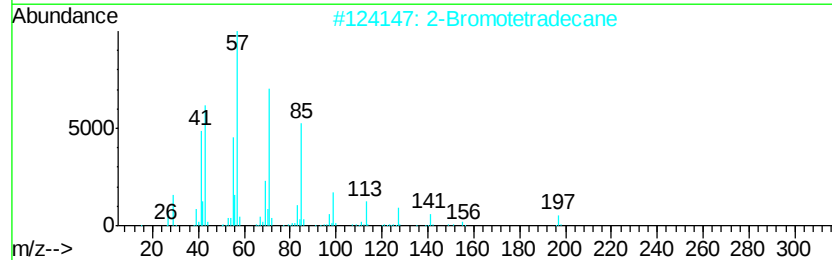
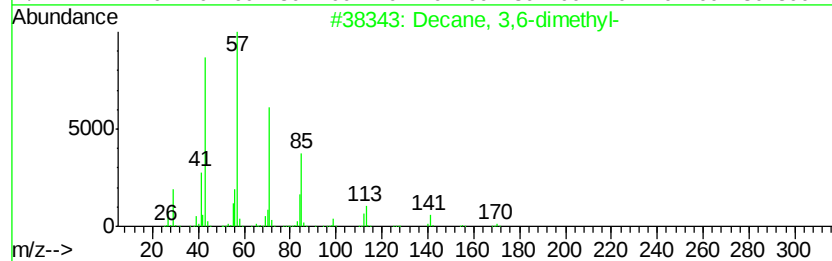
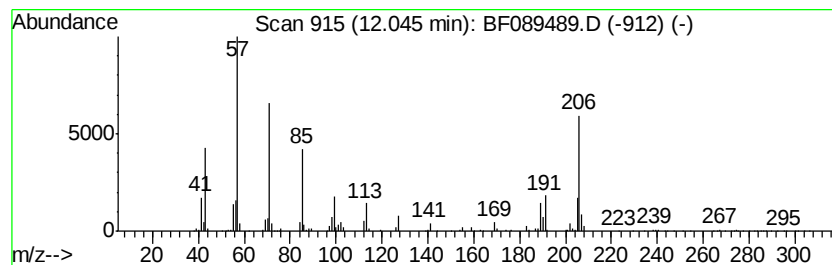
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Decane, 3,6-dimethyl- Concentration Rank 19

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.05 | 25.65 ng | 1584690 | Phenanthrene-d10 | 11.05 |

| Hit# | of | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Decane, 3,6-dimethyl-              | 170 | C12H26   | 017312-53-7 | 55   |
| 2    |    | 2-Bromotetradecane                 | 276 | C14H29Br | 074036-95-6 | 49   |
| 3    |    | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42   | 000638-36-8 | 49   |
| 4    |    | Hexadecane                         | 226 | C16H34   | 000544-76-3 | 49   |
| 5    |    | Octadecane, 1-iodo-                | 380 | C18H37I  | 000629-93-6 | 49   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

