

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
 Data File : BF085289.D  
 Acq On : 9 Mar 2016 6:58  
 Operator : UM/SJ  
 Sample : H1712-06  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-43-COMP-0.5-2.5

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-BF030316.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         | 5.167       | 250           | 252         | 262          | rVB      | 680528         | 1114277       | 25.42%          | 2.538%        |
| 2         | 5.579       | 284           | 288         | 290          | rBV      | 3502992        | 4383129       | 100.00%         | 9.982%        |
| 3         | 6.584       | 373           | 376         | 379          | rBV      | 2822762        | 3849713       | 87.83%          | 8.767%        |
| 4         | 6.733       | 386           | 389         | 391          | rBV      | 3891864        | 4375318       | 99.82%          | 9.964%        |
| 5         | 6.962       | 407           | 409         | 411          | rBV      | 1211216        | 1020002       | 23.27%          | 2.323%        |
| 6         | 7.122       | 420           | 423         | 425          | rVB      | 3460018        | 3428960       | 78.23%          | 7.809%        |
| 7         | 7.522       | 455           | 458         | 461          | rBV      | 2330605        | 2464501       | 56.23%          | 5.612%        |
| 8         | 7.602       | 463           | 465         | 468          | rVB      | 76892          | 97226         | 2.22%           | 0.221%        |
| 9         | 8.242       | 519           | 521         | 524          | rBV      | 1348963        | 1266798       | 28.90%          | 2.885%        |
| 10        | 9.328       | 613           | 616         | 618          | rBV      | 2859385        | 3891663       | 88.79%          | 8.862%        |
| 11        | 9.659       | 640           | 645         | 648          | rBV3     | 24309          | 49648         | 1.13%           | 0.113%        |
| 12        | 9.716       | 648           | 650         | 651          | rBV      | 739649         | 731753        | 16.69%          | 1.666%        |
| 13        | 9.865       | 661           | 663         | 665          | rBV      | 36886          | 47350         | 1.08%           | 0.108%        |
| 14        | 9.933       | 666           | 669         | 671          | rVB      | 134159         | 143137        | 3.27%           | 0.326%        |
| 15        | 10.002      | 673           | 675         | 677          | rVB      | 1397651        | 1201960       | 27.42%          | 2.737%        |
| 16        | 10.162      | 686           | 689         | 690          | rBV      | 38071          | 53237         | 1.21%           | 0.121%        |
| 17        | 10.253      | 695           | 697         | 699          | rVB3     | 39047          | 48486         | 1.11%           | 0.110%        |
| 18        | 10.425      | 711           | 712         | 714          | rVB      | 217820         | 198996        | 4.54%           | 0.453%        |
| 19        | 10.653      | 729           | 732         | 735          | rBV      | 71239          | 138398        | 3.16%           | 0.315%        |
| 20        | 10.791      | 741           | 744         | 746          | rBV      | 2129119        | 2193264       | 50.04%          | 4.995%        |
| 21        | 10.905      | 752           | 754         | 760          | rBV      | 190933         | 290379        | 6.62%           | 0.661%        |
| 22        | 11.099      | 768           | 771         | 772          | rBV2     | 33919          | 59890         | 1.37%           | 0.136%        |
| 23        | 11.351      | 789           | 793         | 794          | rBV      | 131024         | 131772        | 3.01%           | 0.300%        |
| 24        | 11.374      | 794           | 795         | 798          | rVV2     | 36306          | 50917         | 1.16%           | 0.116%        |
| 25        | 11.488      | 802           | 805         | 806          | rVV      | 1030398        | 1055666       | 24.08%          | 2.404%        |
| 26        | 11.556      | 808           | 811         | 813          | rVB      | 100295         | 118586        | 2.71%           | 0.270%        |
| 27        | 11.716      | 822           | 825         | 829          | rBV4     | 25912          | 62938         | 1.44%           | 0.143%        |
| 28        | 11.774      | 829           | 830         | 832          | rVB      | 55246          | 47318         | 1.08%           | 0.108%        |
| 29        | 12.002      | 846           | 850         | 854          | rBV2     | 240271         | 417828        | 9.53%           | 0.952%        |
| 30        | 12.094      | 857           | 858         | 862          | rVV      | 95689          | 159632        | 3.64%           | 0.364%        |
| 31        | 12.185      | 862           | 866         | 868          | rVV      | 36655          | 58698         | 1.34%           | 0.134%        |
| 32        | 12.276      | 870           | 874         | 878          | rBV5     | 51781          | 122455        | 2.79%           | 0.279%        |
| 33        | 12.528      | 894           | 896         | 898          | rBV      | 49509          | 72688         | 1.66%           | 0.166%        |
| 34        | 12.619      | 902           | 904         | 907          | rVB      | 45091          | 49328         | 1.13%           | 0.112%        |

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 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-BF030316.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 12.699 | 908  | 911  | 913  | rBV  | 465235  | 541446  | 12.35% | 1.233% |
| 36 | 12.791 | 917  | 919  | 920  | rBV2 | 196745  | 204432  | 4.66%  | 0.466% |
| 37 | 12.825 | 920  | 922  | 923  | rVB  | 63173   | 68244   | 1.56%  | 0.155% |
| 38 | 12.928 | 928  | 931  | 933  | rBV  | 431862  | 544341  | 12.42% | 1.240% |
| 39 | 12.974 | 933  | 935  | 937  | rVB3 | 28538   | 46092   | 1.05%  | 0.105% |
| 40 | 13.065 | 941  | 943  | 946  | rVB  | 2240407 | 2532445 | 57.78% | 5.767% |
| 41 | 13.145 | 946  | 950  | 953  | rBV3 | 52398   | 82726   | 1.89%  | 0.188% |
| 42 | 13.248 | 953  | 959  | 961  | rBV2 | 82628   | 150694  | 3.44%  | 0.343% |
| 43 | 13.351 | 966  | 968  | 972  | rVB2 | 64834   | 100758  | 2.30%  | 0.229% |
| 44 | 13.477 | 978  | 979  | 981  | rBV  | 52250   | 53836   | 1.23%  | 0.123% |
| 45 | 13.545 | 983  | 985  | 988  | rVB3 | 29744   | 58168   | 1.33%  | 0.132% |
| 46 | 13.637 | 990  | 993  | 1000 | rBV5 | 23766   | 99051   | 2.26%  | 0.226% |
| 47 | 13.751 | 1002 | 1003 | 1006 | rVB  | 61560   | 77165   | 1.76%  | 0.176% |
| 48 | 13.865 | 1010 | 1013 | 1015 | rBV2 | 48983   | 93245   | 2.13%  | 0.212% |
| 49 | 13.922 | 1017 | 1018 | 1020 | rVV2 | 40768   | 57354   | 1.31%  | 0.131% |
| 50 | 13.957 | 1020 | 1021 | 1026 | rVB  | 215595  | 314783  | 7.18%  | 0.717% |
| 51 | 14.117 | 1031 | 1035 | 1041 | rBV2 | 980109  | 1714974 | 39.13% | 3.905% |
| 52 | 14.220 | 1042 | 1044 | 1046 | rBV2 | 40676   | 69182   | 1.58%  | 0.158% |
| 53 | 14.505 | 1067 | 1069 | 1071 | rBV  | 73881   | 85373   | 1.95%  | 0.194% |
| 54 | 14.597 | 1074 | 1077 | 1079 | rBV3 | 86192   | 160613  | 3.66%  | 0.366% |
| 55 | 14.871 | 1099 | 1101 | 1104 | rBV  | 106431  | 132280  | 3.02%  | 0.301% |
| 56 | 14.974 | 1108 | 1110 | 1114 | rBV3 | 53284   | 83310   | 1.90%  | 0.190% |
| 57 | 15.157 | 1123 | 1126 | 1130 | rBV  | 282859  | 556395  | 12.69% | 1.267% |
| 58 | 15.237 | 1130 | 1133 | 1134 | rVV  | 97516   | 152980  | 3.49%  | 0.348% |
| 59 | 15.260 | 1134 | 1135 | 1137 | rVB  | 63526   | 81603   | 1.86%  | 0.186% |
| 60 | 15.465 | 1150 | 1153 | 1155 | rVB  | 136969  | 203247  | 4.64%  | 0.463% |
| 61 | 15.523 | 1155 | 1158 | 1161 | rVV  | 229896  | 278695  | 6.36%  | 0.635% |
| 62 | 15.591 | 1161 | 1164 | 1169 | rVB  | 698238  | 1015437 | 23.17% | 2.312% |
| 63 | 16.048 | 1202 | 1204 | 1207 | rBV  | 59750   | 88522   | 2.02%  | 0.202% |
| 64 | 16.105 | 1207 | 1209 | 1212 | rBV4 | 39354   | 78045   | 1.78%  | 0.178% |
| 65 | 17.043 | 1288 | 1291 | 1297 | rVB2 | 91810   | 228998  | 5.22%  | 0.521% |
| 66 | 17.226 | 1305 | 1307 | 1309 | rBV  | 29979   | 62429   | 1.42%  | 0.142% |
| 67 | 17.488 | 1327 | 1330 | 1339 | rVB  | 84093   | 200834  | 4.58%  | 0.457% |
| 68 | 17.648 | 1341 | 1344 | 1349 | rVV6 | 52721   | 185820  | 4.24%  | 0.423% |
| 69 | 17.728 | 1349 | 1351 | 1358 | rVB3 | 44178   | 112550  | 2.57%  | 0.256% |

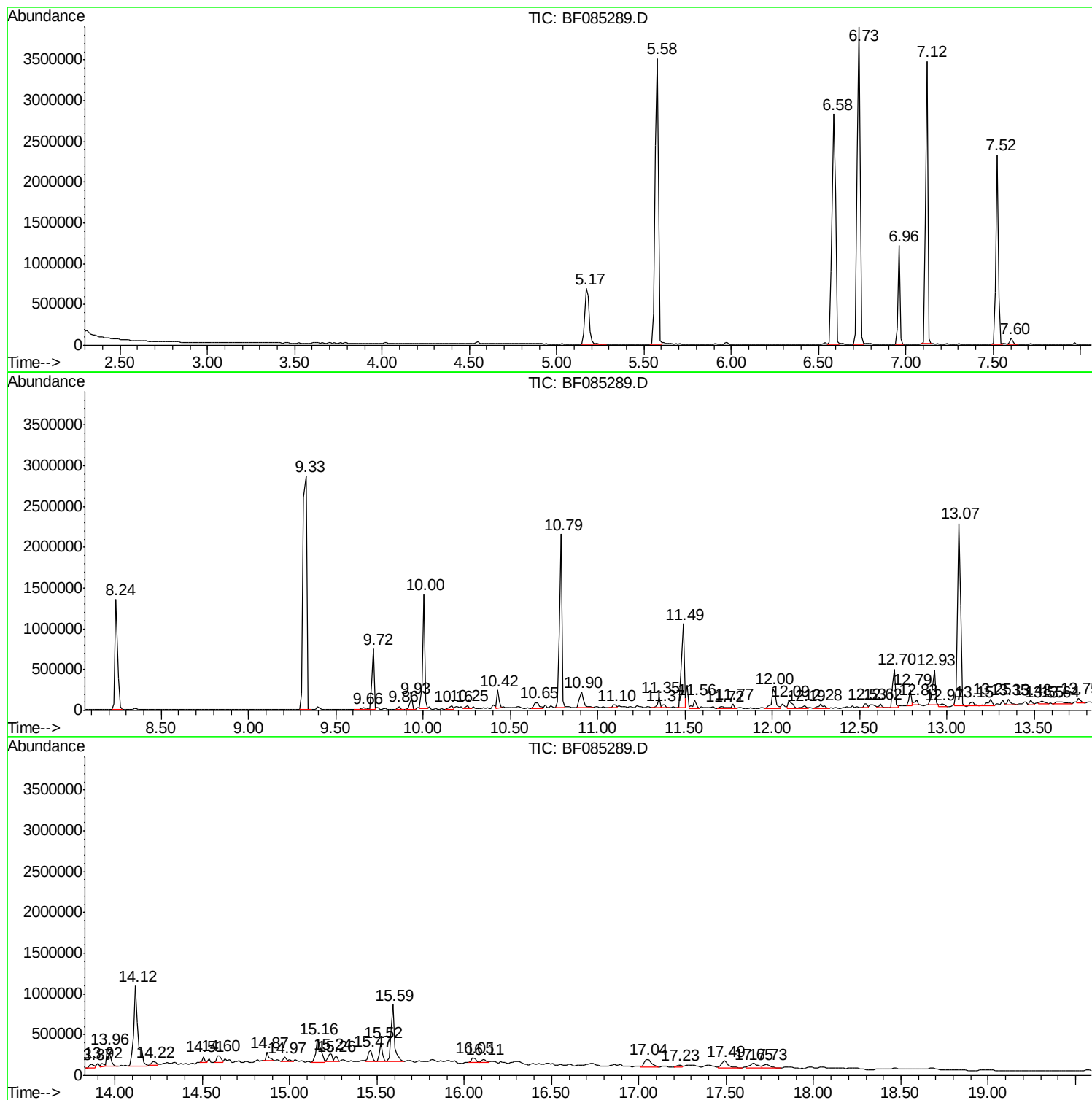
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Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-43-COMP-0.5-2.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

