

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
 Data File : BF091558.D  
 Acq On : 13 Dec 2016 17:13  
 Operator : UM/SJ  
 Sample : H5922-01 5X  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 WC120716

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-BF120916.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|--------|----------|------------|----------|-----------|-------|-------------|------------|--------------|------------|
| 1      | 4.967    | 250        | 252      | 255       | rBV   | 589362      | 618282     | 9.38%        | 0.687%     |
| 2      | 5.402    | 288        | 290      | 293       | rBV   | 1199028     | 1519221    | 23.05%       | 1.688%     |
| 3      | 6.442    | 379        | 381      | 384       | rBV   | 1589068     | 1533560    | 23.27%       | 1.704%     |
| 4      | 6.579    | 391        | 393      | 395       | rBV   | 2029519     | 1692422    | 25.68%       | 1.880%     |
| 5      | 6.807    | 411        | 413      | 416       | rBV   | 1680525     | 1734460    | 26.32%       | 1.927%     |
| 6      | 6.967    | 425        | 427      | 429       | rBV   | 1559824     | 1336551    | 20.28%       | 1.485%     |
| 7      | 7.367    | 460        | 462      | 465       | rVB   | 997258      | 938026     | 14.23%       | 1.042%     |
| 8      | 8.099    | 523        | 526      | 528       | rBV   | 2689897     | 2447676    | 37.14%       | 2.719%     |
| 9      | 8.488    | 558        | 560      | 561       | rBV   | 85673       | 97794      | 1.48%        | 0.109%     |
| 10     | 8.522    | 561        | 563      | 565       | rBV2  | 49323       | 85520      | 1.30%        | 0.095%     |
| 11     | 8.590    | 566        | 569      | 571       | rVV2  | 49938       | 101749     | 1.54%        | 0.113%     |
| 12     | 8.625    | 571        | 572      | 573       | rBV   | 151613      | 133463     | 2.03%        | 0.148%     |
| 13     | 8.693    | 575        | 578      | 582       | rVV3  | 151946      | 454426     | 6.90%        | 0.505%     |
| 14     | 8.785    | 584        | 586      | 587       | rBV   | 73771       | 105132     | 1.60%        | 0.117%     |
| 15     | 8.865    | 590        | 593      | 595       | rBV2  | 85351       | 153652     | 2.33%        | 0.171%     |
| 16     | 8.922    | 595        | 598      | 599       | rBV3  | 146863      | 208329     | 3.16%        | 0.231%     |
| 17     | 9.025    | 604        | 607      | 608       | rVV2  | 187273      | 375973     | 5.70%        | 0.418%     |
| 18     | 9.082    | 608        | 612      | 615       | rVV   | 433566      | 1175544    | 17.84%       | 1.306%     |
| 19     | 9.173    | 618        | 620      | 622       | rVV   | 2335492     | 2219677    | 33.68%       | 2.466%     |
| 20     | 9.208    | 622        | 623      | 626       | rVV2  | 273284      | 529779     | 8.04%        | 0.589%     |
| 21     | 9.288    | 629        | 630      | 632       | rVV2  | 274792      | 440690     | 6.69%        | 0.490%     |
| 22     | 9.345    | 633        | 635      | 636       | rVV2  | 250464      | 407555     | 6.18%        | 0.453%     |
| 23     | 9.368    | 636        | 637      | 639       | rVV2  | 450000      | 648711     | 9.84%        | 0.721%     |
| 24     | 9.425    | 640        | 642      | 645       | rVV   | 287941      | 536171     | 8.14%        | 0.596%     |
| 25     | 9.528    | 649        | 651      | 652       | rVV   | 132207      | 183358     | 2.78%        | 0.204%     |
| 26     | 9.573    | 652        | 655      | 661       | rVB4  | 405007      | 965915     | 14.66%       | 1.073%     |
| 27     | 9.791    | 672        | 674      | 675       | rVB   | 123779      | 143177     | 2.17%        | 0.159%     |
| 28     | 9.848    | 677        | 679      | 683       | rVV   | 2579783     | 2385026    | 36.19%       | 2.649%     |
| 29     | 9.928    | 685        | 686      | 688       | rVV2  | 171749      | 227155     | 3.45%        | 0.252%     |
| 30     | 10.099   | 699        | 701      | 702       | rVB   | 177069      | 241602     | 3.67%        | 0.268%     |
| 31     | 10.202   | 709        | 710      | 711       | rBV   | 169509      | 122786     | 1.86%        | 0.136%     |
| 32     | 10.351   | 721        | 723      | 724       | rVB2  | 419999      | 542797     | 8.24%        | 0.603%     |
| 33     | 10.396   | 724        | 727      | 730       | rBV3  | 531891      | 1110754    | 16.85%       | 1.234%     |
| 34     | 10.522   | 735        | 738      | 740       | rBV3  | 693059      | 1469763    | 22.30%       | 1.633%     |

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
 Data File : BF091558.D  
 Acq On : 13 Dec 2016 17:13  
 Operator : UM/SJ  
 Sample : H5922-01 5X  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 WC120716

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-BF120916.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 35 | 10.556 | 740  | 741  | 742  | rVV  | 696953  | 568072  | 8.62%   | 0.631% |
| 36 | 10.602 | 742  | 745  | 746  | rVV  | 582309  | 1088084 | 16.51%  | 1.209% |
| 37 | 10.636 | 746  | 748  | 751  | rVB2 | 697522  | 1425265 | 21.63%  | 1.583% |
| 38 | 10.716 | 753  | 755  | 757  | rVV  | 1684306 | 1752800 | 26.60%  | 1.947% |
| 39 | 10.751 | 757  | 758  | 760  | rVB  | 550100  | 494195  | 7.50%   | 0.549% |
| 40 | 10.808 | 760  | 763  | 766  | rBV2 | 1278676 | 2534370 | 38.45%  | 2.815% |
| 41 | 10.865 | 766  | 768  | 770  | rBV  | 1674672 | 1459724 | 22.15%  | 1.622% |
| 42 | 10.911 | 770  | 772  | 774  | rBV2 | 785328  | 939067  | 14.25%  | 1.043% |
| 43 | 11.002 | 777  | 780  | 785  | rVB2 | 629987  | 1525137 | 23.14%  | 1.694% |
| 44 | 11.082 | 785  | 787  | 789  | rVB2 | 319701  | 419960  | 6.37%   | 0.467% |
| 45 | 11.116 | 789  | 790  | 791  | rVB  | 434845  | 298086  | 4.52%   | 0.331% |
| 46 | 11.151 | 791  | 793  | 794  | rBV  | 603989  | 846658  | 12.85%  | 0.941% |
| 47 | 11.208 | 794  | 798  | 799  | rBV  | 1291187 | 2387639 | 36.23%  | 2.652% |
| 48 | 11.288 | 803  | 805  | 807  | rBV  | 1332168 | 2014829 | 30.57%  | 2.238% |
| 49 | 11.334 | 807  | 809  | 810  | rBV  | 1729777 | 2332698 | 35.39%  | 2.591% |
| 50 | 11.391 | 812  | 814  | 816  | rVV  | 1746439 | 2727826 | 41.39%  | 3.030% |
| 51 | 11.436 | 816  | 818  | 819  | rVV  | 1261994 | 2001240 | 30.37%  | 2.223% |
| 52 | 11.459 | 819  | 820  | 822  | rVV2 | 1201791 | 1697408 | 25.76%  | 1.886% |
| 53 | 11.528 | 824  | 826  | 827  | rVV2 | 1447335 | 1865808 | 28.31%  | 2.073% |
| 54 | 11.562 | 827  | 829  | 831  | rVV  | 2278571 | 3481426 | 52.82%  | 3.867% |
| 55 | 11.608 | 831  | 833  | 834  | rVV2 | 1405114 | 2139626 | 32.47%  | 2.377% |
| 56 | 11.654 | 834  | 837  | 839  | rVV  | 3328865 | 5886634 | 89.32%  | 6.539% |
| 57 | 11.699 | 839  | 841  | 842  | rVV  | 1504204 | 2149917 | 32.62%  | 2.388% |
| 58 | 11.745 | 842  | 845  | 848  | rVV3 | 2197994 | 6590495 | 100.00% | 7.321% |
| 59 | 11.791 | 848  | 849  | 850  | rVV  | 1773214 | 1808521 | 27.44%  | 2.009% |
| 60 | 11.825 | 850  | 852  | 856  | rVV  | 1928739 | 4477281 | 67.94%  | 4.974% |
| 61 | 11.894 | 856  | 858  | 861  | rVB2 | 888005  | 1489636 | 22.60%  | 1.655% |
| 62 | 12.145 | 879  | 880  | 881  | rBV  | 777444  | 776209  | 11.78%  | 0.862% |
| 63 | 12.911 | 946  | 947  | 950  | rVB  | 1167485 | 1272561 | 19.31%  | 1.414% |
| 64 | 13.962 | 1037 | 1039 | 1043 | rVB  | 1611906 | 2075245 | 31.49%  | 2.305% |
| 65 | 14.980 | 1126 | 1128 | 1131 | rVB  | 263362  | 438051  | 6.65%   | 0.487% |
| 66 | 15.380 | 1160 | 1163 | 1169 | rVB  | 1188259 | 1785992 | 27.10%  | 1.984% |
| 67 | 16.454 | 1255 | 1257 | 1261 | rVB2 | 194183  | 384491  | 5.83%   | 0.427% |