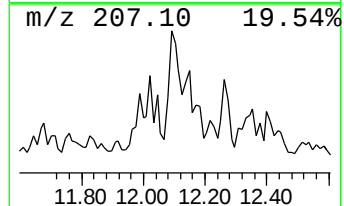
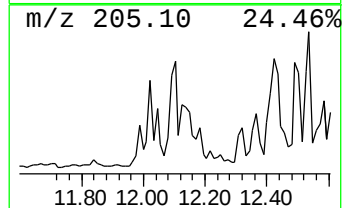
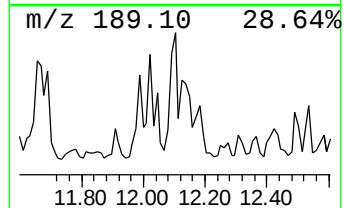
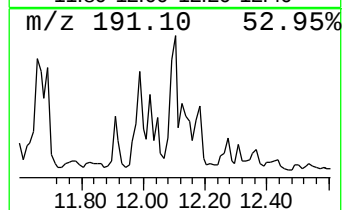
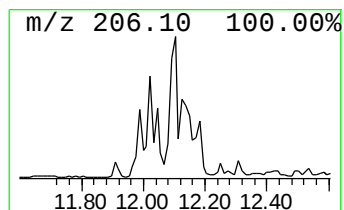
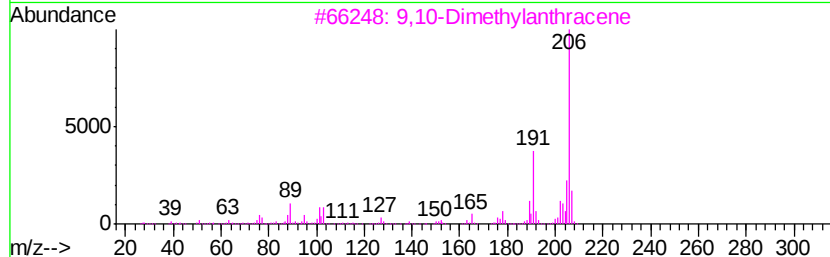
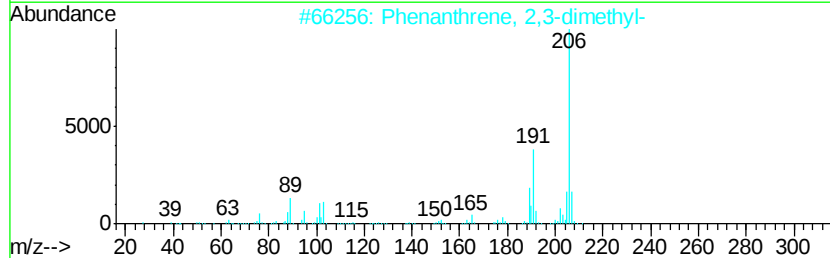
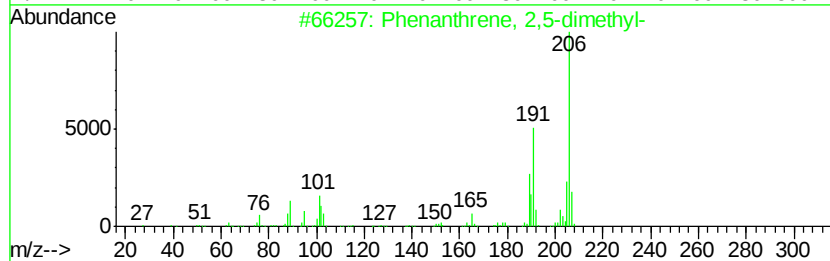
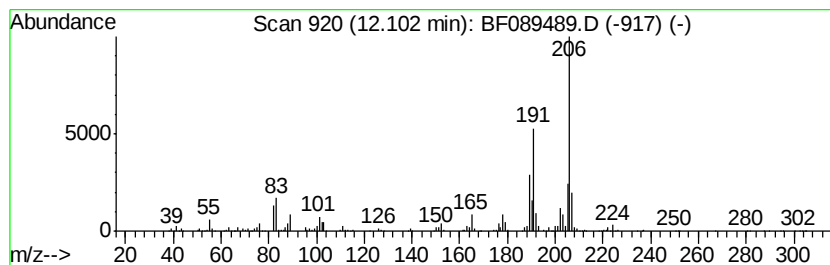
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Phenanthrene, 2,5-dimethyl- Concentration Rank 16

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.10 | 28.19 ng | 1742180 | Phenanthrene-d10 | 11.05 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 93   |
| 2    |    | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 93   |
| 3    |    | 9,10-Dimethylanthracene     | 206 | C16H14  | 000781-43-1 | 91   |
| 4    |    | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 91   |
| 5    |    | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 91   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