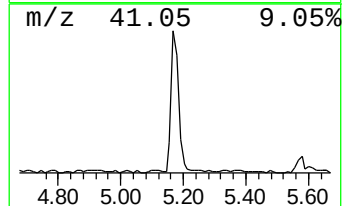
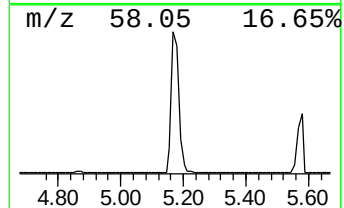
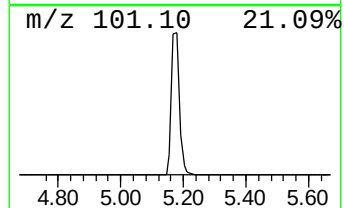
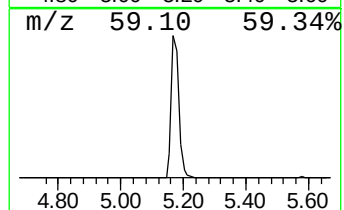
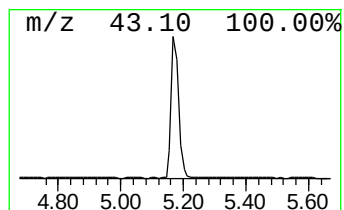
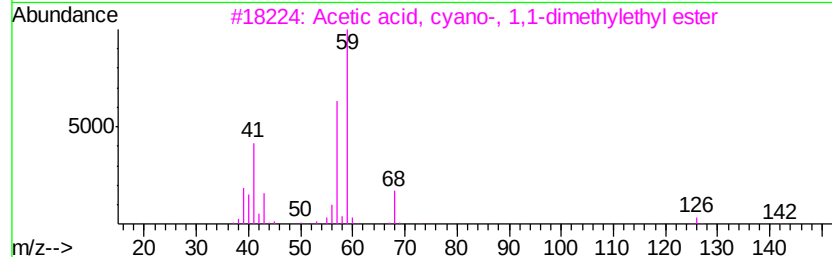
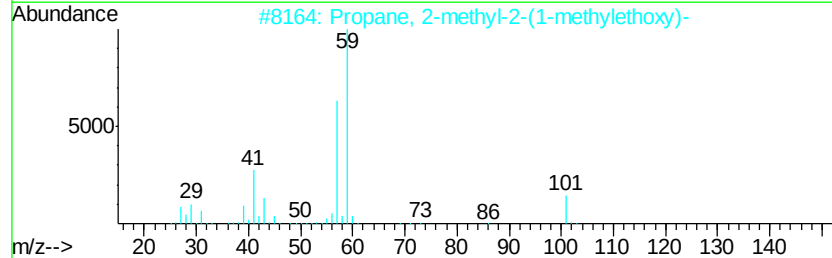
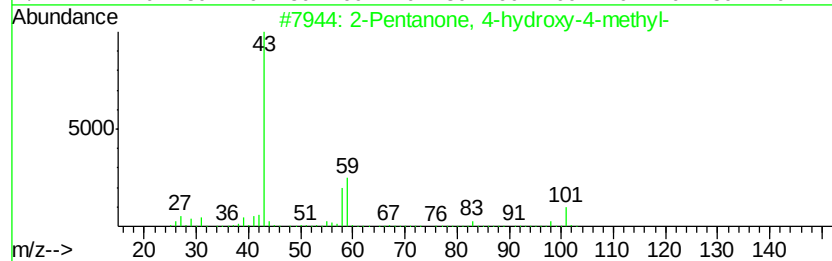
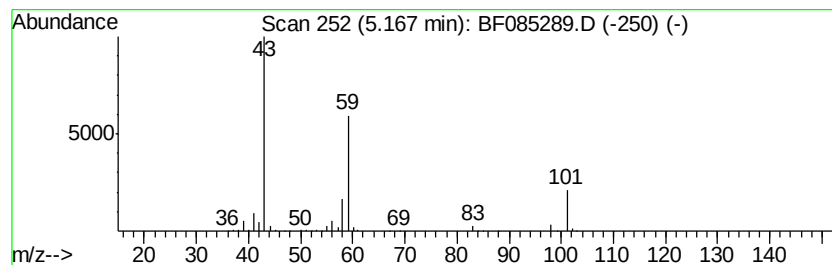
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.17 | 21.85 ng | 1114280 | 1,4-Dichlorobenzene-d4 | 6.96 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 64   |
| 2    |    | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O   | 017348-59-3 | 23   |
| 3    |    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 4    |    | 5-Hexen-2-one                       | 98  | C6H10O   | 000109-49-9 | 9    |
| 5    |    | Butane, 1-ethoxy-                   | 102 | C6H14O   | 000628-81-9 | 9    |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

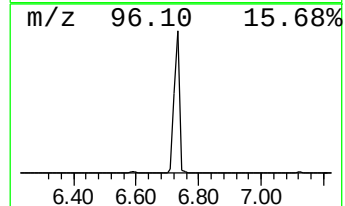
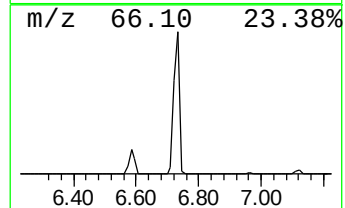
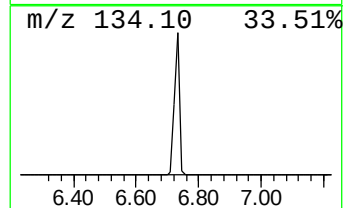
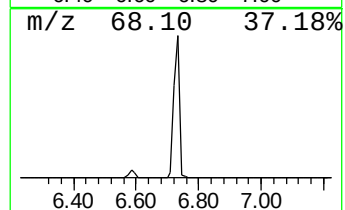
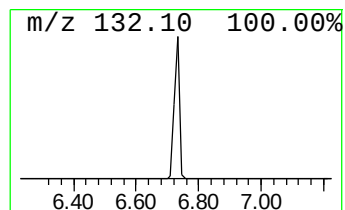
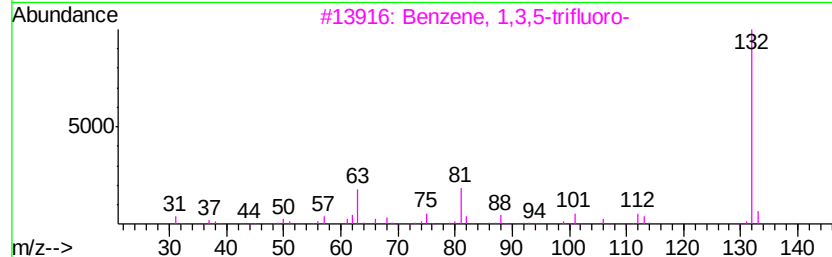
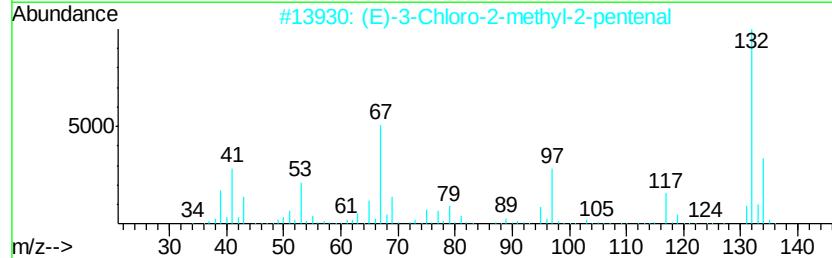
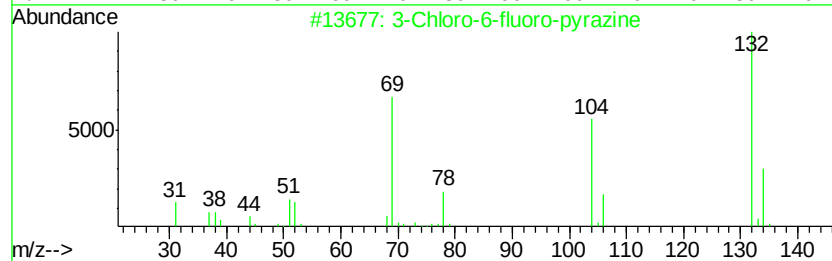
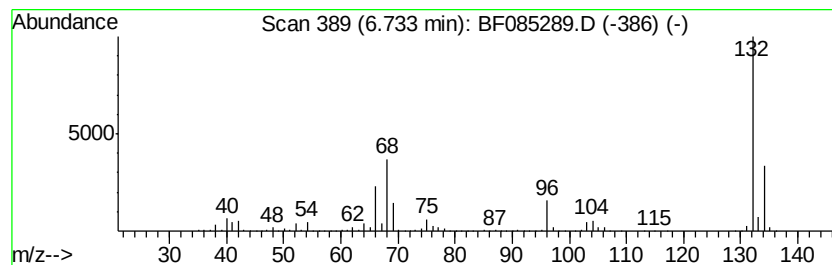
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 unknown6.73 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.73 | 85.79 ng | 4375320 | 1,4-Dichlorobenzene-d4 | 6.96 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 25   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    |    | Benzene, 1,3,5-trifluoro-        | 132 | C6H3F3    | 000372-38-3  | 9    |
| 4    |    | 4-Chloro-2-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-00-5  | 9    |
| 5    |    | 4-Chloro-6-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-01-6  | 9    |



