

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
 Data File : BF085645.D  
 Acq On : 22 Mar 2016 4:21  
 Operator : UM/SJ  
 Sample : H1840-03  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PJ-SB-11-1.5-2.0

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-BF031816.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      | 4.504    | 191        | 194      | 197       | rVB   | 130237      | 150536     | 5.33%        | 0.475%     |
| 2      | 5.030    | 237        | 240      | 249       | rVB   | 226881      | 322826     | 11.43%       | 1.019%     |
| 3      | 5.441    | 274        | 276      | 279       | rBV   | 2220598     | 2485994    | 87.98%       | 7.847%     |
| 4      | 6.481    | 364        | 367      | 369       | rBV   | 2376117     | 2825562    | 100.00%      | 8.919%     |
| 5      | 6.607    | 376        | 378      | 381       | rBV   | 2513239     | 2598741    | 91.97%       | 8.203%     |
| 6      | 6.836    | 396        | 398      | 401       | rBV   | 576195      | 684185     | 24.21%       | 2.160%     |
| 7      | 6.996    | 410        | 412      | 415       | rBV   | 2314719     | 2164213    | 76.59%       | 6.831%     |
| 8      | 7.407    | 445        | 448      | 450       | rBV   | 1833624     | 1789872    | 63.35%       | 5.650%     |
| 9      | 7.499    | 454        | 456      | 459       | rVV   | 37378       | 54057      | 1.91%        | 0.171%     |
| 10     | 8.127    | 508        | 511      | 513       | rBV   | 1040834     | 914857     | 32.38%       | 2.888%     |
| 11     | 9.202    | 602        | 605      | 608       | rBV   | 3032285     | 2730677    | 96.64%       | 8.620%     |
| 12     | 9.282    | 610        | 612      | 617       | rVB   | 24018       | 35419      | 1.25%        | 0.112%     |
| 13     | 9.533    | 633        | 634      | 638       | rBV3  | 24110       | 51367      | 1.82%        | 0.162%     |
| 14     | 9.602    | 638        | 640      | 642       | rBV   | 338288      | 470022     | 16.63%       | 1.484%     |
| 15     | 9.739    | 650        | 652      | 654       | rBV   | 83290       | 67849      | 2.40%        | 0.214%     |
| 16     | 9.819    | 656        | 659      | 660       | rVB   | 45682       | 62574      | 2.21%        | 0.198%     |
| 17     | 9.876    | 662        | 664      | 666       | rVV   | 883655      | 835354     | 29.56%       | 2.637%     |
| 18     | 10.082   | 680        | 682      | 683       | rBV2  | 32092       | 36901      | 1.31%        | 0.116%     |
| 19     | 10.127   | 683        | 686      | 688       | rVB3  | 18661       | 38630      | 1.37%        | 0.122%     |
| 20     | 10.196   | 688        | 692      | 693       | rBV2  | 20777       | 30011      | 1.06%        | 0.095%     |
| 21     | 10.288   | 698        | 700      | 701       | rBV   | 36150       | 46632      | 1.65%        | 0.147%     |
| 22     | 10.310   | 701        | 702      | 704       | rVB   | 107024      | 77854      | 2.76%        | 0.246%     |
| 23     | 10.425   | 710        | 712      | 715       | rVB2  | 24058       | 41254      | 1.46%        | 0.130%     |
| 24     | 10.528   | 719        | 721      | 724       | rVV2  | 36427       | 60197      | 2.13%        | 0.190%     |
| 25     | 10.665   | 730        | 733      | 736       | rBV   | 1447866     | 1468540    | 51.97%       | 4.636%     |
| 26     | 10.790   | 742        | 744      | 748       | rBV2  | 96583       | 174041     | 6.16%        | 0.549%     |
| 27     | 10.973   | 758        | 760      | 761       | rBV   | 33396       | 39000      | 1.38%        | 0.123%     |
| 28     | 11.122   | 769        | 773      | 776       | rVB3  | 20897       | 37201      | 1.32%        | 0.117%     |
| 29     | 11.236   | 779        | 783      | 784       | rBV2  | 58893       | 84633      | 3.00%        | 0.267%     |
| 30     | 11.259   | 784        | 785      | 788       | rVB   | 46601       | 44677      | 1.58%        | 0.141%     |
| 31     | 11.362   | 792        | 794      | 795       | rBV   | 846419      | 761674     | 26.96%       | 2.404%     |
| 32     | 11.430   | 798        | 800      | 802       | rVB2  | 42054       | 65269      | 2.31%        | 0.206%     |
| 33     | 11.476   | 802        | 804      | 805       | rBV2  | 24040       | 28862      | 1.02%        | 0.091%     |
| 34     | 11.590   | 811        | 814      | 817       | rBV3  | 27121       | 59043      | 2.09%        | 0.186%     |