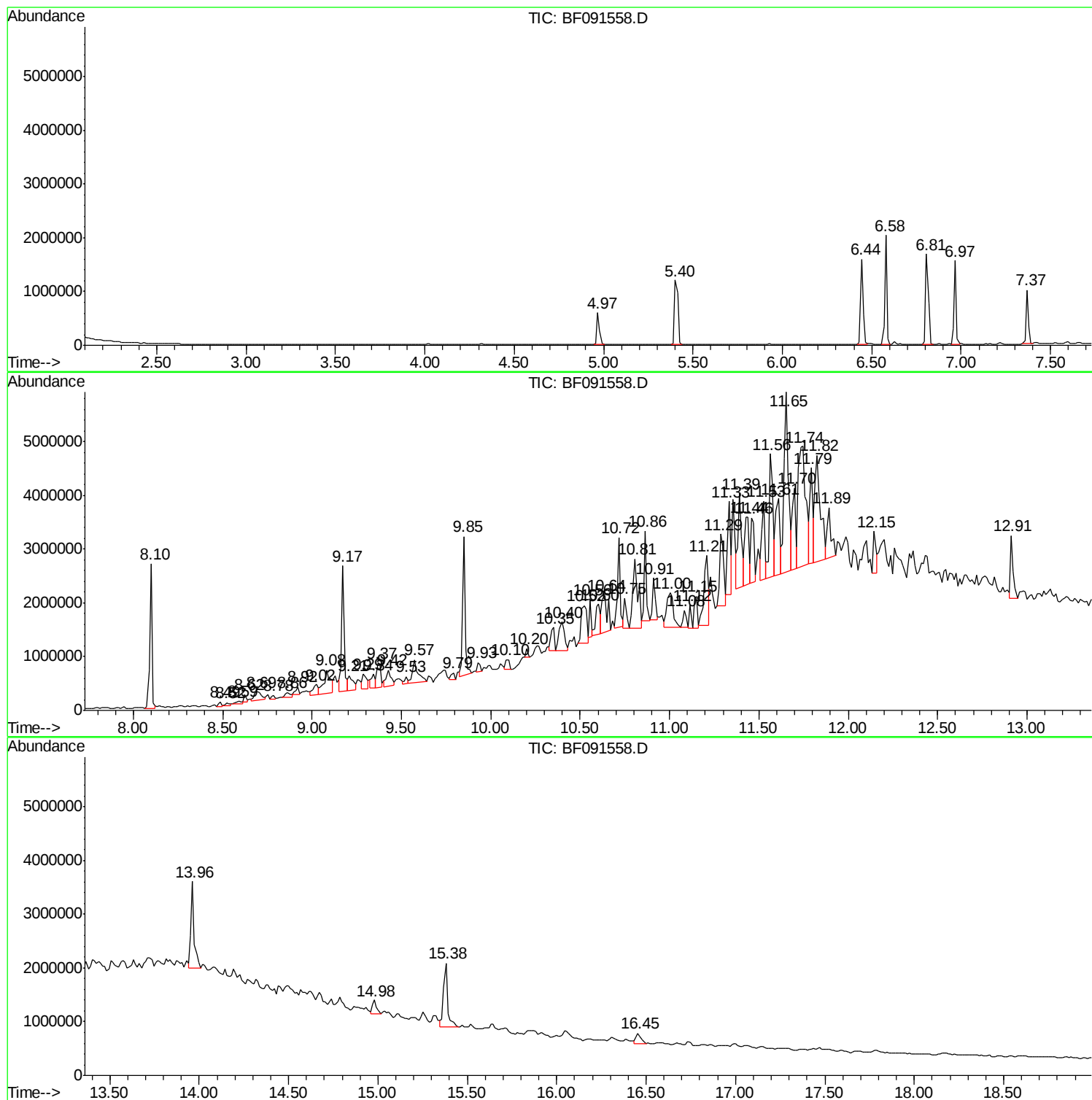
Sum of corrected areas: 90021647

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

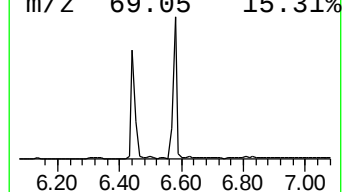
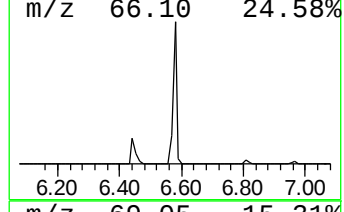
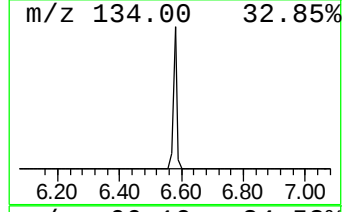
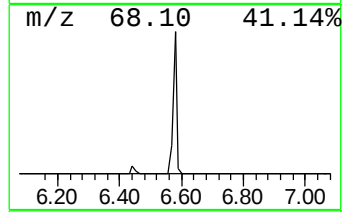
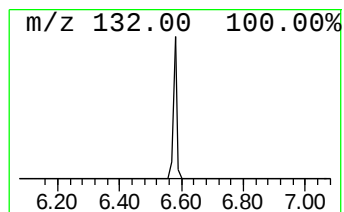
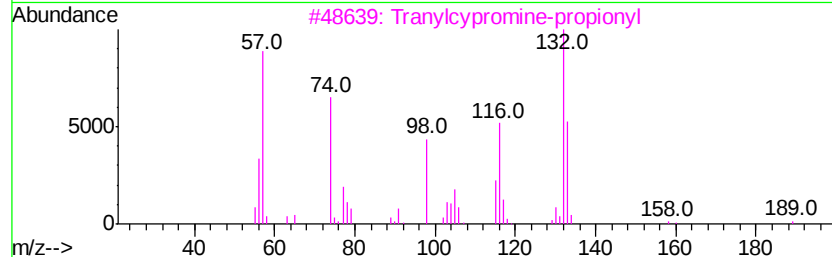
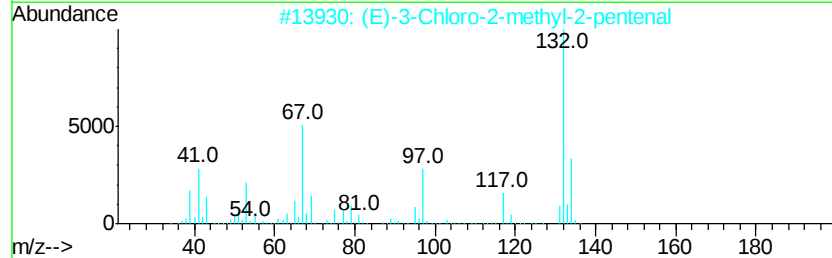
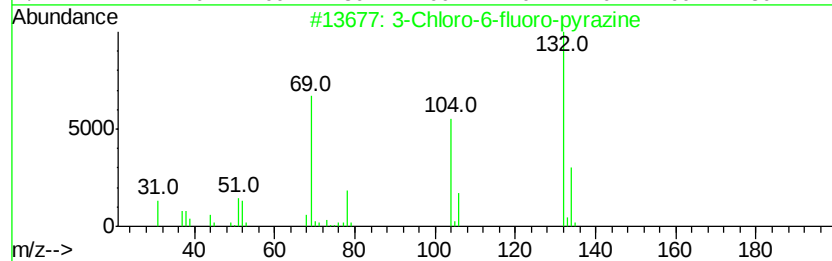
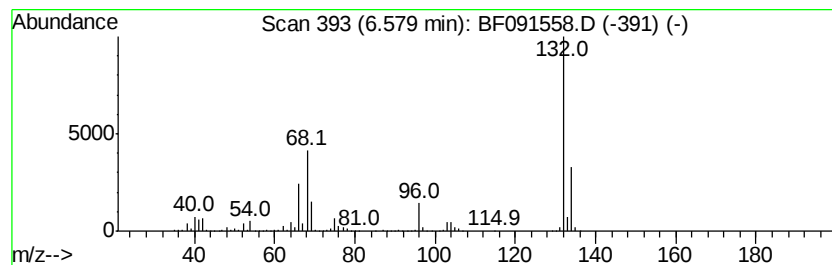
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 unknown6.58 Concentration Rank 7

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.58 | 19.52 ng | 1692420 | 1,4-Dichlorobenzene-d4 | 6.81 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 16   |
| 3    |    | Tranvlcyproline-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 4    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    | 5-Fluoro-2-chloropyrimidine         | 132 | C4H2ClFN2 | 062802-42-0  | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

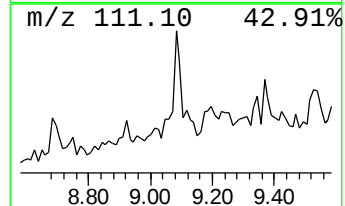
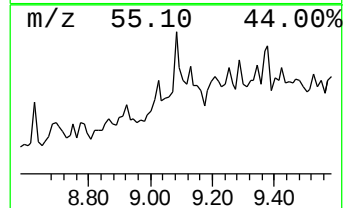
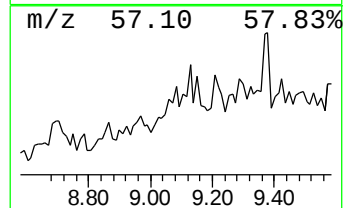
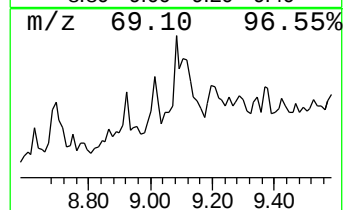
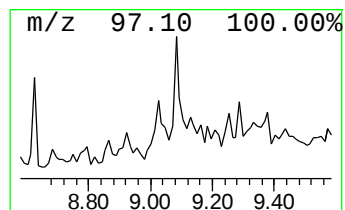
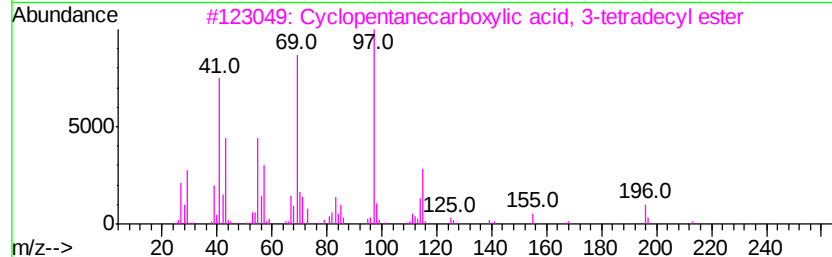
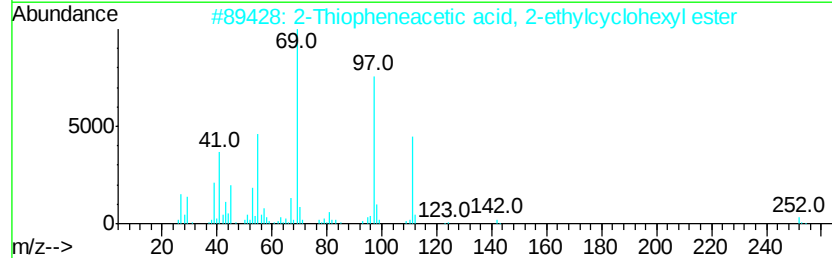
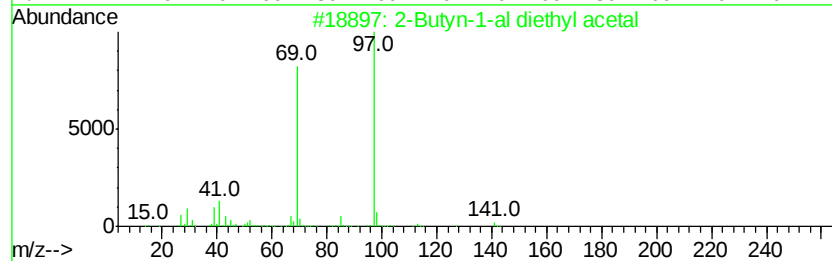
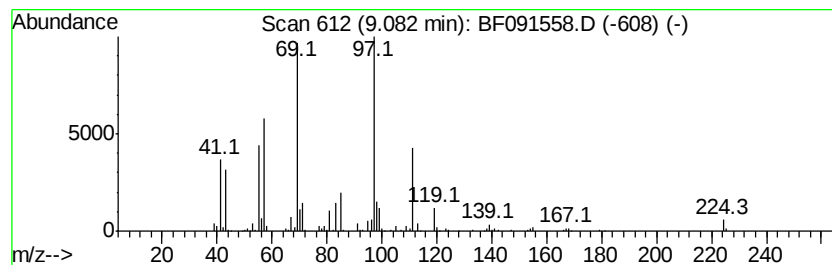
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 2-Butyn-1-al diethyl acetal Concentration Rank 20

| R.T. | EstConc | Area    | Relative to ISTD | R.T. |
|------|---------|---------|------------------|------|
| 9.08 | 9.86 ng | 1175540 | Acenaphthene-d10 | 9.85 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 2-Butyn-1-al diethyl acetal         | 142 | C8H14O2   | 002806-97-5  | 50   |
| 2    |    | 2-Thiopheneacetic acid, 2-ethylc... | 252 | C14H20O2S | 1000278-96-1 | 50   |
| 3    |    | Cyclopentanecarboxylic acid, 3-t... | 310 | C20H38O2  | 1000280-58-8 | 42   |
| 4    |    | Cyclopentanecarboxylic acid, 4-t... | 296 | C19H36O2  | 1000280-58-6 | 42   |
| 5    |    | Cyclopentanecarboxylic acid, 4-h... | 338 | C22H42O2  | 1000282-77-5 | 40   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