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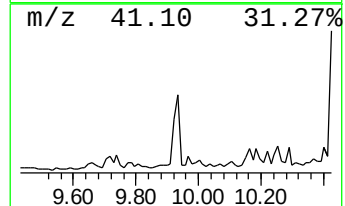
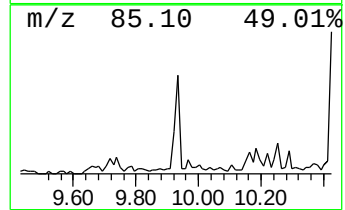
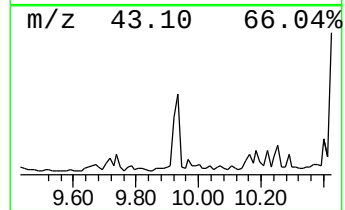
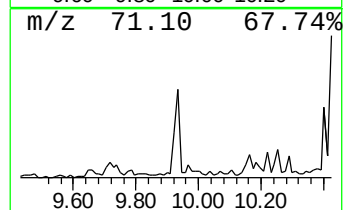
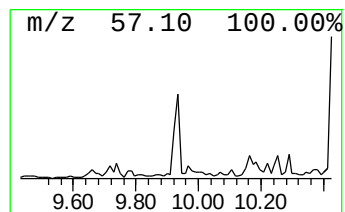
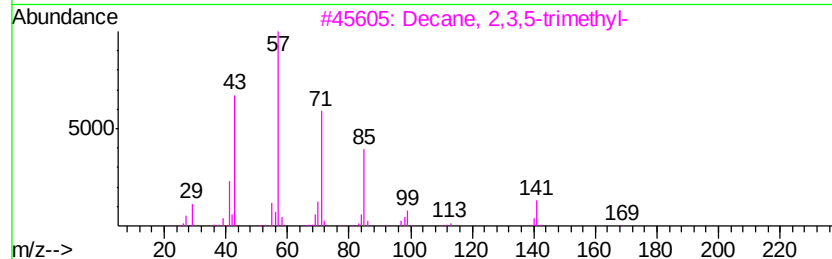
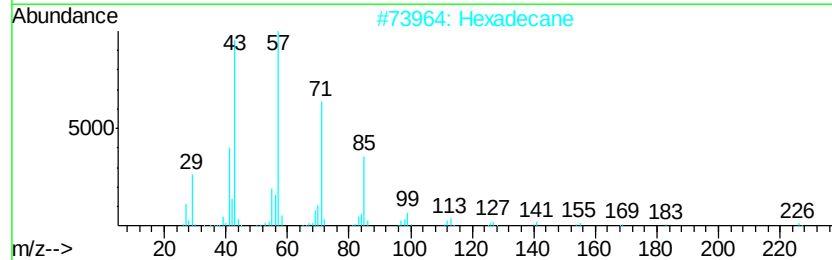
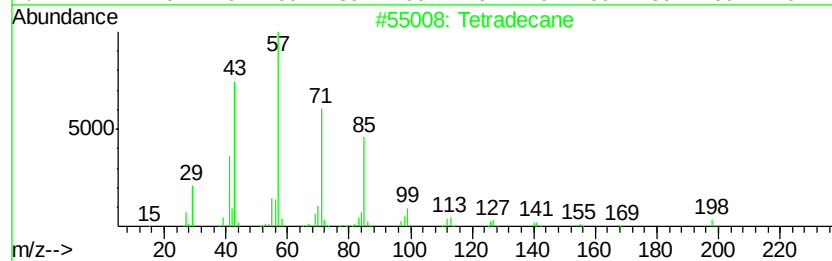
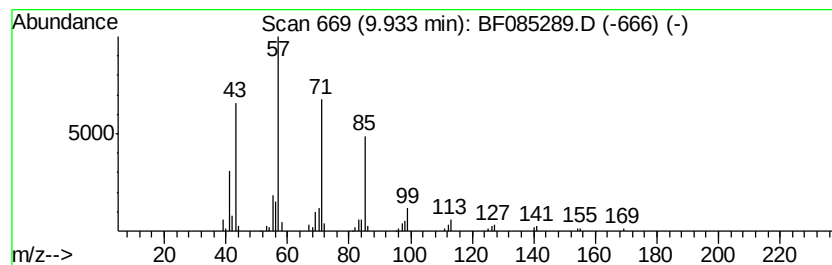
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Tetradecane Concentration Rank 15

| R.T. | EstConc | Area   | Relative to ISTD | R.T.  |
|------|---------|--------|------------------|-------|
| 9.93 | 2.38 ng | 143137 | Acenaphthene-d10 | 10.00 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Tetradecane              | 198 | C14H30  | 000629-59-4 | 87   |
| 2    |    | Hexadecane               | 226 | C16H34  | 000544-76-3 | 86   |
| 3    |    | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 83   |
| 4    |    | Eicosane                 | 282 | C20H42  | 000112-95-8 | 72   |
| 5    |    | Pentadecane              | 212 | C15H32  | 000629-62-9 | 64   |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

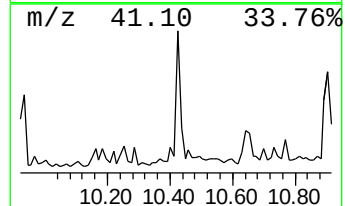
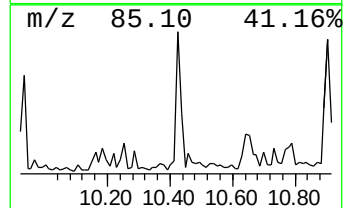
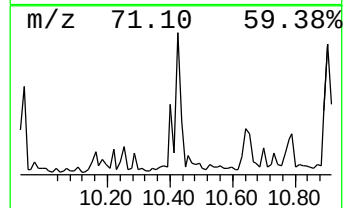
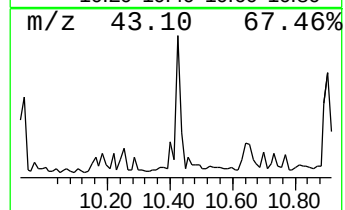
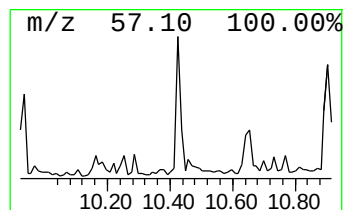
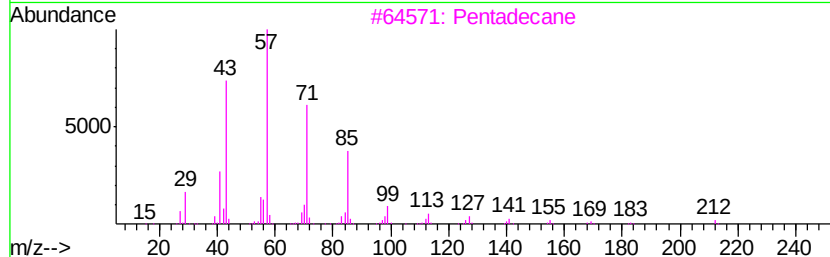
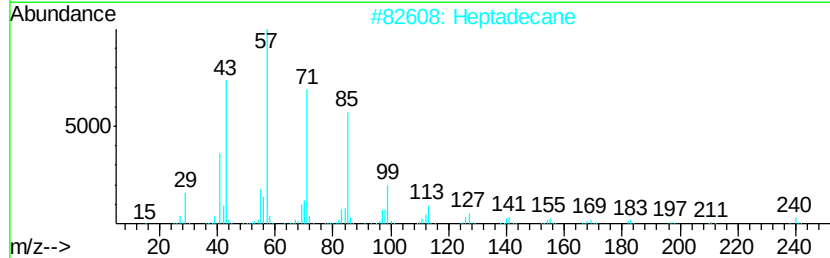
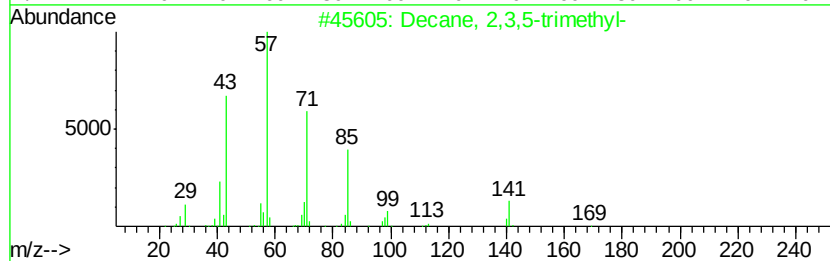
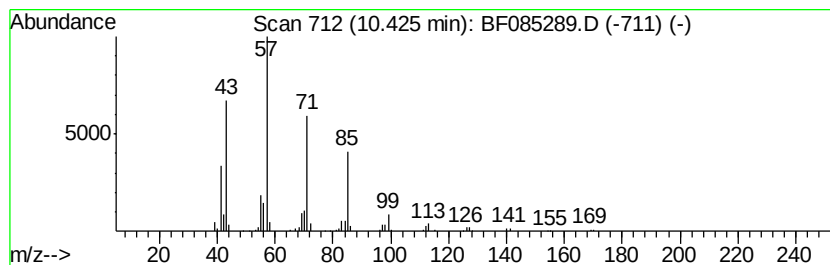
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Decane, 2,3,5-trimethyl- Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.42 | 3.31 ng | 198996 | Acenaphthene-d10 | 10.00 |

| Hit# of | 5                        | Tentative ID | MW     | MolForm     | CAS# | Qual |
|---------|--------------------------|--------------|--------|-------------|------|------|
| 1       | Decane, 2,3,5-trimethyl- | 184          | C13H28 | 062238-11-3 | 83   |      |
| 2       | Heptadecane              | 240          | C17H36 | 000629-78-7 | 83   |      |
| 3       | Pentadecane              | 212          | C15H32 | 000629-62-9 | 83   |      |
| 4       | Nonane, 5-butyl-         | 184          | C13H28 | 017312-63-9 | 80   |      |
| 5       | Hexadecane               | 226          | C16H34 | 000544-76-3 | 72   |      |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

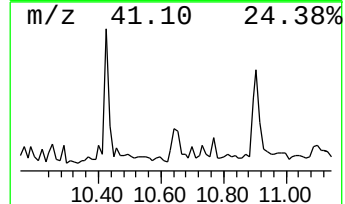
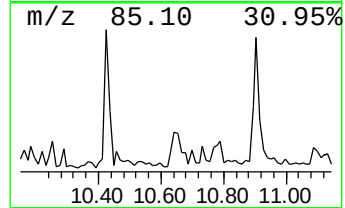
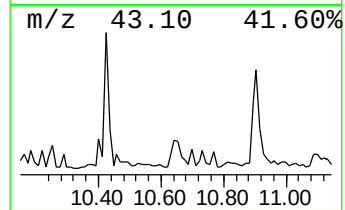
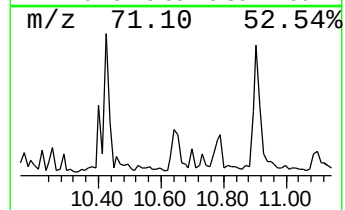
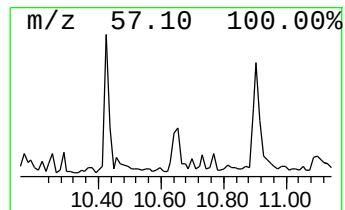
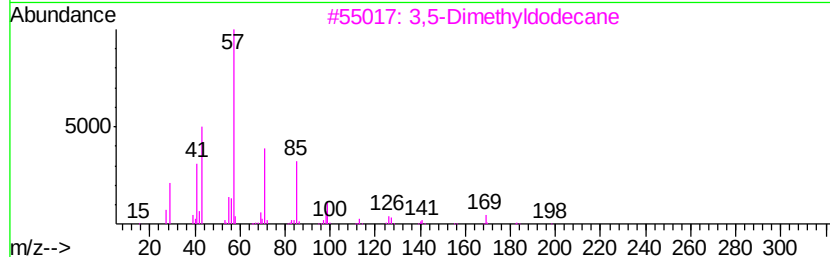
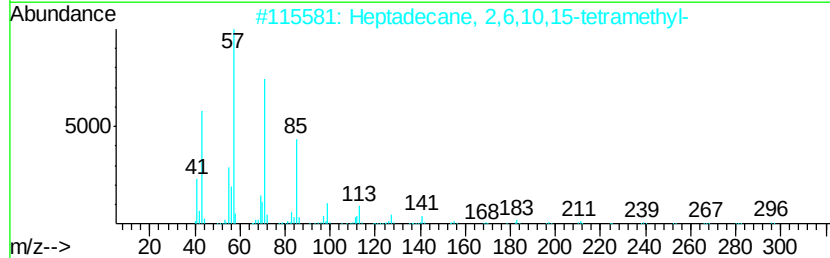
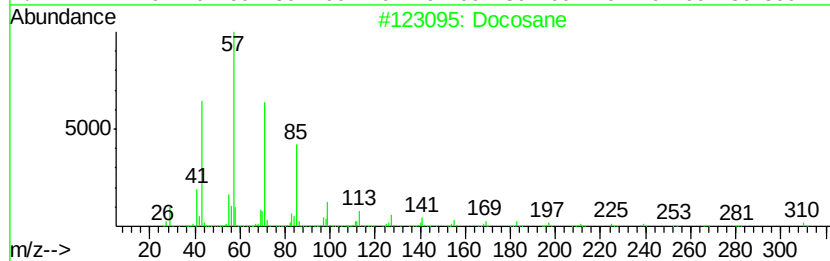
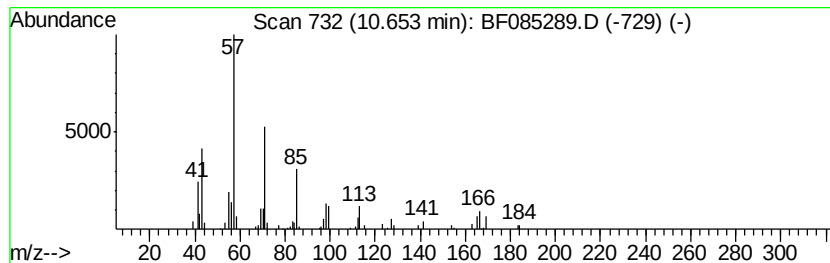
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ClientSampleId :  
SB-43-COMP-0.5-2.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Docosane Concentration Rank 17