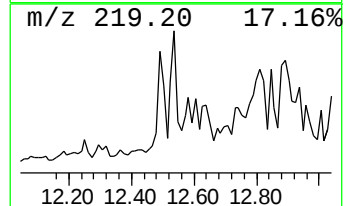
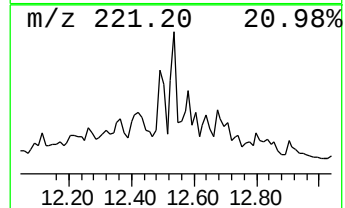
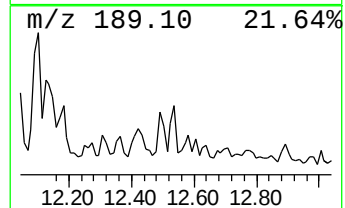
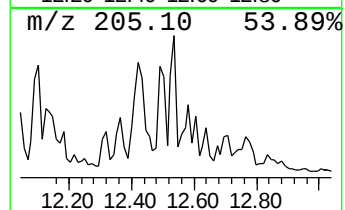
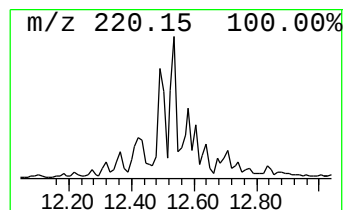
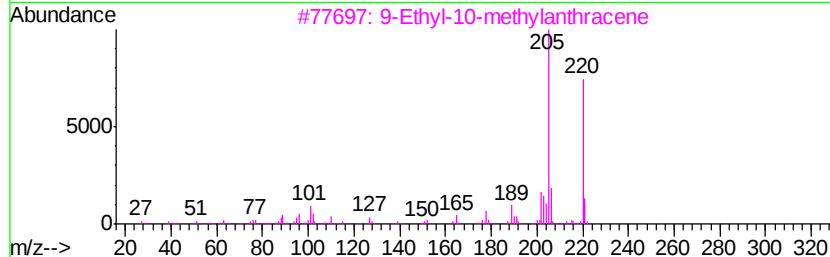
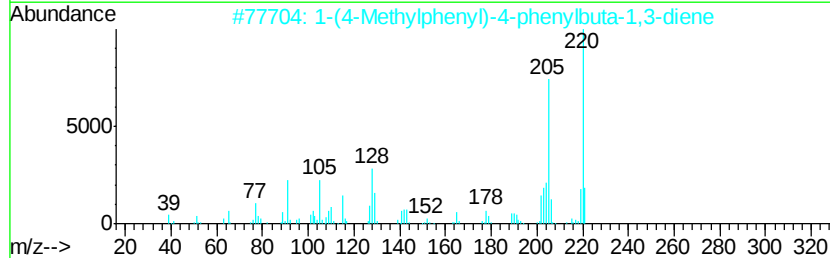
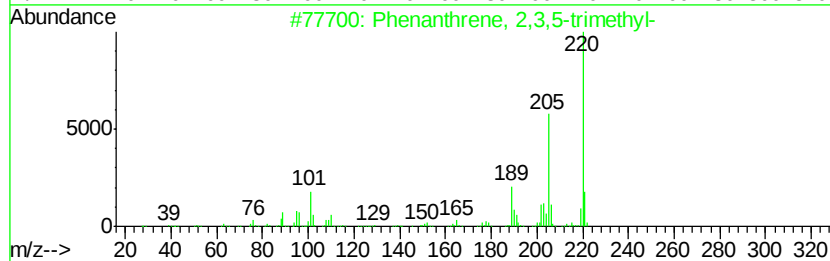
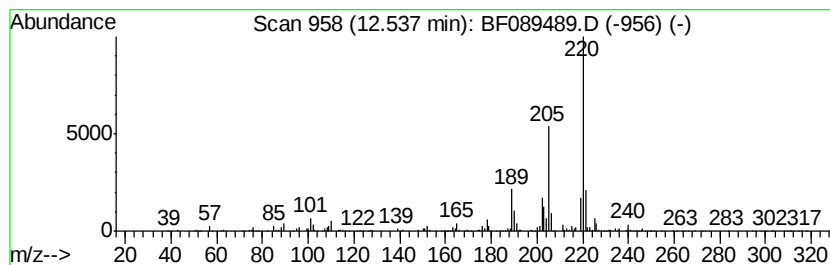
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Phenanthrene, 2,3,5-trimethyl- Concentration Rank 13