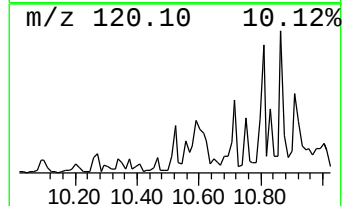
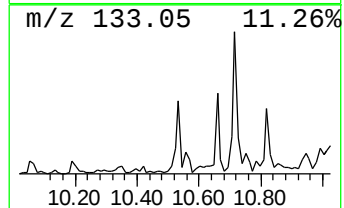
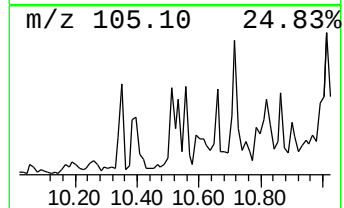
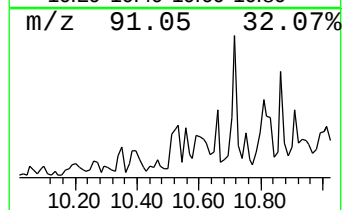
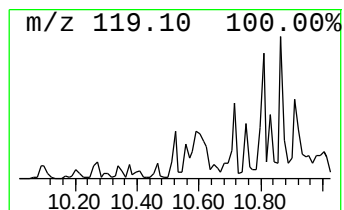
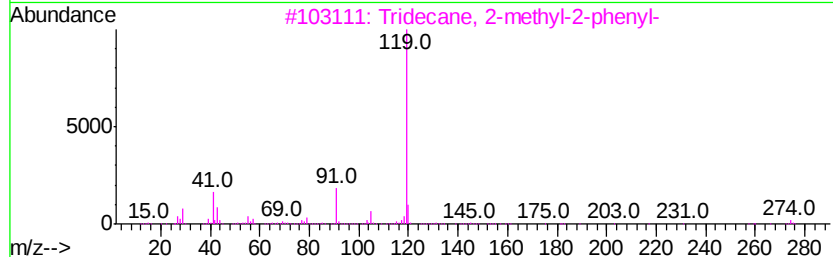
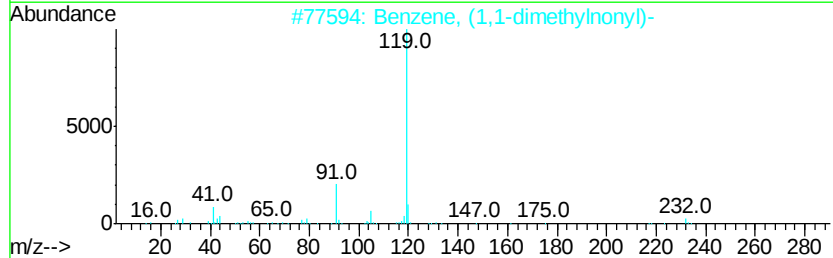
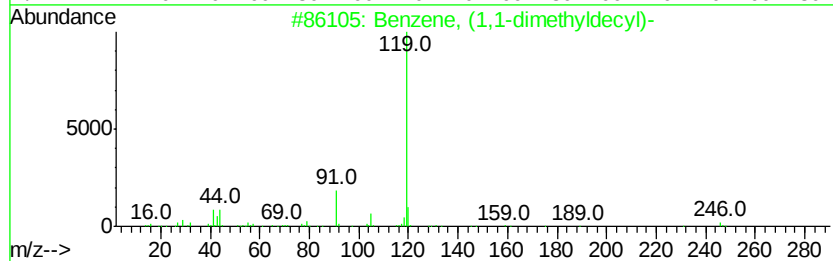
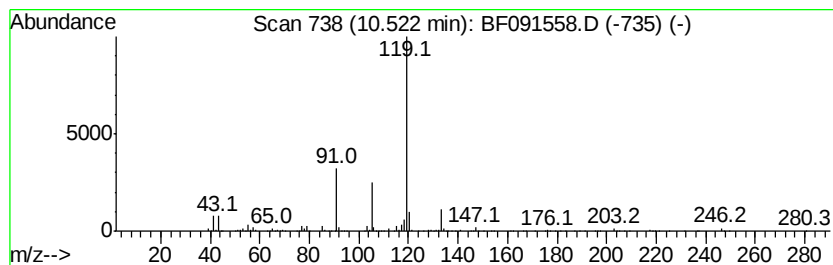
Instrument :  
BNA\_F  
ClientSampled :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Benzene, (1,1-dimethyldecyl)- Concentration Rank 19

| R.T.    | EstConc  | Area                                | Relative to ISTD | R.T.            |
|---------|----------|-------------------------------------|------------------|-----------------|
| 10.52   | 12.32 ng | 1469760                             | Acenaphthene-d10 | 9.85            |
| Hit# of | 5        | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |          | Benzene, (1,1-dimethyldecyl)-       | 246 C18H30       | 027854-40-6 64  |
| 2       |          | Benzene, (1,1-dimethylnonyl)-       | 232 C17H28       | 055191-25-8 53  |
| 3       |          | Tridecane, 2-methyl-2-phenyl-       | 274 C20H34       | 027854-41-7 53  |
| 4       |          | m-Toluylic acid, oct-3-en-2-yl e... | 246 C16H22O2     | 1000292-61-2 53 |
| 5       |          | p-Menthane, 2,3-dibromo-8-phenyl-   | 372 C16H22Br2    | 1000152-61-6 50 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

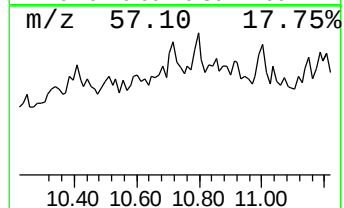
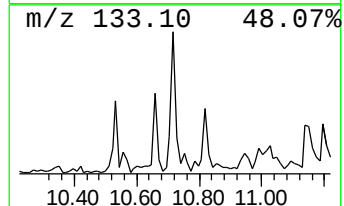
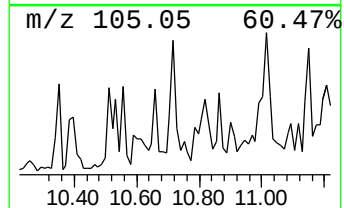
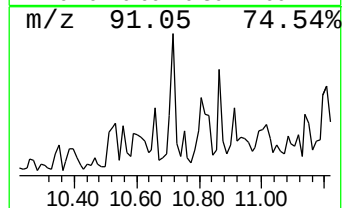
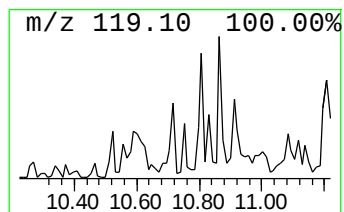
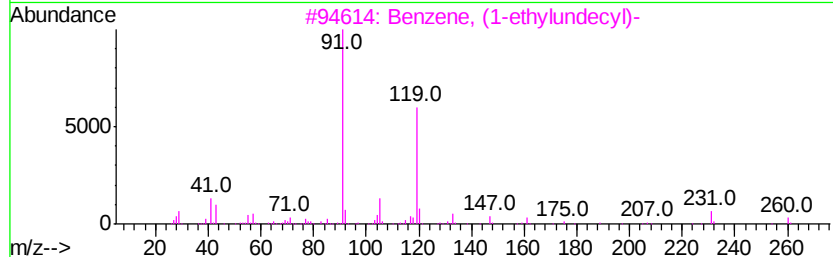
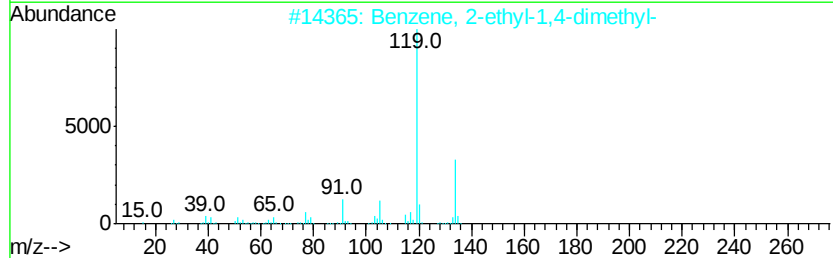
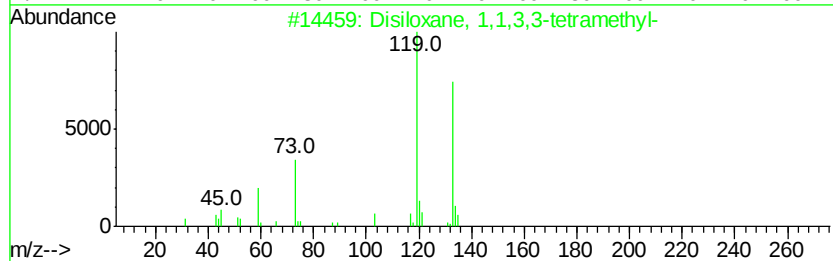
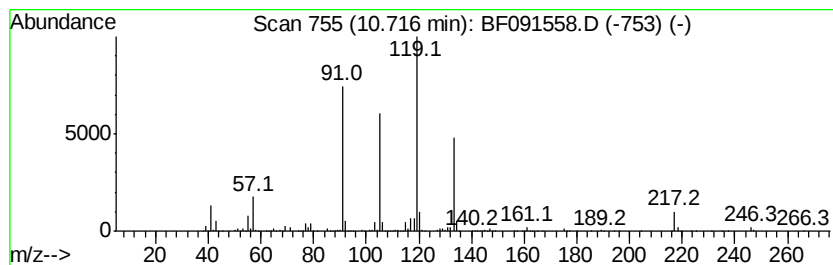
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Disiloxane, 1,1,3,3-tetrame... Concentration Rank 14