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.           |
|---------|-------------------------------------|--------------|------------------|----------------|
| 10.65   | 2.30 ng                             | 138398       | Acenaphthene-d10 | 10.00          |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | Docosane                            |              | 310 C22H46       | 000629-97-0 72 |
| 2       | Heptadecane, 2,6,10,15-tetramethyl- |              | 296 C21H44       | 054833-48-6 68 |
| 3       | 3,5-Dimethyldodecane                |              | 198 C14H30       | 107770-99-0 64 |
| 4       | 2-Bromo dodecane                    |              | 248 C12H25Br     | 013187-99-0 62 |
| 5       | Decane, 6-ethyl-2-methyl-           |              | 184 C13H28       | 062108-21-8 59 |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
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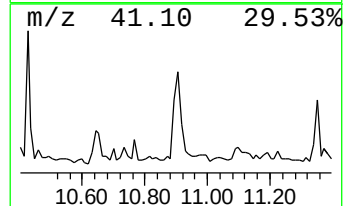
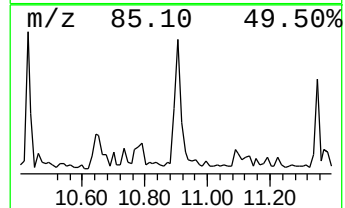
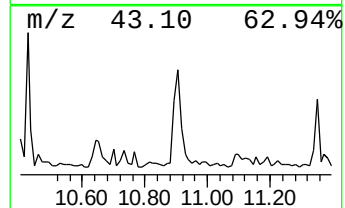
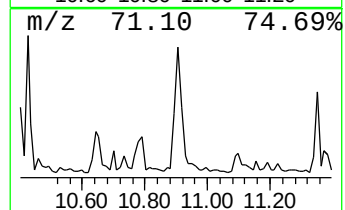
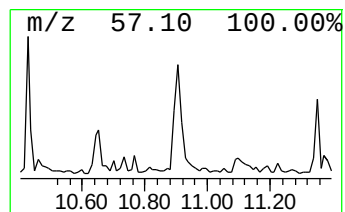
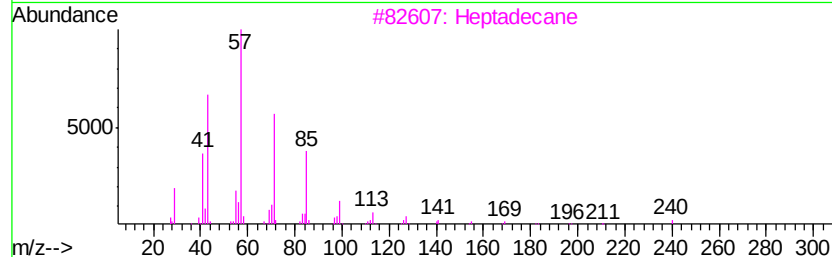
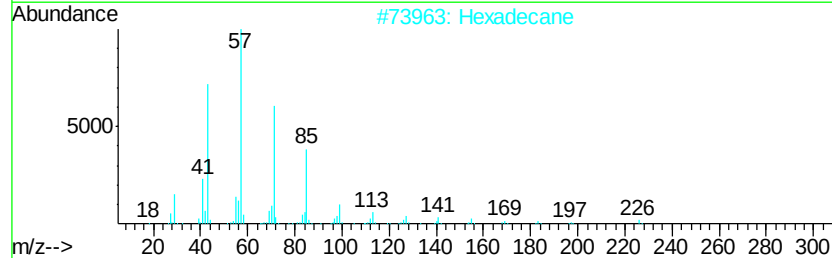
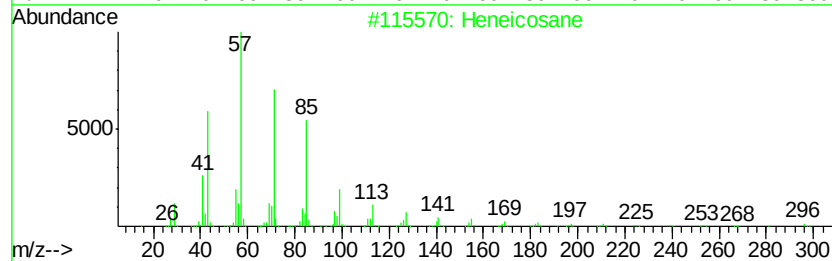
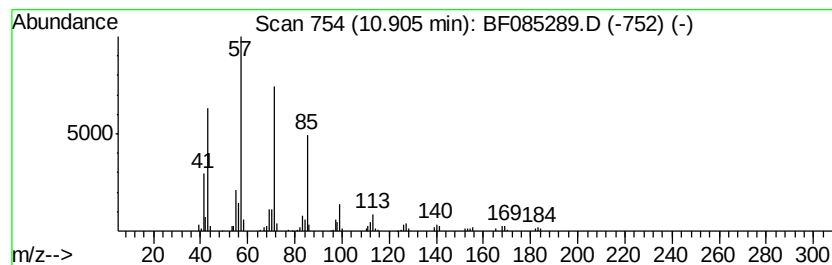
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Heneicosane Concentration Rank 5

| R.T.      | EstConc      | Area   | Relative to ISTD |             | R.T.  |
|-----------|--------------|--------|------------------|-------------|-------|
| 10.90     | 5.50 ng      | 290379 | Phenanthrene-d10 |             | 11.49 |
| Hit# of 5 | Tentative ID | MW     | MolForm          | CAS#        | Qual  |
| 1         | Heneicosane  | 296    | C21H44           | 000629-94-7 | 91    |
| 2         | Hexadecane   | 226    | C16H34           | 000544-76-3 | 90    |
| 3         | Heptadecane  | 240    | C17H36           | 000629-78-7 | 90    |
| 4         | Pentadecane  | 212    | C15H32           | 000629-62-9 | 90    |
| 5         | Dodecane     | 170    | C12H26           | 000112-40-3 | 87    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

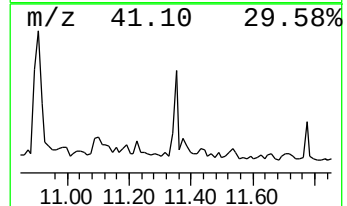
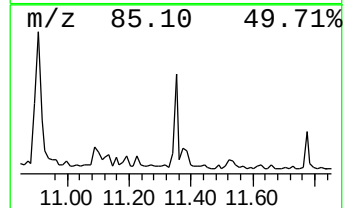
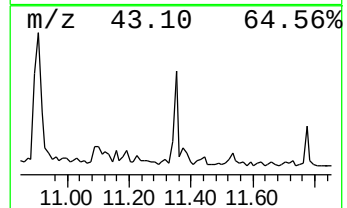
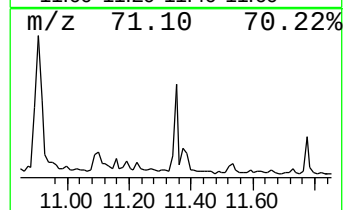
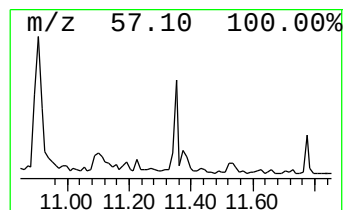
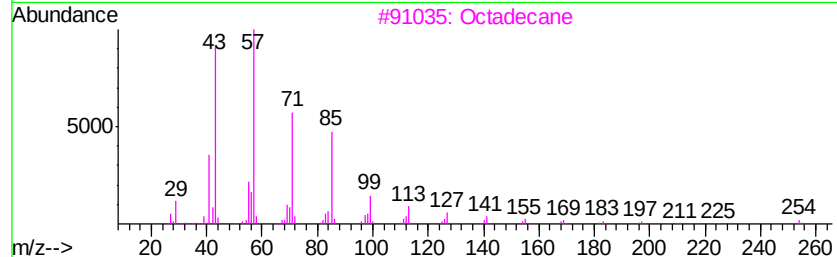
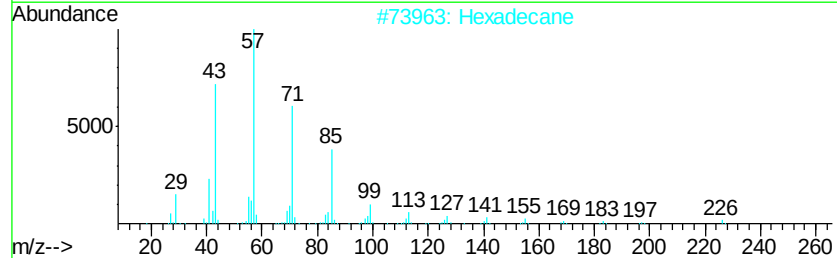
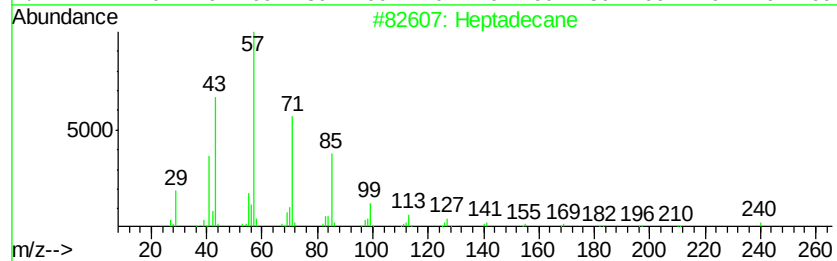
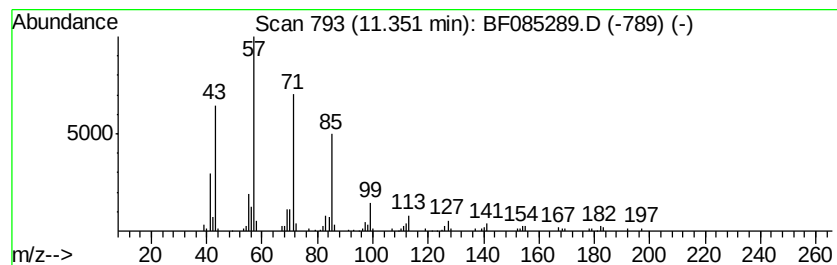
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 Heptadecane Concentration Rank 14

| R.T.      | EstConc      | Area   | Relative to ISTD |             | R.T.  |
|-----------|--------------|--------|------------------|-------------|-------|
| 11.35     | 2.50 ng      | 131772 | Phenanthrene-d10 |             | 11.49 |
| Hit# of 5 | Tentative ID | MW     | MolForm          | CAS#        | Qual  |
| 1         | Heptadecane  | 240    | C17H36           | 000629-78-7 | 91    |
| 2         | Hexadecane   | 226    | C16H34           | 000544-76-3 | 91    |
| 3         | Octadecane   | 254    | C18H38           | 000593-45-3 | 91    |
| 4         | Pentadecane  | 212    | C15H32           | 000629-62-9 | 90    |
| 5         | Tetradecane  | 198    | C14H30           | 000629-59-4 | 90    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-43-COMP-0.5-2.5