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 Peak Location: TOP

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Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |           |         |         |        |        |
|----|--------|------|------|-----------|---------|---------|--------|--------|
| 35 | 11.659 | 818  | 820  | 821 rVV   | 62176   | 57320   | 2.03%  | 0.181% |
| 36 | 11.693 | 821  | 823  | 826 rVB   | 30183   | 34859   | 1.23%  | 0.110% |
| 37 | 11.865 | 835  | 838  | 839 rBV   | 91747   | 91657   | 3.24%  | 0.289% |
| 38 | 11.888 | 839  | 840  | 843 rVV   | 250183  | 267543  | 9.47%  | 0.845% |
| 39 | 11.968 | 846  | 847  | 852 rVB   | 115832  | 207709  | 7.35%  | 0.656% |
| 40 | 12.059 | 852  | 855  | 857 rBV2  | 26685   | 48048   | 1.70%  | 0.152% |
| 41 | 12.162 | 860  | 864  | 865 rBV   | 42476   | 68303   | 2.42%  | 0.216% |
| 42 | 12.311 | 874  | 877  | 878 rBV2  | 28589   | 48898   | 1.73%  | 0.154% |
| 43 | 12.333 | 878  | 879  | 883 rVB   | 31401   | 51737   | 1.83%  | 0.163% |
| 44 | 12.413 | 883  | 886  | 887 rBV   | 97975   | 119045  | 4.21%  | 0.376% |
| 45 | 12.448 | 887  | 889  | 892 rVB   | 69996   | 106557  | 3.77%  | 0.336% |
| 46 | 12.493 | 892  | 893  | 897 rVB   | 49713   | 65646   | 2.32%  | 0.207% |
| 47 | 12.573 | 897  | 900  | 902 rBV   | 434272  | 422320  | 14.95% | 1.333% |
| 48 | 12.619 | 902  | 904  | 907 rVB2  | 21142   | 38630   | 1.37%  | 0.122% |
| 49 | 12.676 | 907  | 909  | 915 rBV4  | 105975  | 217929  | 7.71%  | 0.688% |
| 50 | 12.802 | 915  | 920  | 921 rVV   | 344045  | 486921  | 17.23% | 1.537% |
| 51 | 12.859 | 923  | 925  | 927 rVB   | 37077   | 49950   | 1.77%  | 0.158% |
| 52 | 12.951 | 930  | 933  | 935 rVB   | 1310876 | 1831650 | 64.82% | 5.782% |
| 53 | 13.019 | 936  | 939  | 943 rBV3  | 62917   | 108778  | 3.85%  | 0.343% |
| 54 | 13.122 | 945  | 948  | 950 rVB   | 75799   | 146945  | 5.20%  | 0.464% |
| 55 | 13.179 | 950  | 953  | 956 rBV2  | 75770   | 160379  | 5.68%  | 0.506% |
| 56 | 13.225 | 956  | 957  | 959 rBV   | 49472   | 67400   | 2.39%  | 0.213% |
| 57 | 13.328 | 961  | 966  | 967 rBV   | 46023   | 73242   | 2.59%  | 0.231% |
| 58 | 13.351 | 967  | 968  | 970 rBV   | 60532   | 57204   | 2.02%  | 0.181% |
| 59 | 13.431 | 972  | 975  | 980 rBV   | 185796  | 501481  | 17.75% | 1.583% |
| 60 | 13.522 | 980  | 983  | 985 rBV2  | 103311  | 172785  | 6.12%  | 0.545% |
| 61 | 13.636 | 990  | 993  | 995 rVB2  | 56674   | 75208   | 2.66%  | 0.237% |
| 62 | 13.739 | 999  | 1002 | 1004 rBV2 | 56159   | 110058  | 3.90%  | 0.347% |
| 63 | 13.842 | 1009 | 1011 | 1016 rVB  | 460246  | 586809  | 20.77% | 1.852% |
| 64 | 13.991 | 1021 | 1024 | 1030 rBV2 | 713536  | 1157066 | 40.95% | 3.652% |
| 65 | 14.105 | 1030 | 1034 | 1036 rBV4 | 54004   | 132034  | 4.67%  | 0.417% |
| 66 | 14.162 | 1037 | 1039 | 1042 rVB  | 171875  | 179691  | 6.36%  | 0.567% |
| 67 | 14.379 | 1054 | 1058 | 1060 rBV  | 85470   | 146913  | 5.20%  | 0.464% |
| 68 | 14.414 | 1060 | 1061 | 1063 rVV2 | 52551   | 61464   | 2.18%  | 0.194% |
| 69 | 14.471 | 1063 | 1066 | 1068 rVV3 | 151615  | 232839  | 8.24%  | 0.735% |
| 70 | 14.505 | 1068 | 1069 | 1073 rVB2 | 58230   | 87370   | 3.09%  | 0.276% |
| 71 | 14.768 | 1090 | 1092 | 1093 rBV  | 99046   | 111953  | 3.96%  | 0.353% |