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 10.72 | 15.03 ng | 1752800 | Phenanthrene-d10 | 11.33 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|-----------|-------------------------------------|-----|-----------|-------------|------|
| 1         | Disiloxane, 1,1,3,3-tetramethyl-    | 134 | C4H14OSi2 | 003277-26-7 | 59   |
| 2         | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14    | 001758-88-9 | 53   |
| 3         | Benzene, (1-ethylundecyl)-          | 260 | C19H32    | 004534-52-5 | 53   |
| 4         | Cyclohexanone, 5-methyl-2-(1-met... | 230 | C16H22O   | 097275-48-4 | 50   |
| 5         | Benzene, (1,1-dimethyldecyl)-       | 246 | C18H30    | 027854-40-6 | 49   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

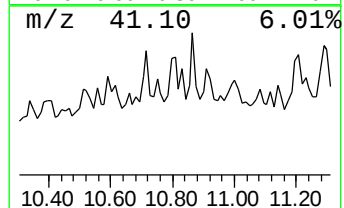
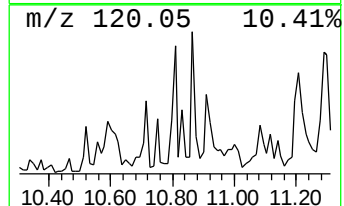
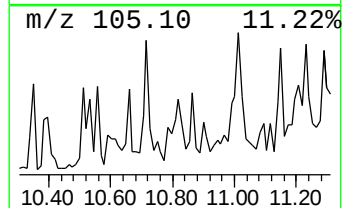
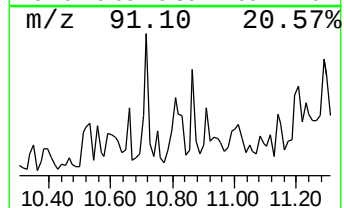
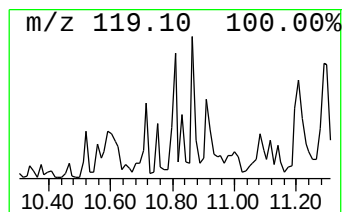
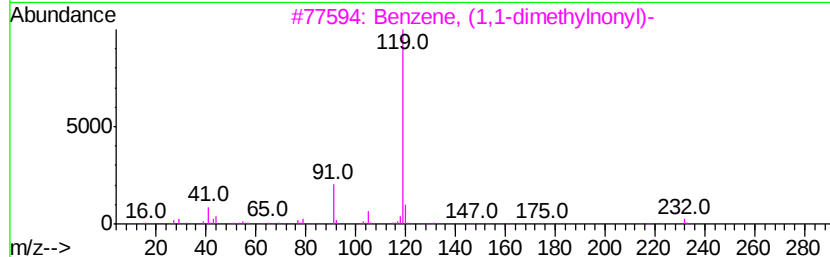
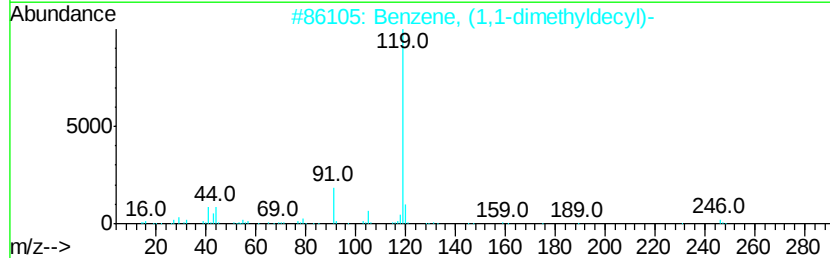
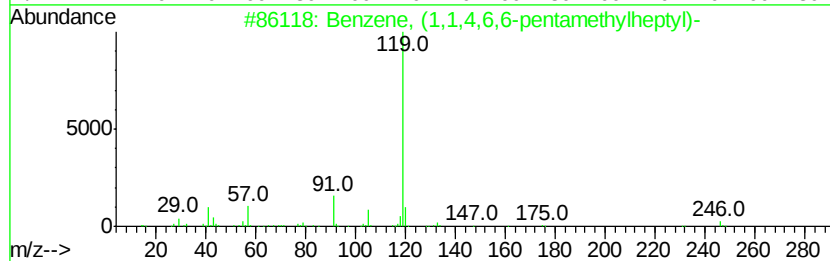
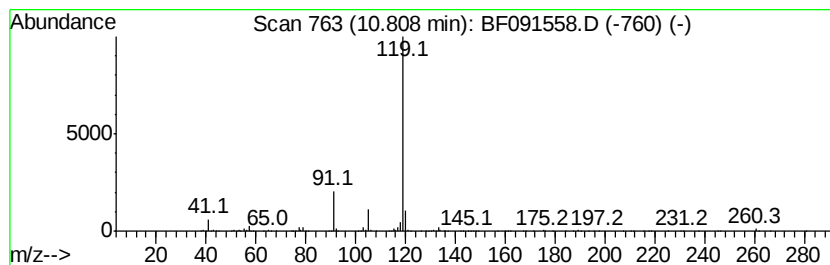
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Benzene, (1,1,4,6,6-pentame... Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 10.81 | 21.73 ng | 2534370 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (1,1,4,6,6-pentamethylh... | 246 | C18H30  | 055134-07-1 | 83   |
| 2    |    | Benzene, (1,1-dimethyldecyl)-       | 246 | C18H30  | 027854-40-6 | 83   |
| 3    |    | Benzene, (1,1-dimethylnonyl)-       | 232 | C17H28  | 055191-25-8 | 83   |
| 4    |    | Pentadecane, 2-methyl-2-phenyl-     | 302 | C22H38  | 029138-94-1 | 83   |
| 5    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

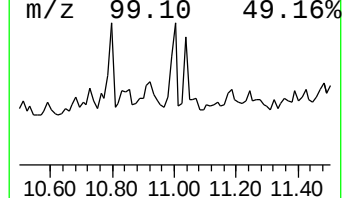
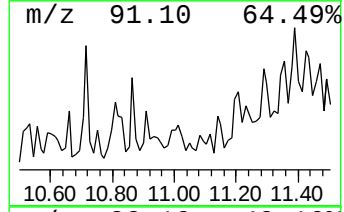
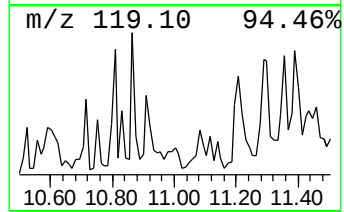
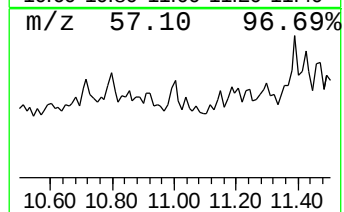
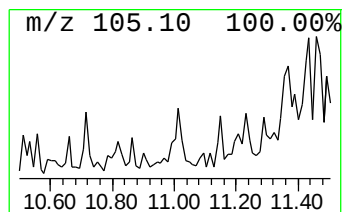
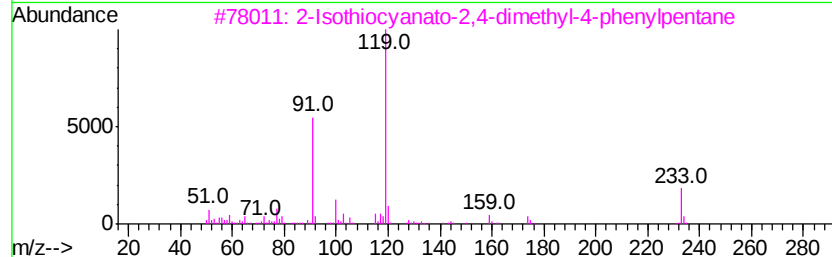
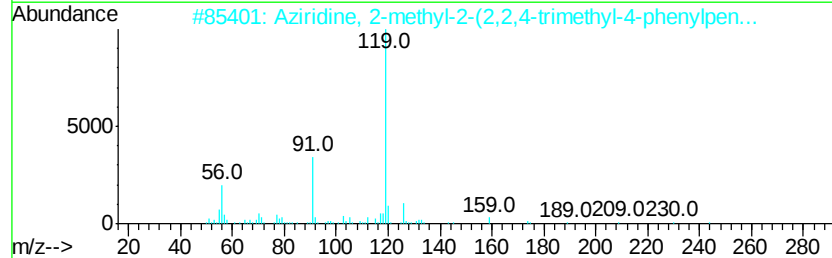
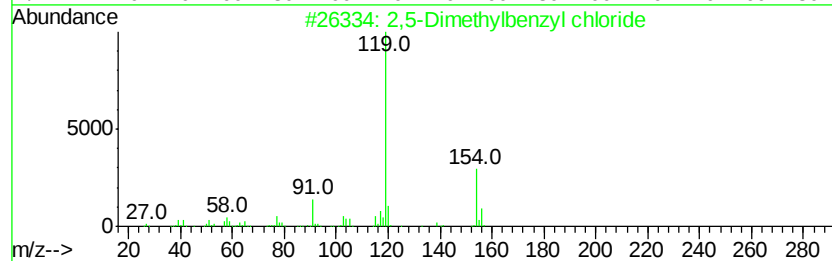
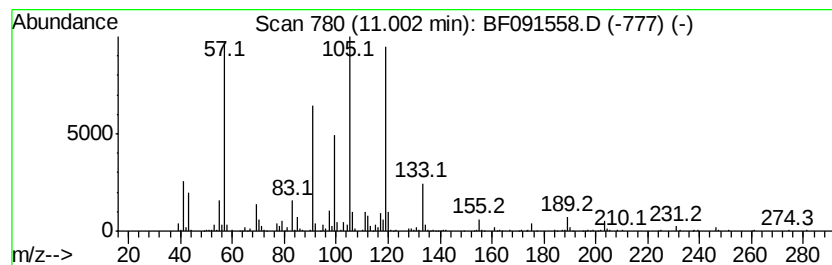
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 unknown11.00 Concentration Rank 16

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.00 | 13.08 ng | 1525140 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 2,5-Dimethylbenzyl chloride         | 154 | C9H11Cl  | 000824-45-3  | 35   |
| 2    |    | Aziridine, 2-methyl-2-(2,2,4-tri... | 245 | C17H27N  | 1000293-28-5 | 35   |
| 3    |    | 2-Isothiocyanto-2,4-dimethyl-4-...  | 233 | C14H19NS | 1000293-27-6 | 35   |
| 4    |    | .alpha.-Bromomesitylene             | 198 | C9H11Br  | 027129-86-8  | 30   |
| 5    |    | Benzene, (1,1-dimethyldecyl)-       | 246 | C18H30   | 027854-40-6  | 30   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