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.54 | 33.73 ng | 577168 | Chrysene-d12     | 13.63 |

| Hit# | of | Tentative ID                         | MW  | MolForm    | CAS#         | Qual |
|------|----|--------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Phenanthrene, 2,3,5-trimethyl-       | 220 | C17H16     | 003674-73-5  | 83   |
| 2    |    | 1-(4-Methylphenyl)-4-phenylbuta-...  | 220 | C17H16     | 037985-11-8  | 72   |
| 3    |    | 9-Ethyl-10-methylanthracene          | 220 | C17H16     | 019713-49-6  | 68   |
| 4    |    | Phenmethylnitrile, .alpha.,.alpha... | 220 | C11H12N2O3 | 1000128-94-6 | 53   |
| 5    |    | 10-Methylanthracene-9-carboxalde...  | 220 | C16H12O    | 007072-00-6  | 50   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\  
Data File : BF089489.D  
Acq On : 1 Aug 2016 2:03  
Operator : UM/SJ  
Sample : H4307-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

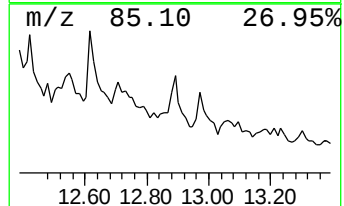
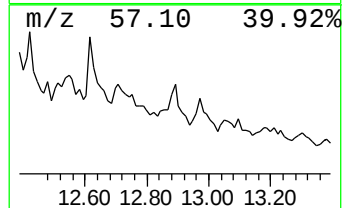
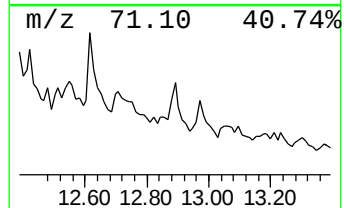
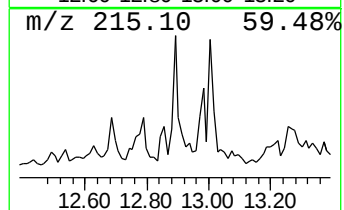
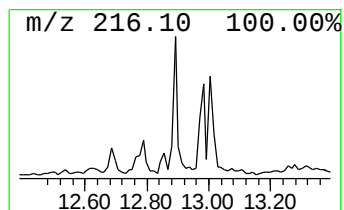
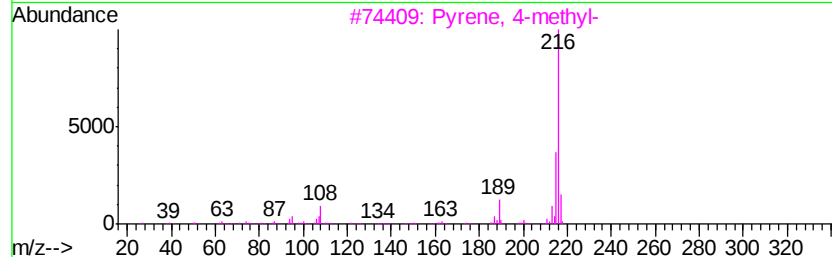
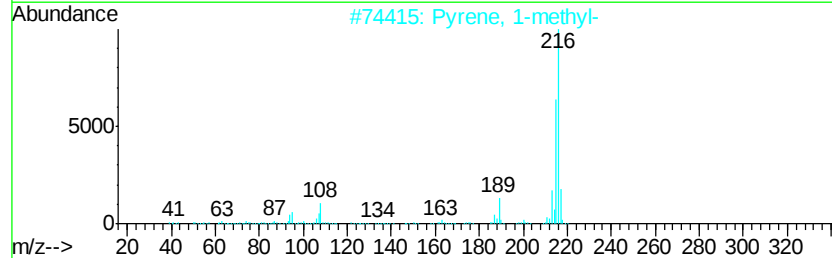
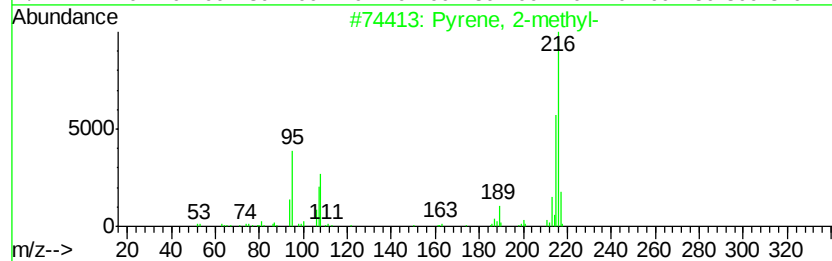
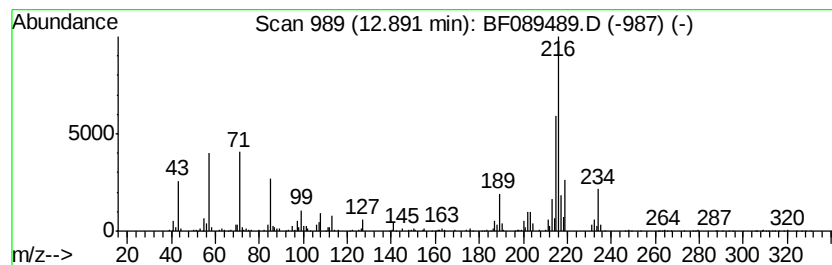
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Pyrene, 2-methyl- Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.89 | 60.34 ng | 1032310 | Chrysene-d12     | 13.63 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 70   |
| 2    |    | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 70   |
| 3    |    | Pyrene, 4-methyl-       | 216 | C17H12  | 003353-12-6 | 64   |
| 4    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 50   |
| 5    |    | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 50   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF073116\  
 Data File : BF089489.D  
 Acq On : 1 Aug 2016 2:03  
 Operator : UM/SJ  
 Sample : H4307-01  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 1057-SOUTH