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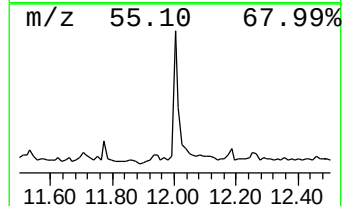
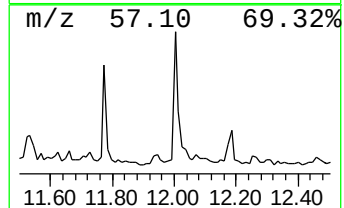
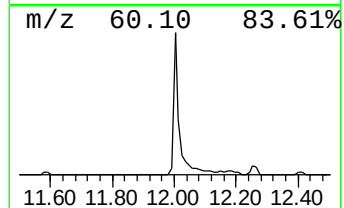
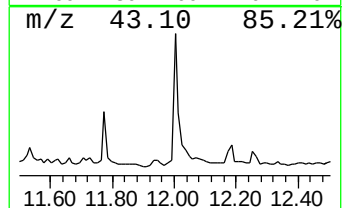
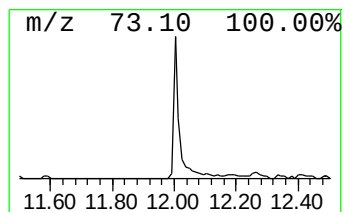
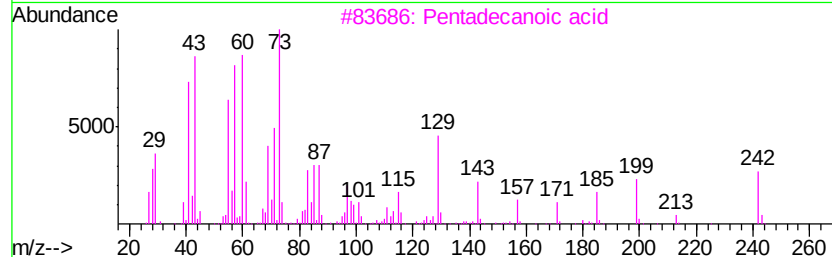
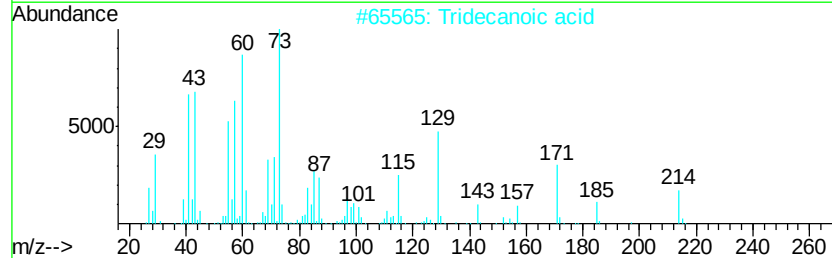
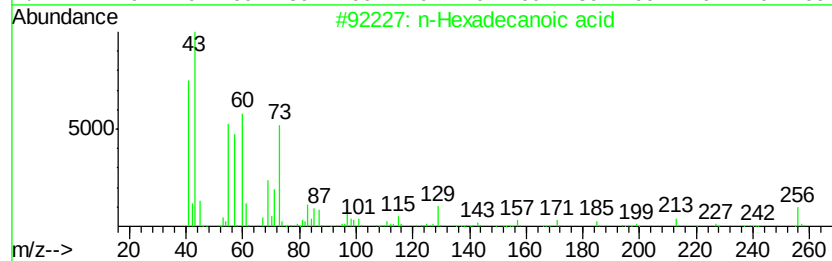
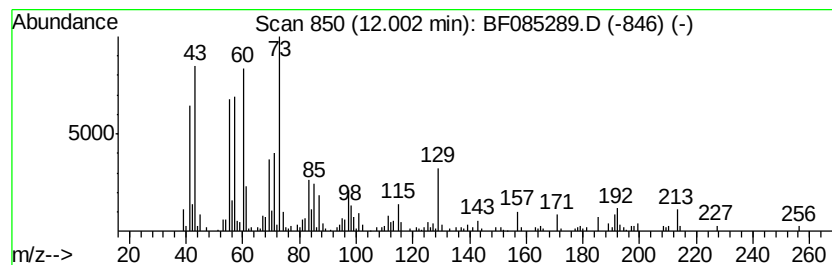
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 n-Hexadecanoic acid Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.00 | 7.92 ng | 417828 | Phenanthrene-d10 | 11.49 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 95   |
| 2    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 90   |
| 3    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 86   |
| 4    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 72   |
| 5    |    | Tetradecane         | 198 | C14H30   | 000629-59-4 | 25   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

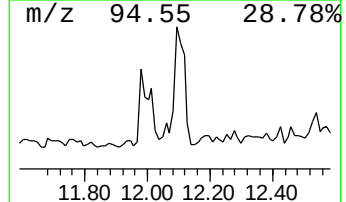
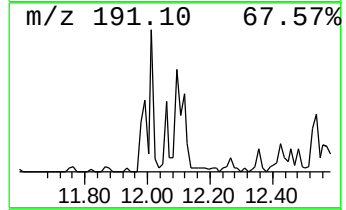
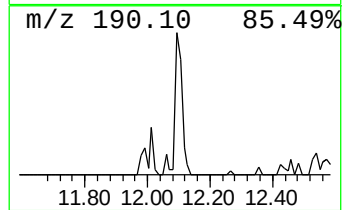
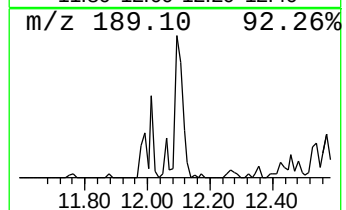
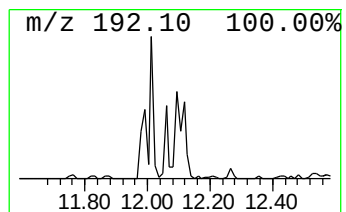
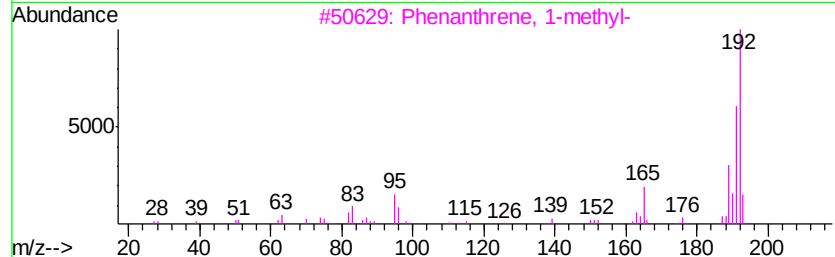
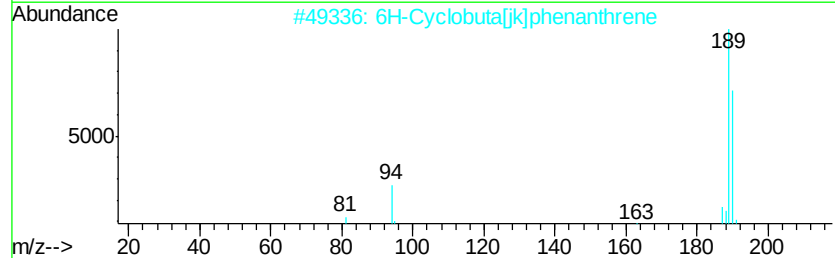
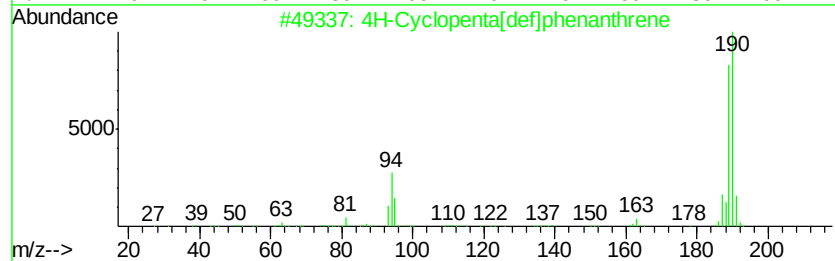
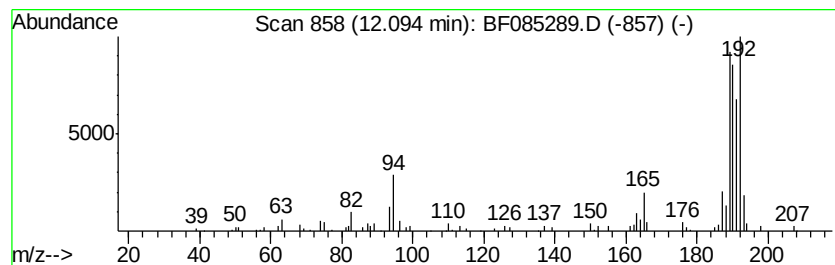
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SB-43-COMP-0.5-2.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 11

| R.T.      | EstConc                        | Area   | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------|--------|------------------|-------------|------|
| 12.09     | 3.02 ng                        | 159632 | Phenanthrene-d10 | 11.49       |      |
| Hit# of 5 | Tentative ID                   | MW     | MolForm          | CAS#        | Qual |
| 1         | 4H-Cyclopenta[def]phenanthrene | 190    | C15H10           | 000203-64-5 | 53   |
| 2         | 6H-Cyclobuta[jk]phenanthrene   | 190    | C15H10           | 083469-43-6 | 49   |
| 3         | Phenanthrene, 1-methyl-        | 192    | C15H12           | 000832-69-9 | 46   |
| 4         | Phenanthrene, 2-methyl-        | 192    | C15H12           | 002531-84-2 | 45   |
| 5         | Anthracene, 1-methyl-          | 192    | C15H12           | 000610-48-0 | 42   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-43-COMP-0.5-2.5