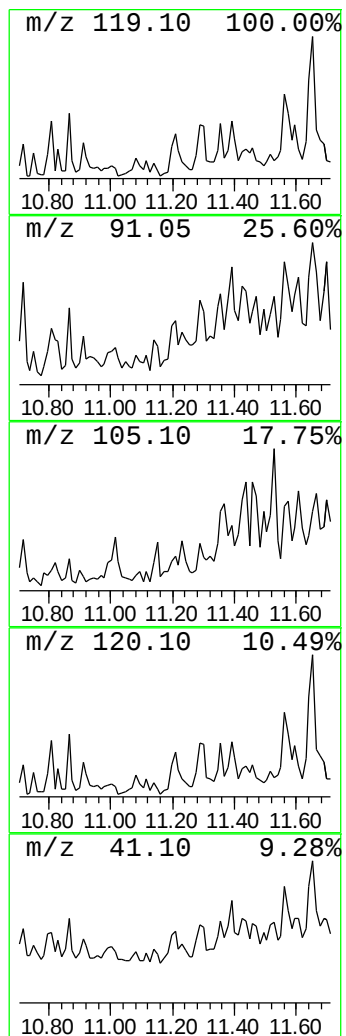
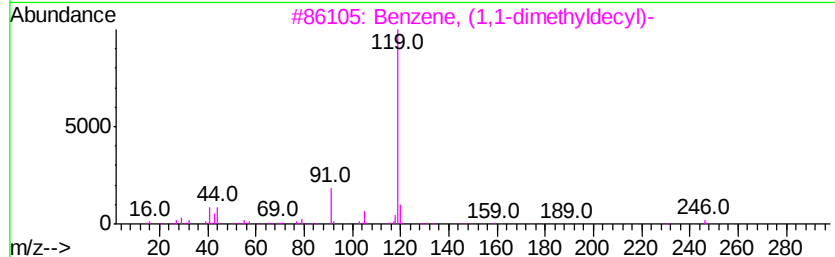
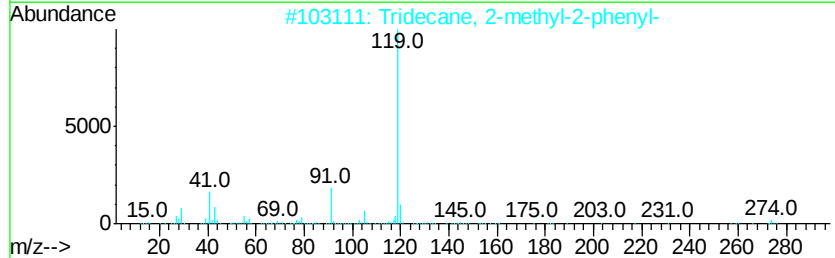
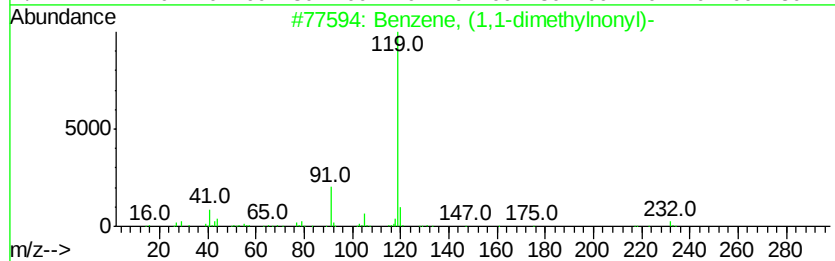
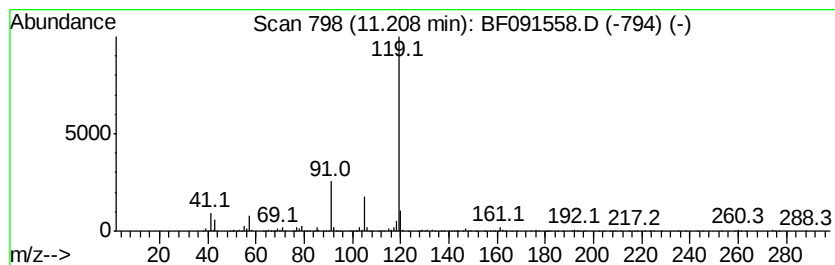
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Benzene, (1,1-dimethylnonyl)- Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.21 | 20.47 ng | 2387640 | Phenanthrene-d10 | 11.33 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|---------|---|-------------------------------------|-----|-------------|--------------|------|
| 1       |   | Benzene, (1,1-dimethylnonyl)-       | 232 | C17H28      | 055191-25-8  | 64   |
| 2       |   | Tridecane, 2-methyl-2-phenyl-       | 274 | C20H34      | 027854-41-7  | 64   |
| 3       |   | Benzene, (1,1-dimethyldecyl)-       | 246 | C18H30      | 027854-40-6  | 64   |
| 4       |   | Benzene, (1,1,4,6,6-pentamethylh... | 246 | C18H30      | 055134-07-1  | 64   |
| 5       |   | Benzamide, N-[2-acetyl-1,1-di(tr... | 383 | C16H15F6N03 | 1000276-52-7 | 59   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

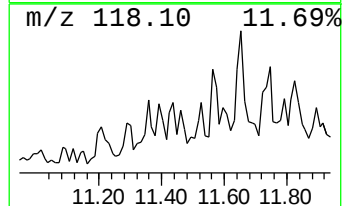
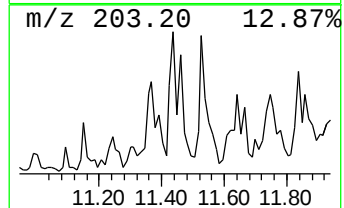
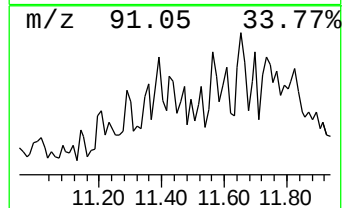
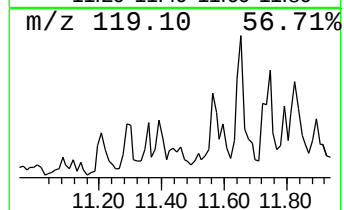
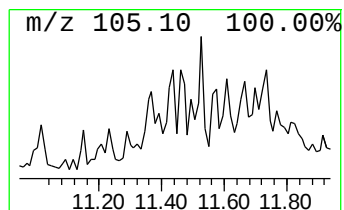
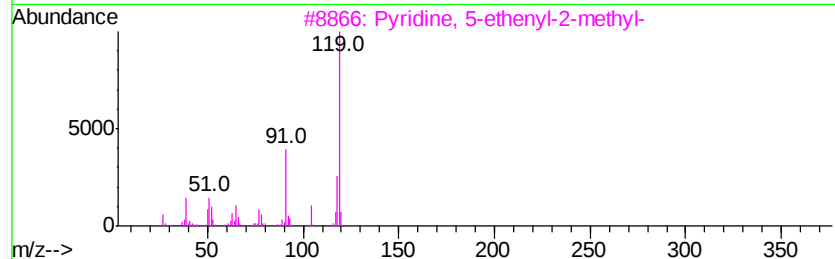
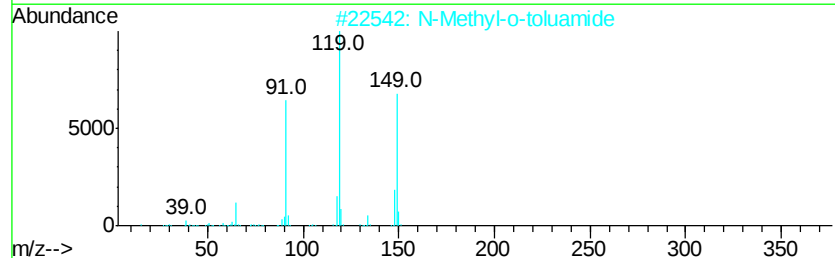
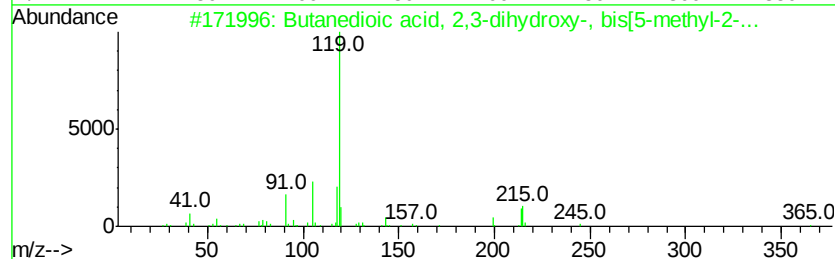
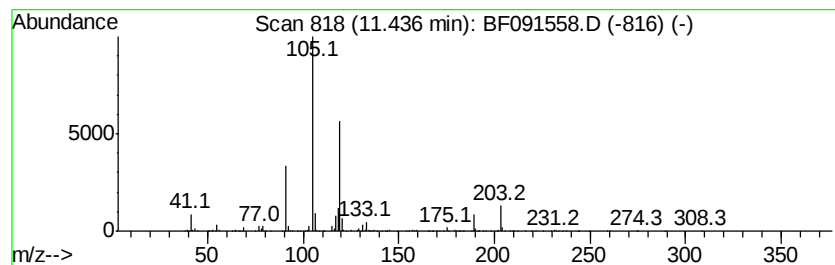
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 unknown11.44 Concentration Rank 10

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.44 | 17.16 ng | 2001240 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Butanedioic acid, 2,3-dihydroxyv-... | 578 | C36H50O6 | 125276-96-2 | 42   |
| 2    |    | N-Methyl-o-toluamide                 | 149 | C9H11NO  | 002170-09-4 | 38   |
| 3    |    | Pyridine, 5-ethenyl-2-methyl-        | 119 | C8H9N    | 000140-76-1 | 32   |
| 4    |    | 1-Chloro-2-methyl-2-phenylpropane    | 168 | C10H13Cl | 000515-40-2 | 27   |
| 5    |    | Cubanol                              | 222 | C15H26O  | 021284-22-0 | 27   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