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Misc :  
ALS Vial : 28 Sample Multiplier: 1

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BNA\_F  
ClientSampleId :  
PJ-SB-11-1.5-2.0

Integration Parameters: rteint.p  
Integrator: RTE  
Smoothing : ON Filtering: 5  
Sampling : 1 Min Area: 3 % of largest Peak  
Start Thrs: 0.2 Max Peaks: 100  
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 72 | 15.008 | 1111 | 1113 | 1117 | rBV  | 192067 | 388442 | 13.75% | 1.226% |
| 73 | 15.088 | 1117 | 1120 | 1121 | rBV2 | 76042  | 85182  | 3.01%  | 0.269% |
| 74 | 15.305 | 1136 | 1139 | 1141 | rVB  | 97229  | 156132 | 5.53%  | 0.493% |
| 75 | 15.362 | 1141 | 1144 | 1147 | rVV  | 160288 | 207849 | 7.36%  | 0.656% |
| 76 | 15.419 | 1147 | 1149 | 1159 | rVB  | 495577 | 716056 | 25.34% | 2.260% |
| 77 | 15.637 | 1166 | 1168 | 1170 | rVB2 | 34400  | 46069  | 1.63%  | 0.145% |
| 78 | 15.911 | 1190 | 1192 | 1199 | rBV4 | 41561  | 114891 | 4.07%  | 0.363% |
| 79 | 16.802 | 1268 | 1270 | 1274 | rVB2 | 57724  | 117278 | 4.15%  | 0.370% |
| 80 | 17.225 | 1304 | 1307 | 1314 | rVB  | 54438  | 125312 | 4.43%  | 0.396% |