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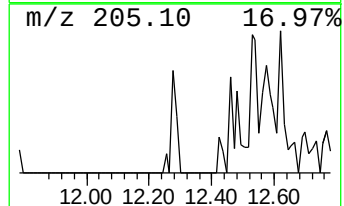
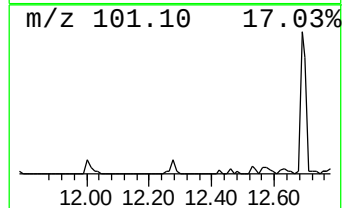
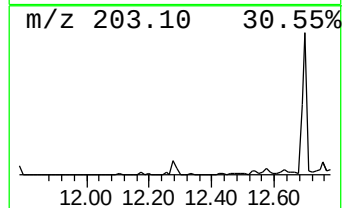
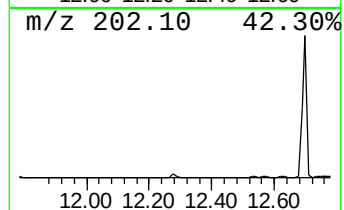
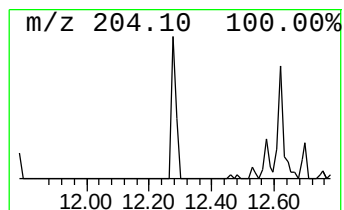
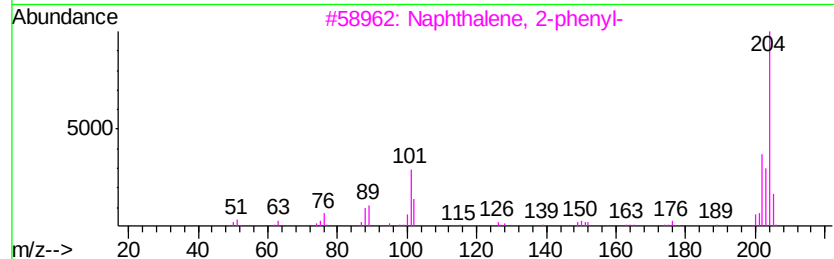
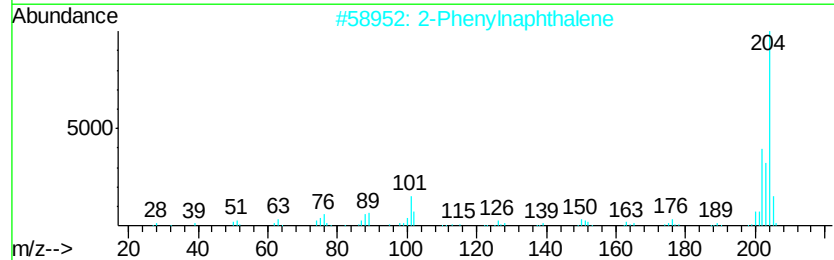
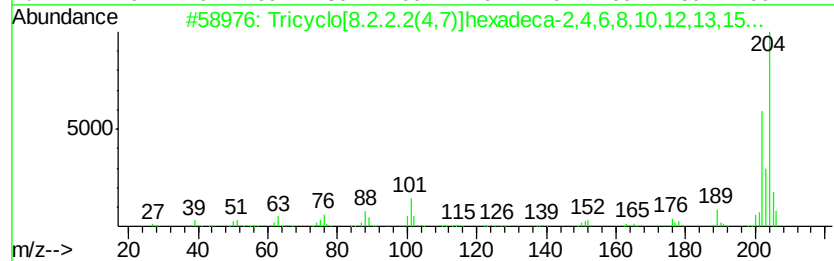
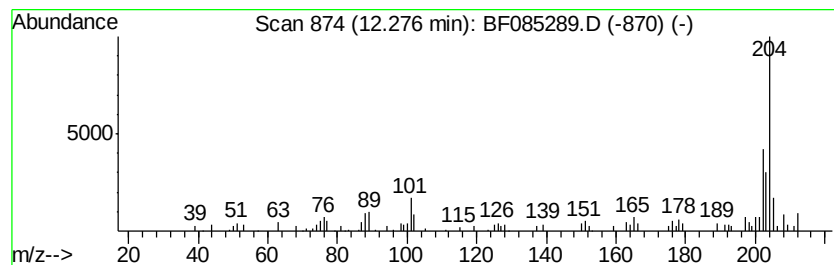
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 Tricyclo[8.2.2.2(4,7)]hexad... Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.28 | 2.32 ng | 122455 | Phenanthrene-d10 | 11.49 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12    | 006572-60-7 | 83   |
| 2    |    | 2-Phenylnaphthalene                 | 204 | C16H12    | 035465-71-5 | 70   |
| 3    |    | Naphthalene, 2-phenyl-              | 204 | C16H12    | 000612-94-2 | 70   |
| 4    |    | 5H-Dibenzo[a,d]cycloheptene, 5-m... | 204 | C16H12    | 002975-79-3 | 60   |
| 5    |    | 9,10-Bis(bromomethyl)anthracene     | 362 | C16H12Br2 | 034373-96-1 | 58   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-43-COMP-0.5-2.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

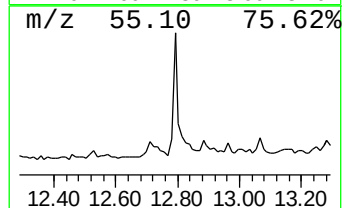
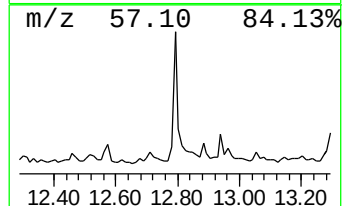
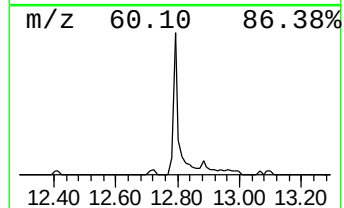
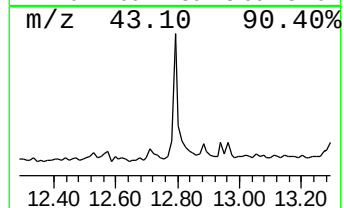
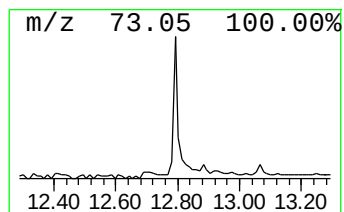
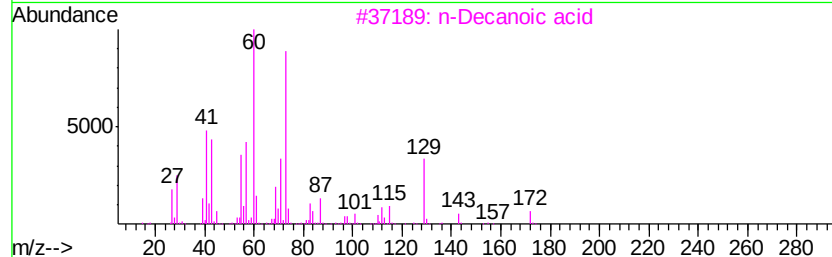
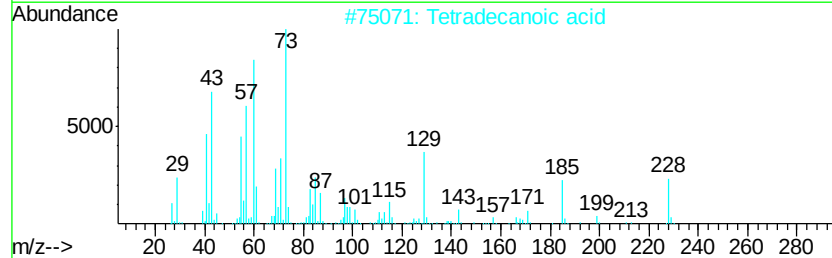
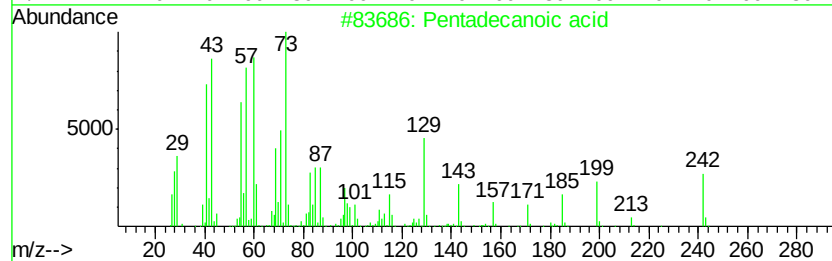
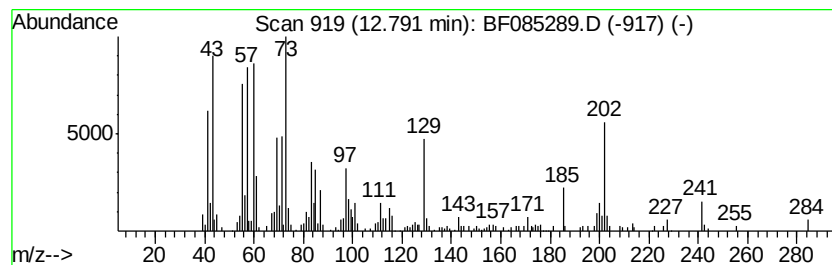
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 Pentadecanoic acid Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.79 | 3.87 ng | 204432 | Phenanthrene-d10 | 11.49 |

| Hit# | of | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------|-----|----------|-------------|------|
| 1    | 5  | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 89   |
| 2    |    | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 72   |
| 3    |    | n-Decanoic acid    | 172 | C10H20O2 | 000334-48-5 | 64   |
| 4    |    | Tridecanoic acid   | 214 | C13H26O2 | 000638-53-9 | 62   |
| 5    |    | Dodecanoic acid    | 200 | C12H24O2 | 000143-07-7 | 50   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

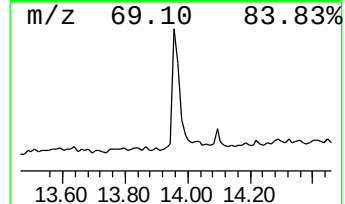
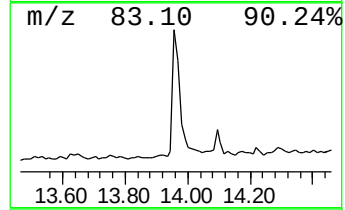
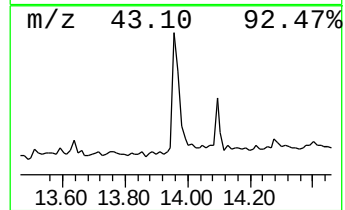
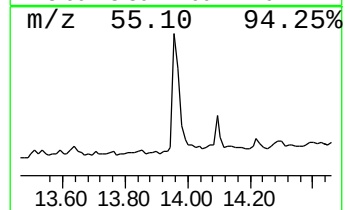
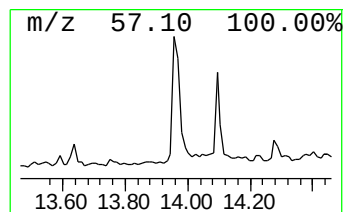
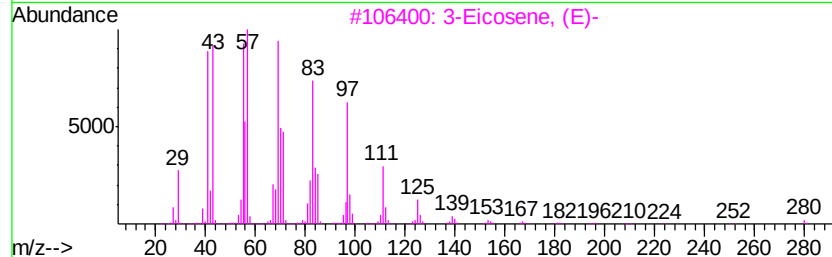
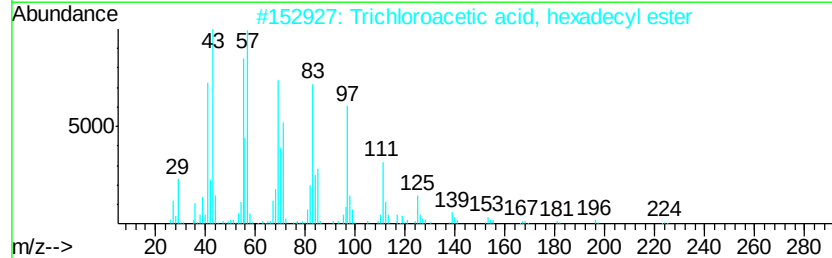
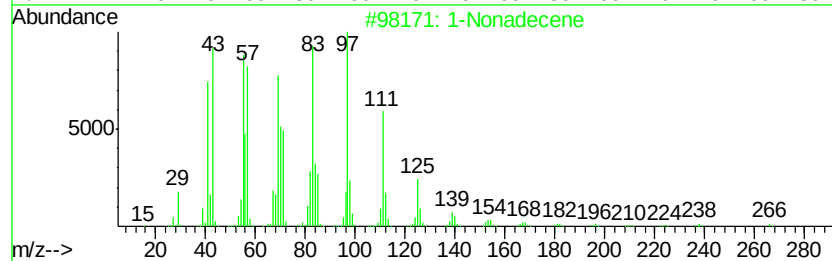
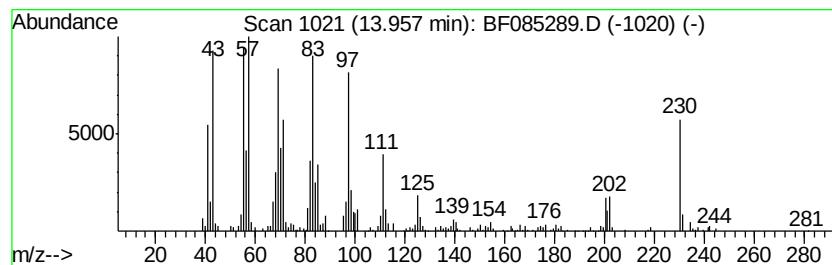
Instrument :  
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ClientSampleId :  
SB-43-COMP-0.5-2.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 1-Nonadecene Concentration Rank 8