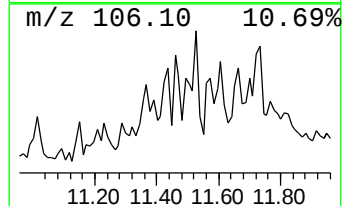
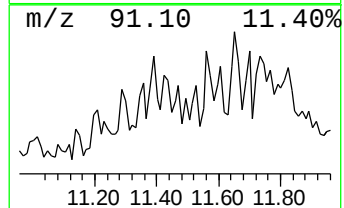
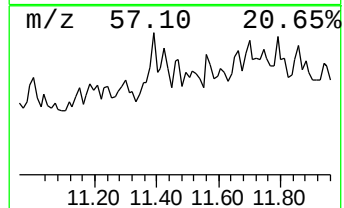
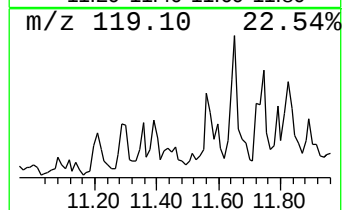
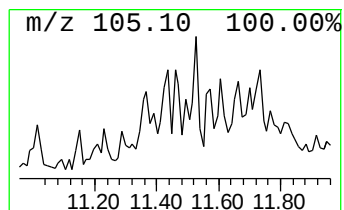
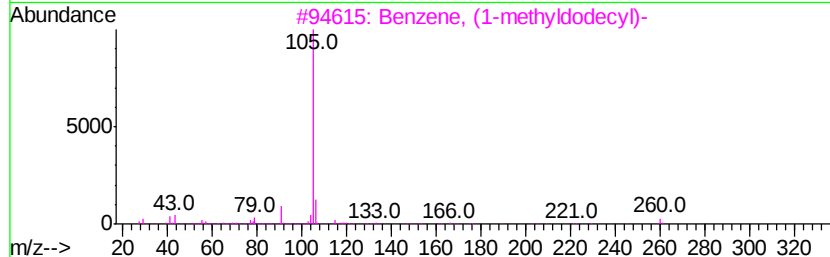
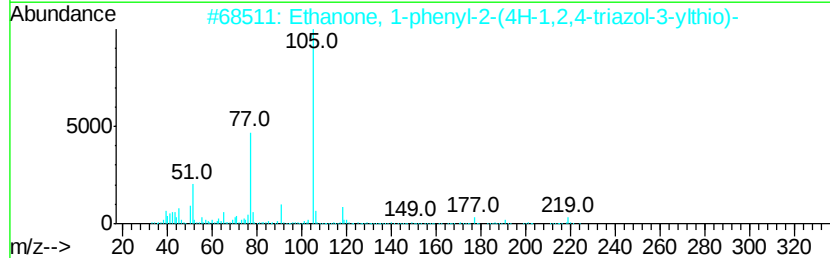
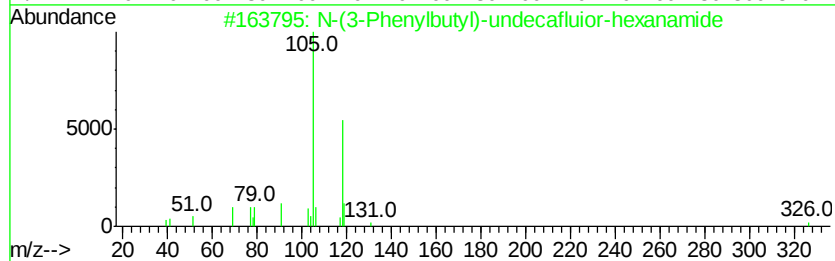
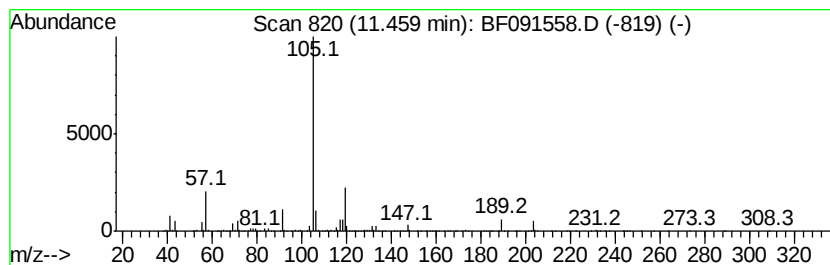
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 unknown11.46 Concentration Rank 15

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.46 | 14.55 ng | 1697410 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|-------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | N-(3-Phenylbutyl)-undecafluor-h...  | 445 | C16H14F11NO | 077970-70-8 | 25   |
| 2    |    | Ethanone, 1-phenyl-2-(4H-1,2,4-t... | 219 | C10H9N3OS   | 032189-00-7 | 23   |
| 3    |    | Benzene, (1-methyldodecyl)-         | 260 | C19H32      | 004534-53-6 | 17   |
| 4    |    | Benzene, (1-methylundecyl)-         | 246 | C18H30      | 002719-61-1 | 17   |
| 5    |    | Benzene, (1-methylnonadecyl)-       | 358 | C26H46      | 002398-66-5 | 17   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

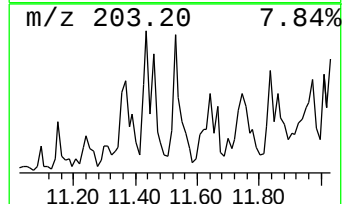
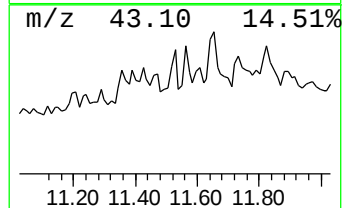
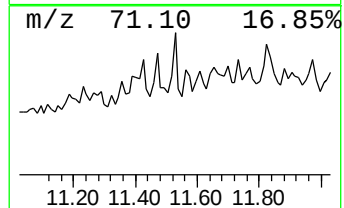
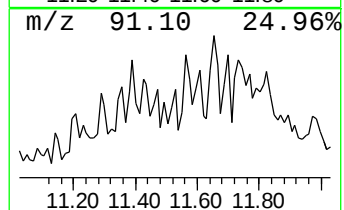
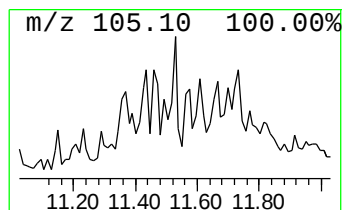
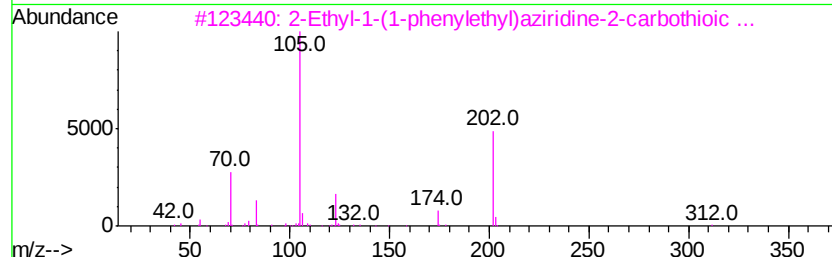
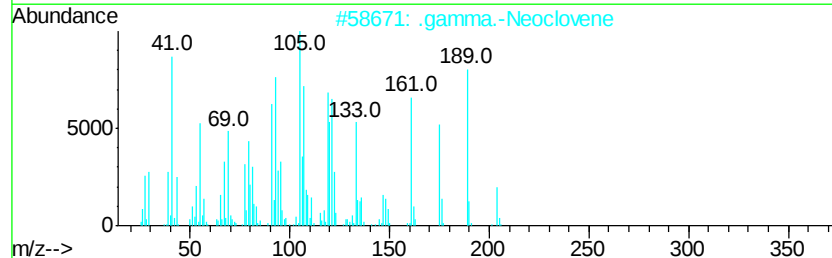
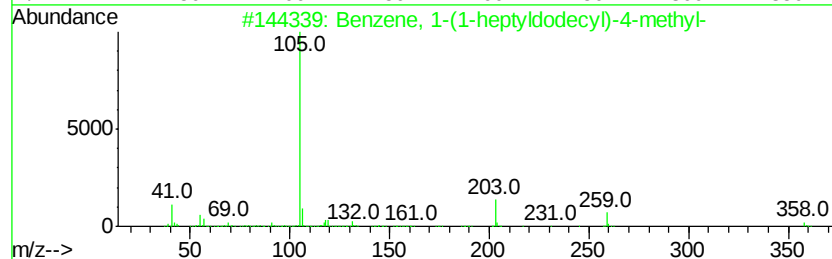
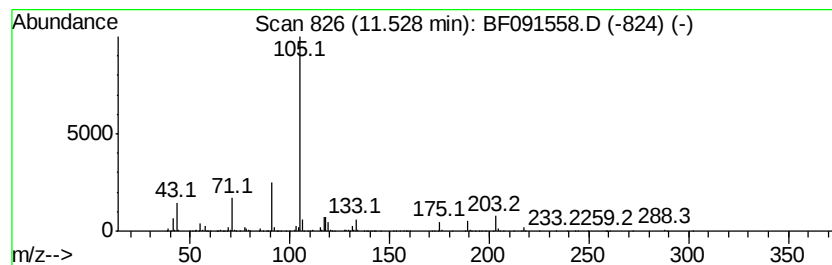
Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 unknown11.53 Concentration Rank 11

| R.T.    | EstConc  | Area                                 | Relative to ISTD | R.T.            |
|---------|----------|--------------------------------------|------------------|-----------------|
| 11.53   | 16.00 ng | 1865810                              | Phenanthrene-d10 | 11.33           |
| Hit# of | 5        | Tentative ID                         | MW MolForm       | CAS# Qual       |
| 1       |          | Benzene, 1-(1-heptylidodecyl)-4-m... | 358 C26H46       | 055191-36-1 27  |
| 2       |          | .gamma.-Neoclovene                   | 204 C15H24       | 1000156-11-7 16 |
| 3       |          | 2-Ethyl-1-(1-phenylethyl)aziridin... | 311 C19H21NOS    | 114950-91-3 12  |
| 4       |          | Benzene, (1-methylundecyl)-          | 246 C18H30       | 002719-61-1 12  |
| 5       |          | 1,1-Cyclopropanedimethanol, 2-me...  | 192 C12H16O2     | 108546-96-9 10  |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