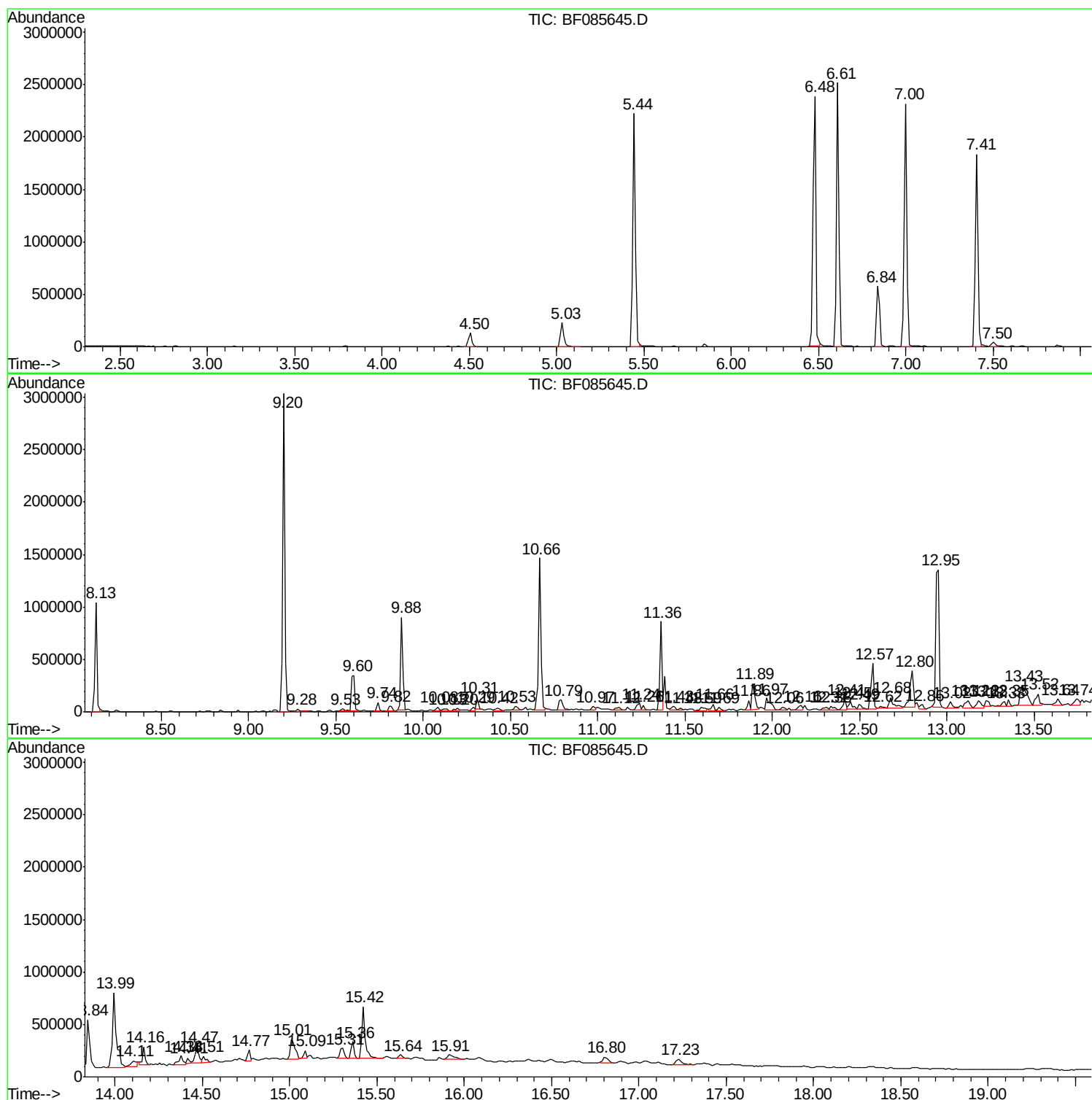
Sum of corrected areas: 31680076

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
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ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.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\BF032116\  
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Misc :  
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Instrument :  
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ClientSampleId :  
PJ-SB-11-1.5-2.0