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072716.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Nonane, 3-methyl-    | 5.74  | 39.9    | ng    | 3221930  | 1                     | 6.50  | 1617170 | 20.0 |
| Heptane, 3-ethyl-... | 5.79  | 31.1    | ng    | 2517770  | 1                     | 6.50  | 1617170 | 20.0 |
| Nonane, 4-methyl-    | 6.01  | 43.3    | ng    | 3497380  | 1                     | 6.50  | 1617170 | 20.0 |
| unknown6.27          | 6.27  | 52.7    | ng    | 4263310  | 1                     | 6.50  | 1617170 | 20.0 |
| Benzene, 1,2,3-tr... | 6.34  | 75.6    | ng    | 6109790  | 1                     | 6.50  | 1617170 | 20.0 |
| unknown6.68          | 6.68  | 65.9    | ng    | 5328700  | 1                     | 6.50  | 1617170 | 20.0 |
| 2-Tolyloxirane       | 6.81  | 45.7    | ng    | 3696270  | 1                     | 6.50  | 1617170 | 20.0 |
| Naphthalene, deca... | 6.91  | 34.7    | ng    | 2802430  | 1                     | 6.50  | 1617170 | 20.0 |
| Benzene, 4-ethyl-... | 7.02  | 29.8    | ng    | 2412570  | 1                     | 6.50  | 1617170 | 20.0 |
| Naphthalene, 1,3-... | 9.27  | 75.3    | ng    | 6274330  | 3                     | 9.59  | 1666180 | 20.0 |
| Naphthalene, 2,3,... | 9.84  | 27.8    | ng    | 2319610  | 3                     | 9.59  | 1666180 | 20.0 |
| Naphthalene, 1,4,... | 9.92  | 27.8    | ng    | 2316580  | 3                     | 9.59  | 1666180 | 20.0 |
| Tridecane, 5-propyl- | 10.26 | 60.0    | ng    | 5002110  | 3                     | 9.59  | 1666180 | 20.0 |
| Bacchotricuneatin c  | 10.54 | 72.7    | ng    | 4492920  | 4                     | 11.05 | 1235820 | 20.0 |
| Heptadecane, 2,6,... | 10.98 | 53.3    | ng    | 3295770  | 4                     | 11.05 | 1235820 | 20.0 |
| Decane, 3,6-dimet... | 12.05 | 25.6    | ng    | 1584690  | 4                     | 11.05 | 1235820 | 20.0 |
| Phenanthrene, 2,5... | 12.10 | 28.2    | ng    | 1742180  | 4                     | 11.05 | 1235820 | 20.0 |
| Phenanthrene, 2,3... | 12.54 | 33.7    | ng    | 577168   | 5                     | 13.63 | 342192  | 20.0 |
| Pyrene, 2-methyl-    | 12.89 | 60.3    | ng    | 1032310  | 5                     | 13.63 | 342192  | 20.0 |