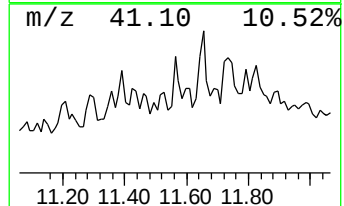
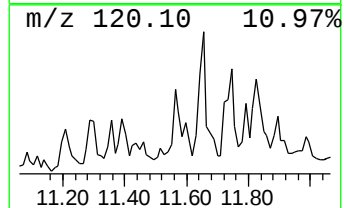
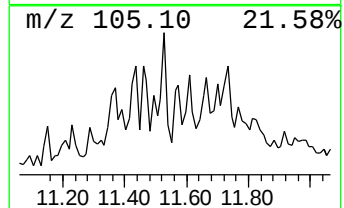
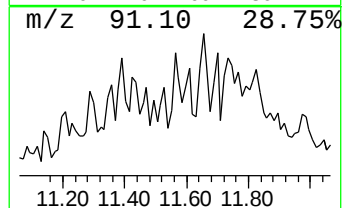
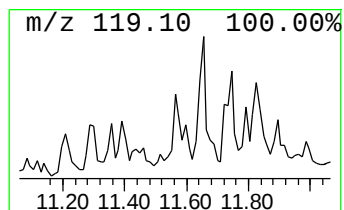
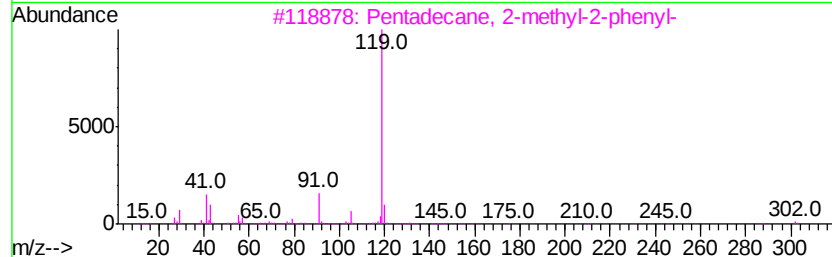
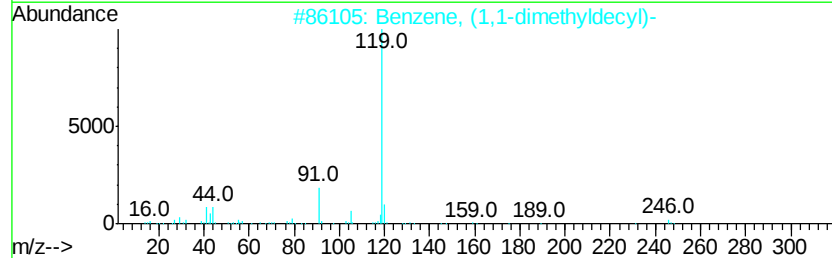
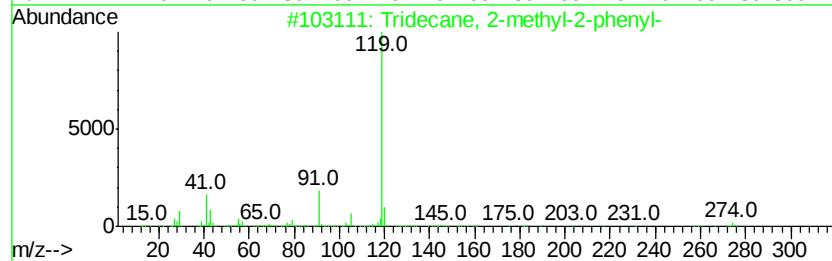
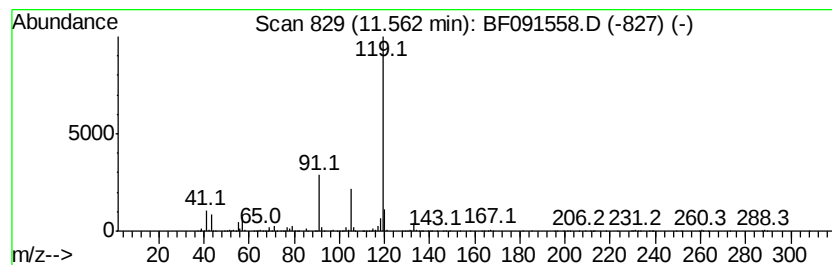
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Tridecane, 2-methyl-2-phenyl- Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.56 | 29.85 ng | 3481430 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane, 2-methyl-2-phenyl-       | 274 | C20H34  | 027854-41-7 | 64   |
| 2    |    | Benzene, (1,1-dimethyldecyl)-       | 246 | C18H30  | 027854-40-6 | 64   |
| 3    |    | Pentadecane, 2-methyl-2-phenyl-     | 302 | C22H38  | 029138-94-1 | 64   |
| 4    |    | Benzene, (1,1-dimethylnonyl)-       | 232 | C17H28  | 055191-25-8 | 64   |
| 5    |    | Benzene, 2-(1-decylundecyl)-1,4-... | 400 | C29H52  | 055373-91-6 | 59   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

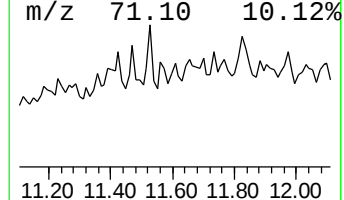
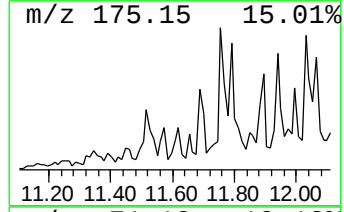
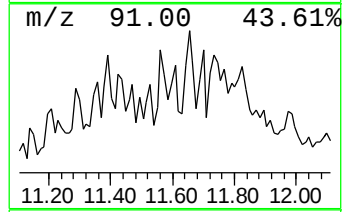
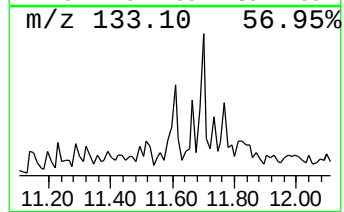
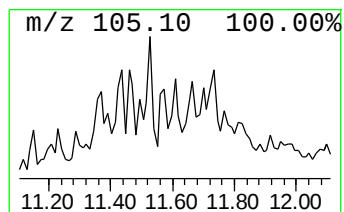
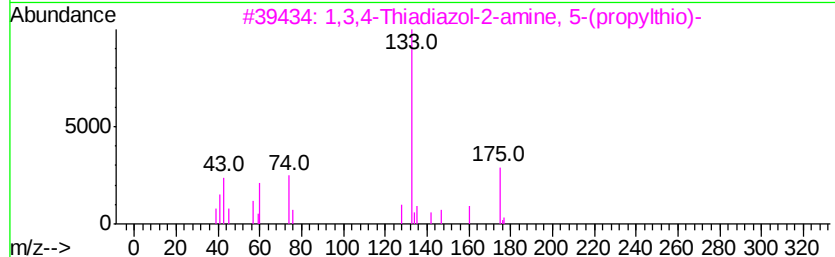
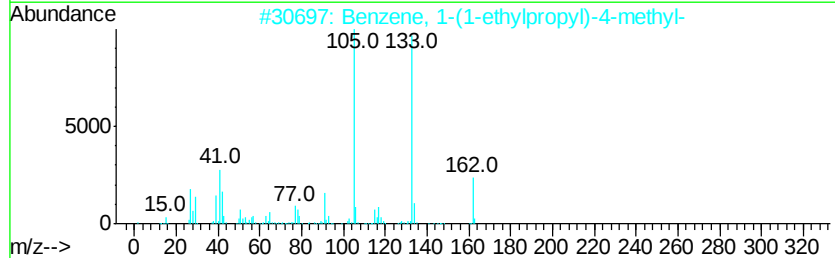
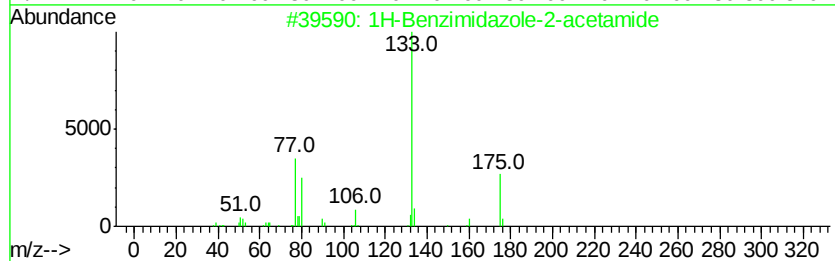
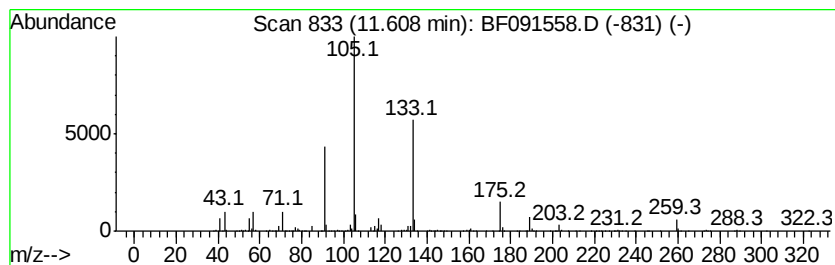
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 unknown11.61 Concentration Rank 9

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.61 | 18.34 ng | 2139630 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 1H-Benzimidazole-2-acetamide        | 175 | C9H9N3O  | 060792-56-5  | 43   |
| 2    |    | Benzene, 1-(1-ethylpropyl)-4-met... | 162 | C12H18   | 022975-58-2  | 42   |
| 3    |    | 1,3,4-Thiadiazol-2-amine, 5-(pro... | 175 | C5H9N3S2 | 030062-49-8  | 38   |
| 4    |    | 4-Indolyl acetate                   | 175 | C10H9NO2 | 005585-96-6  | 37   |
| 5    |    | 3-Phenylpropanoic acid, cyclobut... | 204 | C13H16O2 | 1000282-93-3 | 33   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