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 13.96     | 3.67 ng                             | 314783 | Chrysene-d12     | 14.12        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1-Nonadecene                        | 266    | C19H38           | 018435-45-5  | 90   |
| 2         | Trichloroacetic acid, hexadecyl ... | 386    | C18H33Cl3O2      | 074339-54-1  | 90   |
| 3         | 3-Eicosene, (E)-                    | 280    | C20H40           | 074685-33-9  | 83   |
| 4         | 1-Octadecanol                       | 270    | C18H38O          | 000112-92-5  | 83   |
| 5         | Dichloroacetic acid, heptadecyl ... | 366    | C19H36Cl2O2      | 1000282-98-2 | 81   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-43-COMP-0.5-2.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

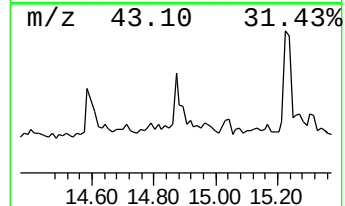
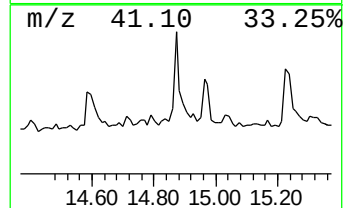
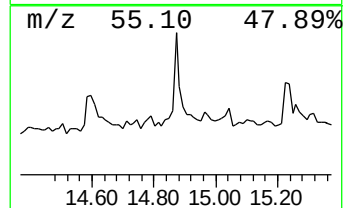
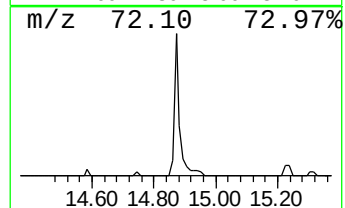
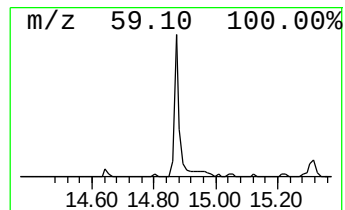
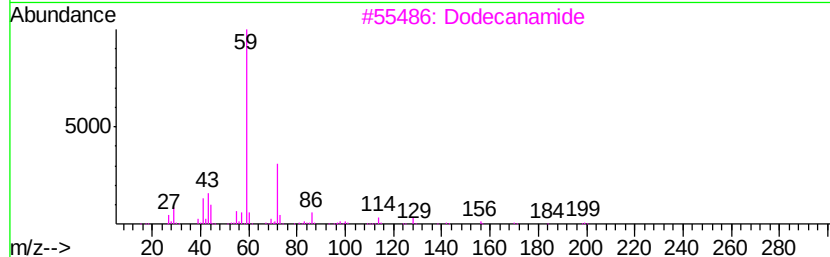
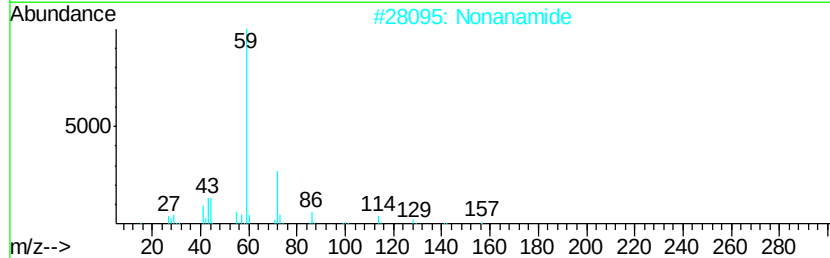
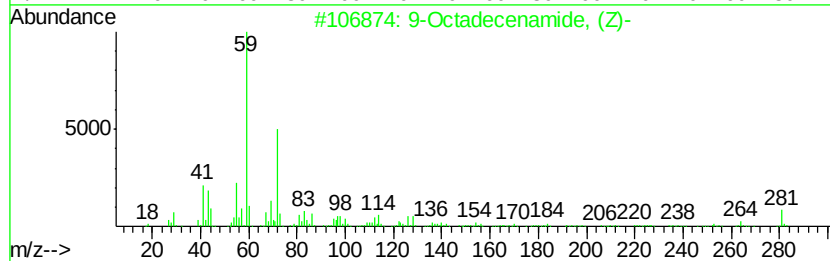
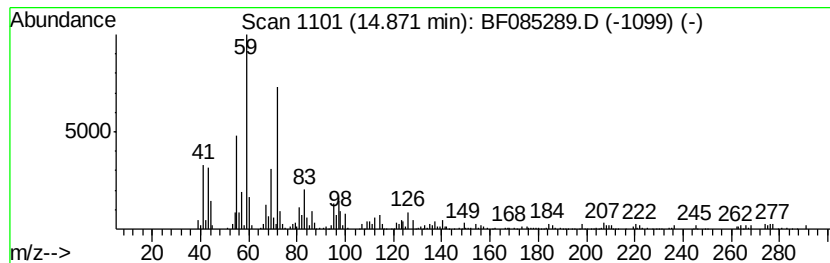
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 9-Octadecenamide, (Z)- Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.87 | 2.61 ng | 132280 | Perylene-d12     | 15.59 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------------|-----|----------|-------------|------|
| 1    |    |   | 9-Octadecenamide, (Z)- | 281 | C18H35NO | 000301-02-0 | 53   |
| 2    |    |   | Nonanamide             | 157 | C9H19NO  | 001120-07-6 | 49   |
| 3    |    |   | Dodecanamide           | 199 | C12H25NO | 001120-16-7 | 47   |
| 4    |    |   | Octanamide             | 143 | C8H17NO  | 000629-01-6 | 47   |
| 5    |    |   | 2-Methyl-2-decanol     | 172 | C11H24O  | 003396-02-9 | 47   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

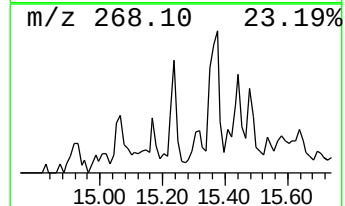
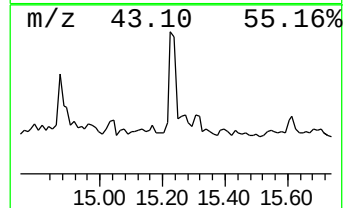
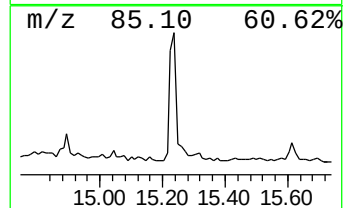
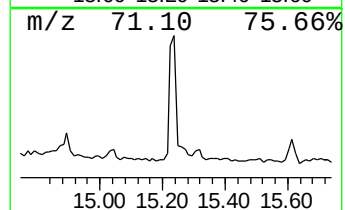
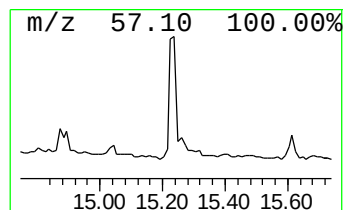
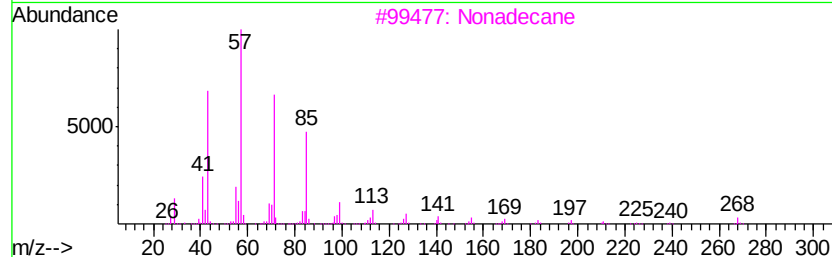
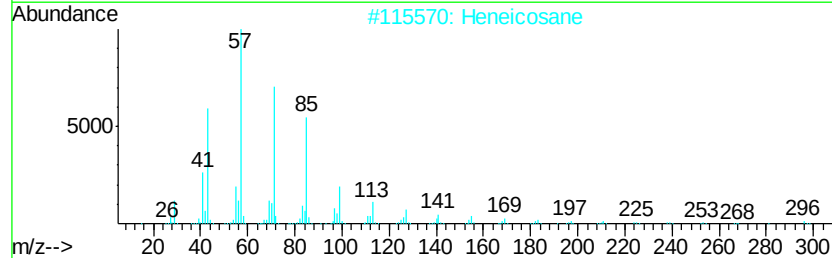
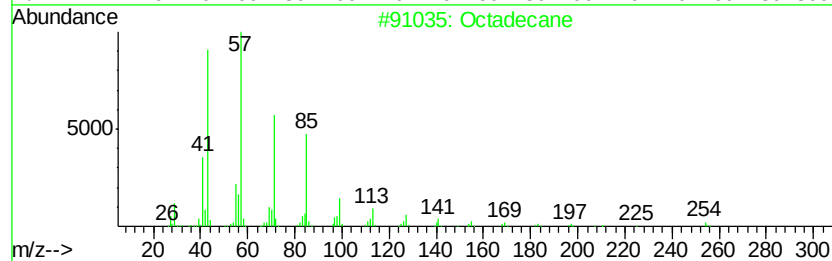
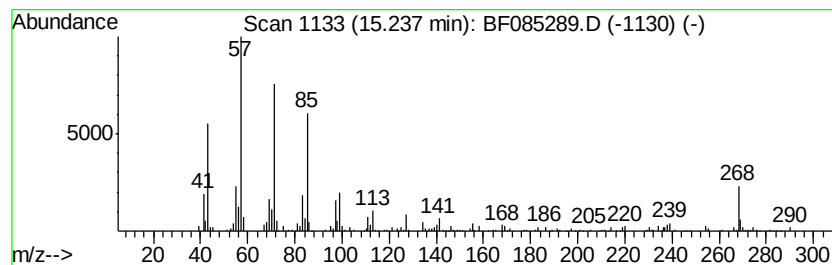
Instrument :  
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ClientSampleId :  
SB-43-COMP-0.5-2.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 Octadecane Concentration Rank 12

| R.T.      | EstConc                | Area   | Relative to ISTD |             | R.T.  |
|-----------|------------------------|--------|------------------|-------------|-------|
| 15.24     | 3.01 ng                | 152980 | Perylene-d12     |             | 15.59 |
| Hit# of 5 | Tentative ID           | MW     | MolForm          | CAS#        | Qual  |
| 1         | Octadecane             | 254    | C18H38           | 000593-45-3 | 90    |
| 2         | Heneicosane            | 296    | C21H44           | 000629-94-7 | 87    |
| 3         | Nonadecane             | 268    | C19H40           | 000629-92-5 | 83    |
| 4         | Pentadecane, 8-heptyl- | 310    | C22H46           | 071005-15-7 | 83    |
| 5         | Eicosane, 10-methyl-   | 296    | C21H44           | 054833-23-7 | 80    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-43-COMP-0.5-2.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