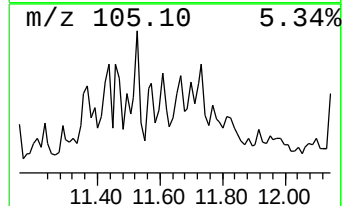
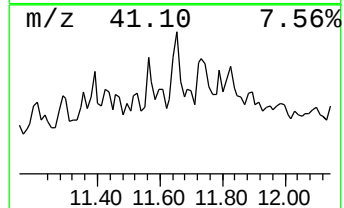
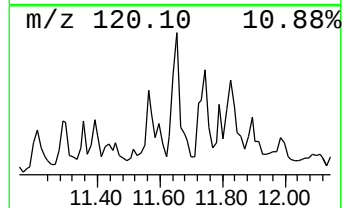
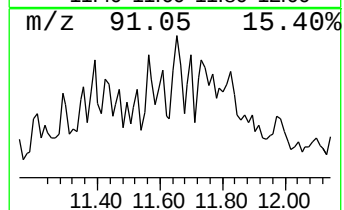
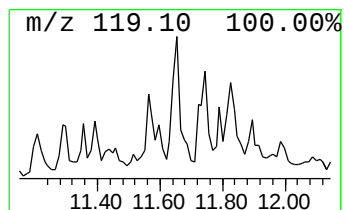
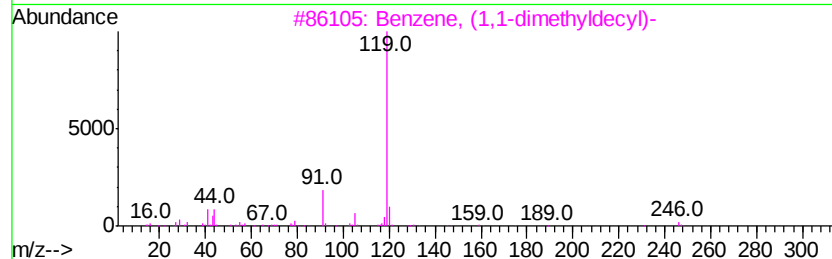
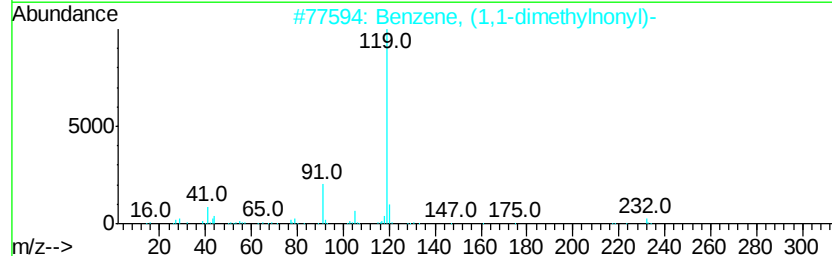
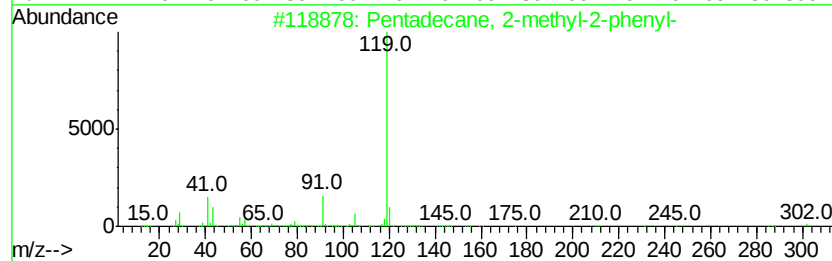
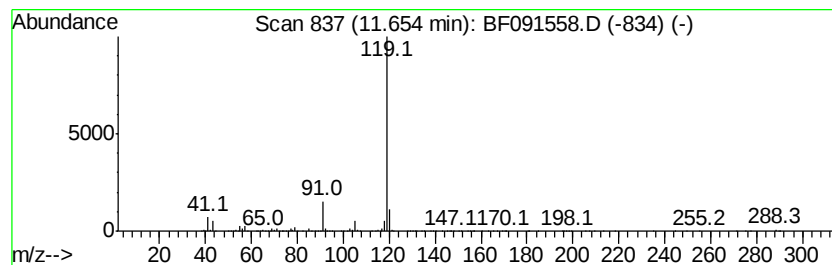
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Pentadecane, 2-methyl-2-phe... Concentration Rank 2

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.65 | 50.47 ng | 5886630 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Pentadecane, 2-methyl-2-phenyl-     | 302 | C22H38   | 029138-94-1 | 83   |
| 2    |    | Benzene, (1,1-dimethylnonyl)-       | 232 | C17H28   | 055191-25-8 | 83   |
| 3    |    | Benzene, (1,1-dimethyldecyl)-       | 246 | C18H30   | 027854-40-6 | 83   |
| 4    |    | Benzene, (1,1,4,6,6-pentamethylh... | 246 | C18H30   | 055134-07-1 | 78   |
| 5    |    | Dicumyl peroxide                    | 270 | C18H22O2 | 000080-43-3 | 78   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF121316\  
Data File : BF091558.D  
Acq On : 13 Dec 2016 17:13  
Operator : UM/SJ  
Sample : H5922-01 5X  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
WC120716

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF120916.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

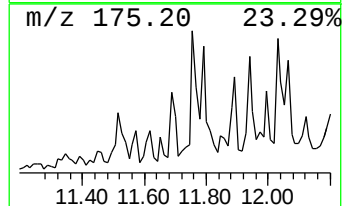
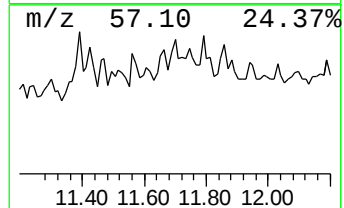
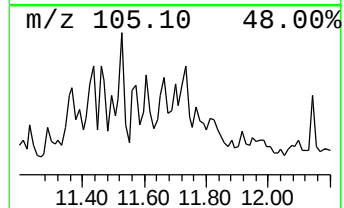
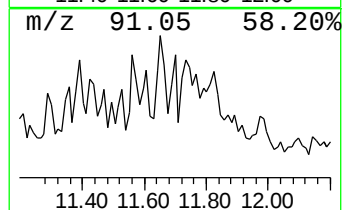
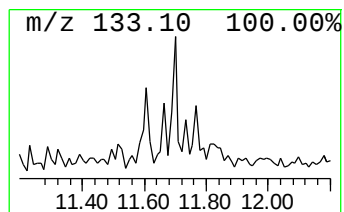
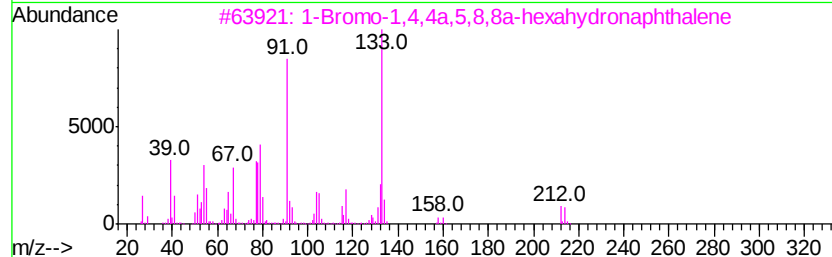
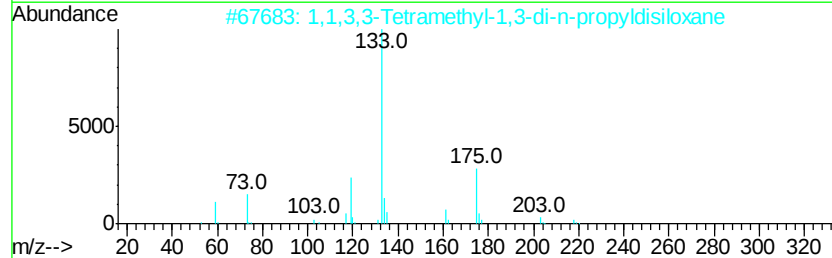
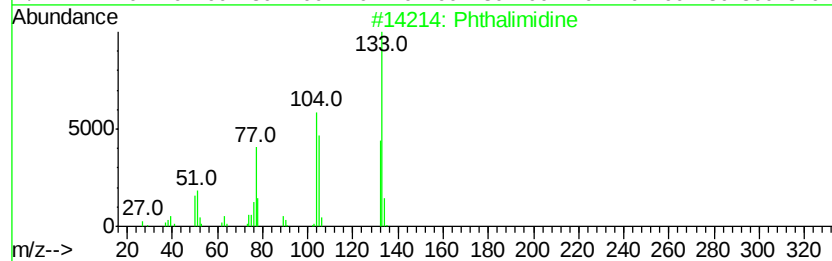
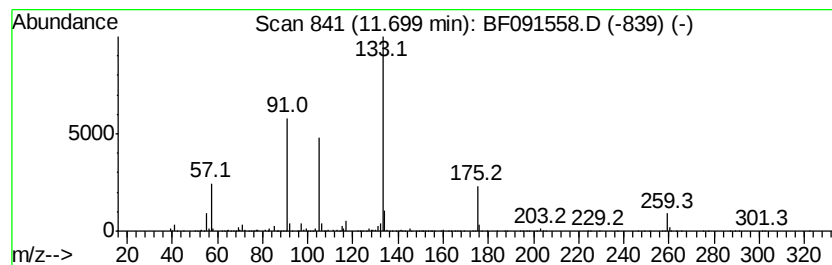
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Phthalimidine Concentration Rank 8

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.70 | 18.43 ng | 2149920 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Phthalimidine                       | 133 | C8H7NO     | 000480-91-1  | 53   |
| 2    |    | 1,1,3,3-Tetramethyl-1,3-di-n-pro... | 218 | C10H26OSi2 | 018001-73-5  | 50   |
| 3    |    | 1-Bromo-1,4,4a,5,8,8a-hexahydron... | 212 | C10H13Br   | 1000192-97-4 | 47   |
| 4    |    | Benzene, 1-(1,1-dimethylethyl)-3... | 148 | C11H16     | 001075-38-3  | 45   |
| 5    |    | 6-Aminoindazole                     | 133 | C7H7N3     | 006967-12-0  | 45   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF121316\  
 Data File : BF091558.D  
 Acq On : 13 Dec 2016 17:13  
 Operator : UM/SJ  
 Sample : H5922-01 5X  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 WC120716

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF120916.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 |
| unknown6.58          | 6.58  | 19.5    | ng    | 1692420  | 1                     | 6.81  | 1734460 | 20.0 |
| 2-Butyn-1-al diet... | 9.08  | 9.9     | ng    | 1175540  | 3                     | 9.85  | 2385030 | 20.0 |
| Benzene, (1,1-dim... | 10.52 | 12.3    | ng    | 1469760  | 3                     | 9.85  | 2385030 | 20.0 |
| Disiloxane, 1,1,3... | 10.72 | 15.0    | ng    | 1752800  | 4                     | 11.33 | 2332700 | 20.0 |
| Benzene, (1,1,4,6... | 10.81 | 21.7    | ng    | 2534370  | 4                     | 11.33 | 2332700 | 20.0 |
| unknown11.00         | 11.00 | 13.1    | ng    | 1525140  | 4                     | 11.33 | 2332700 | 20.0 |
| Benzene, (1,1-dim... | 11.21 | 20.5    | ng    | 2387640  | 4                     | 11.33 | 2332700 | 20.0 |
| unknown11.44         | 11.44 | 17.2    | ng    | 2001240  | 4                     | 11.33 | 2332700 | 20.0 |
| unknown11.46         | 11.46 | 14.6    | ng    | 1697410  | 4                     | 11.33 | 2332700 | 20.0 |
| unknown11.53         | 11.53 | 16.0    | ng    | 1865810  | 4                     | 11.33 | 2332700 | 20.0 |
| Tridecane, 2-meth... | 11.56 | 29.9    | ng    | 3481430  | 4                     | 11.33 | 2332700 | 20.0 |
| unknown11.61         | 11.61 | 18.3    | ng    | 2139630  | 4                     | 11.33 | 2332700 | 20.0 |
| Pentadecane, 2-me... | 11.65 | 50.5    | ng    | 5886630  | 4                     | 11.33 | 2332700 | 20.0 |
| Phthalimidine        | 11.70 | 18.4    | ng    | 2149920  | 4                     | 11.33 | 2332700 | 20.0 |