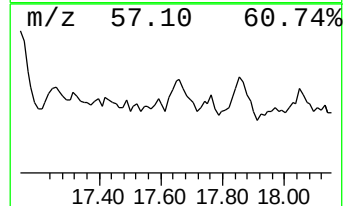
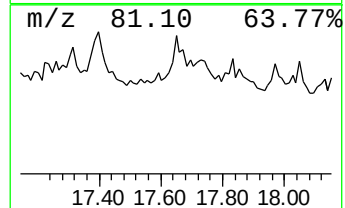
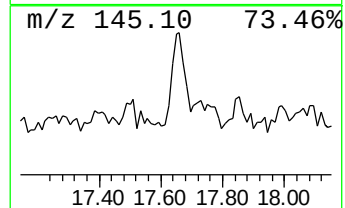
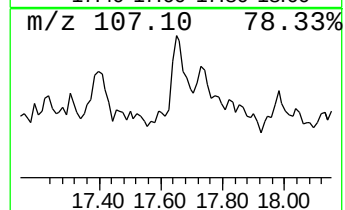
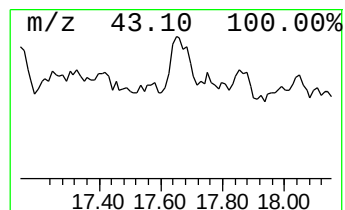
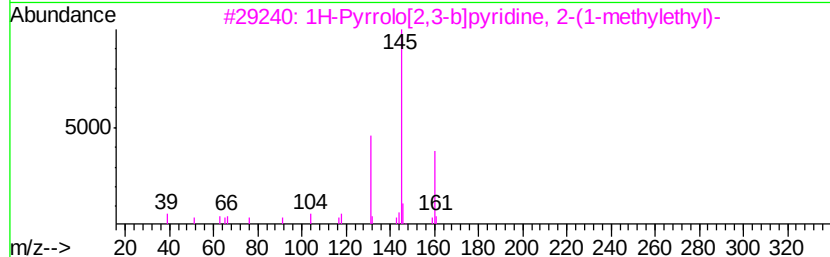
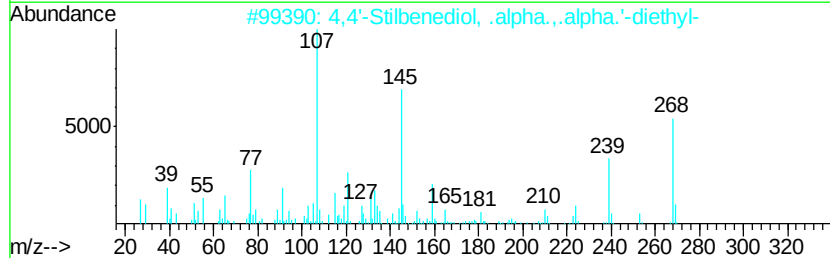
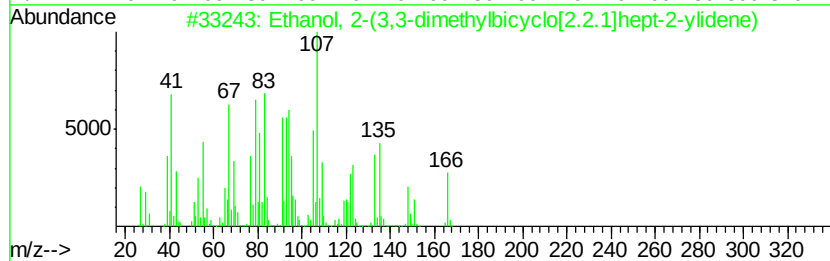
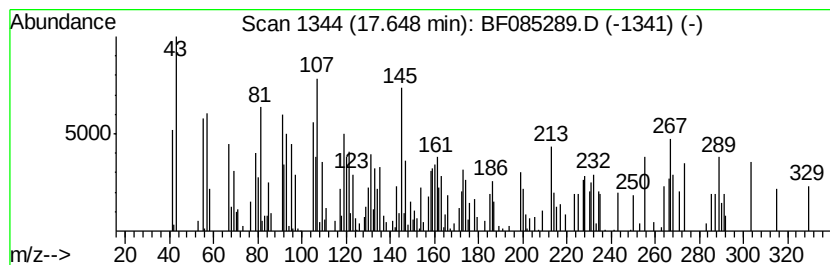
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 unknown17.65 Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.65 | 3.66 ng | 185820 | Perylene-d12     | 15.59 |

| Hit# | of | Tentative ID                         | MW  | MolForm   | CAS#         | Qual |
|------|----|--------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Ethanol, 2-(3,3-dimethylbicyclo[...] | 166 | C11H18O   | 002226-05-3  | 44   |
| 2    |    | 4,4'-Stilbenediol, .alpha.,.alph...  | 268 | C18H20O2  | 006898-97-1  | 25   |
| 3    |    | 1H-Pyrrolo[2,3-b]pyridine, 2-(1-...  | 160 | C10H12N2  | 027257-18-7  | 11   |
| 4    |    | Benzenemethanol, .alpha.-(1-meth...  | 150 | C10H14O   | 014898-86-3  | 11   |
| 5    |    | N-Methylproline, benzyl ester        | 219 | C13H17NO2 | 1000130-46-8 | 11   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-43-COMP-0.5-2.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

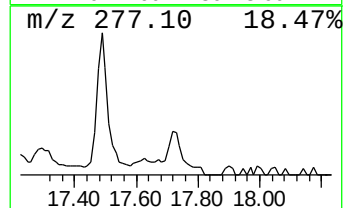
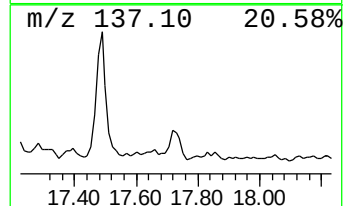
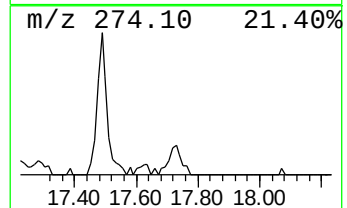
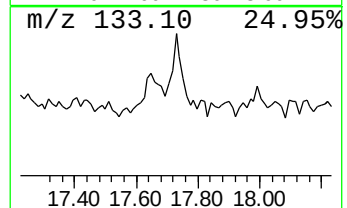
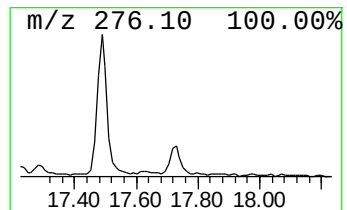
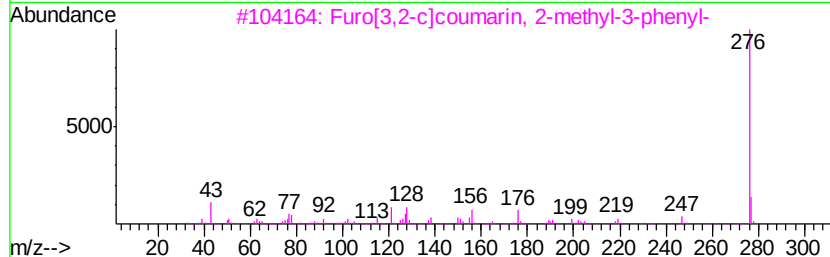
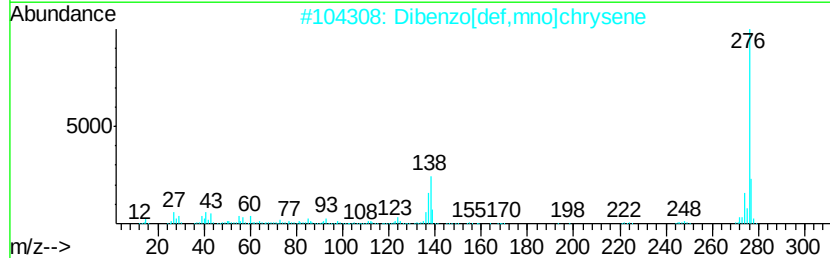
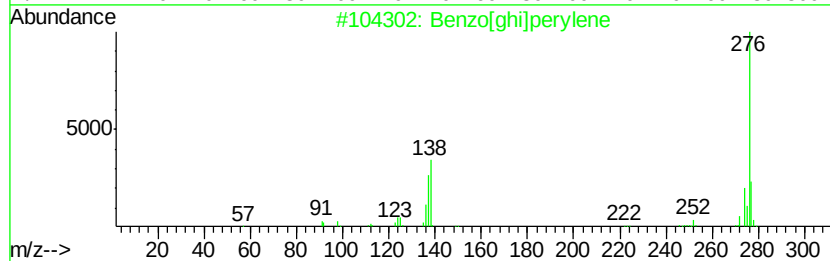
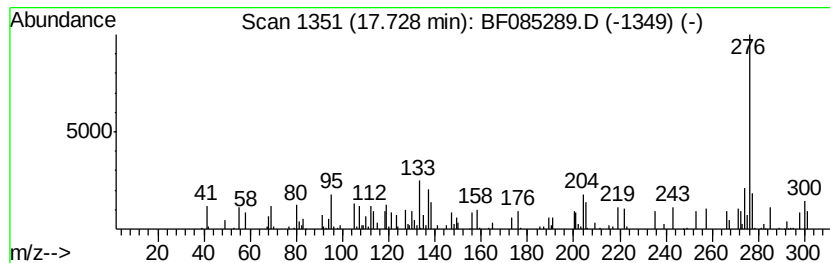
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 unknown17.73 Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.73 | 2.22 ng | 112550 | Perylene-d12     | 15.59 |

| Hit# | of | Tentative ID                         | MW  | MolForm     | CAS#         | Qual |
|------|----|--------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Benzo[ghi]perylene                   | 276 | C22H12      | 000191-24-2  | 43   |
| 2    |    | Dibenzo[def,mno]chrysene             | 276 | C22H12      | 000191-26-4  | 43   |
| 3    |    | Furo[3,2-c]coumarin, 2-methyl-3-...  | 276 | C18H12O3    | 010375-42-5  | 41   |
| 4    |    | Phenol, 2-(4-chlorophenyl)iminome... | 276 | C13H9ClN2O3 | 1000262-90-3 | 38   |
| 5    |    | Indeno[1,2,3-cd]pyrene               | 276 | C22H12      | 000193-39-5  | 38   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF030816\  
Data File : BF085289.D  
Acq On : 9 Mar 2016 6:58  
Operator : UM/SJ  
Sample : H1712-06  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-43-COMP-0.5-2.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name      | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|-----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                       |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2-Pentanone, 4-hy...  | 5.17  | 21.9    | ng    | 1114280  | 1                     | 6.96  | 1020000 | 20.0 |
| unknown6.73           | 6.73  | 85.8    | ng    | 4375320  | 1                     | 6.96  | 1020000 | 20.0 |
| Tetradecane           | 9.93  | 2.4     | ng    | 143137   | 3                     | 10.00 | 1201960 | 20.0 |
| Decane, 2,3,5-tri...  | 10.42 | 3.3     | ng    | 198996   | 3                     | 10.00 | 1201960 | 20.0 |
| Docosane              | 10.65 | 2.3     | ng    | 138398   | 3                     | 10.00 | 1201960 | 20.0 |
| Heneicosane           | 10.90 | 5.5     | ng    | 290379   | 4                     | 11.49 | 1055670 | 20.0 |
| Heptadecane           | 11.35 | 2.5     | ng    | 131772   | 4                     | 11.49 | 1055670 | 20.0 |
| n-Hexadecanoic acid   | 12.00 | 7.9     | ng    | 417828   | 4                     | 11.49 | 1055670 | 20.0 |
| 4H-Cyclopenta[def...  | 12.09 | 3.0     | ng    | 159632   | 4                     | 11.49 | 1055670 | 20.0 |
| Tricyclo[8.2.2.2(...  | 12.28 | 2.3     | ng    | 122455   | 4                     | 11.49 | 1055670 | 20.0 |
| Pentadecanoic acid    | 12.79 | 3.9     | ng    | 204432   | 4                     | 11.49 | 1055670 | 20.0 |
| 1-Nonadecene          | 13.96 | 3.7     | ng    | 314783   | 5                     | 14.12 | 1714970 | 20.0 |
| 9-Octadecenamide, ... | 14.87 | 2.6     | ng    | 132280   | 6                     | 15.59 | 1015440 | 20.0 |
| Octadecane            | 15.24 | 3.0     | ng    | 152980   | 6                     | 15.59 | 1015440 | 20.0 |
| unknown17.65          | 17.65 | 3.7     | ng    | 185820   | 6                     | 15.59 | 1015440 | 20.0 |
| unknown17.73          | 17.73 | 2.2     | ng    | 112550   | 6                     | 15.59 | 1015440 | 20.0 |