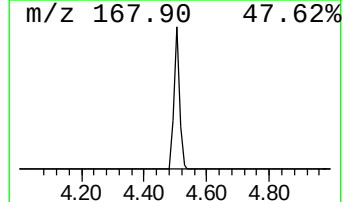
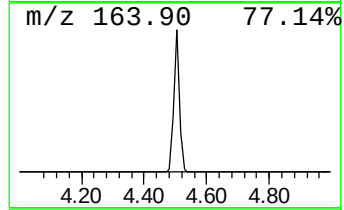
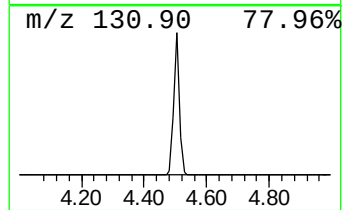
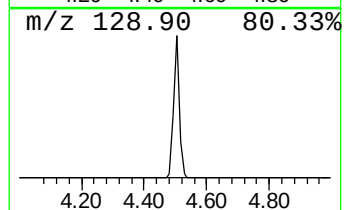
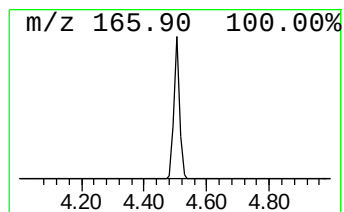
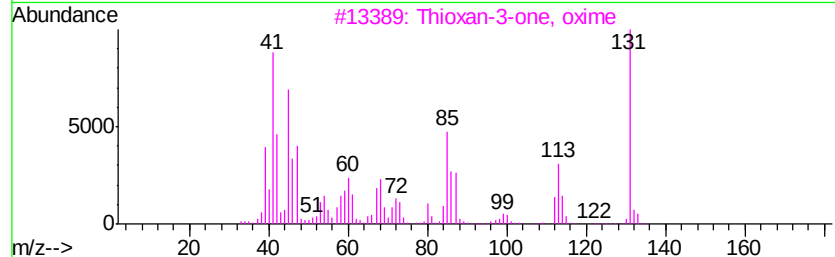
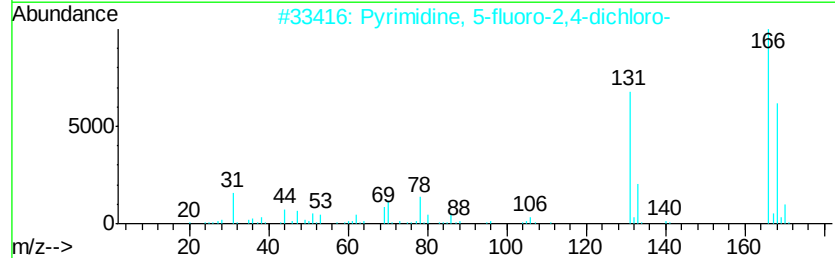
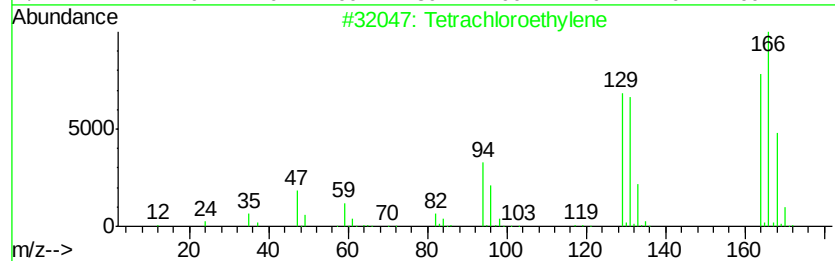
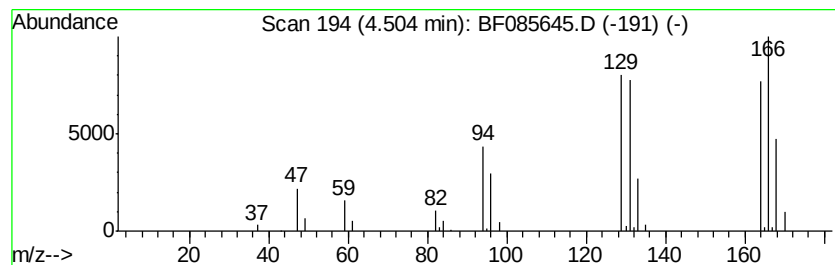
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Tetrachloroethylene Concentration Rank 10

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.50 | 4.40 ng | 150536 | 1,4-Dichlorobenzene-d4 | 6.84 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm   | CAS#        | Qual |
|------|----|---|------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Tetrachloroethylene                | 164 | C2Cl4     | 000127-18-4 | 98   |
| 2    |    |   | Pyrimidine, 5-fluoro-2,4-dichloro- | 166 | C4HCl2FN2 | 002927-71-1 | 25   |
| 3    |    |   | Thioxan-3-one, oxime               | 131 | C5H9NOS   | 058230-51-6 | 10   |
| 4    |    |   | Quinazoline, 4-chloro-             | 164 | C8H5ClN2  | 005190-68-1 | 9    |
| 5    |    |   | 2-Chloroquinoxaline                | 164 | C8H5ClN2  | 001448-87-9 | 9    |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

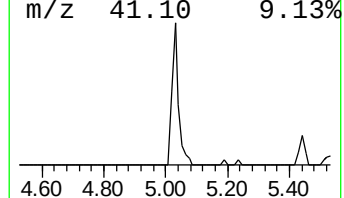
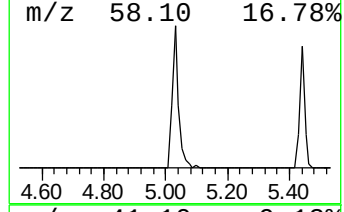
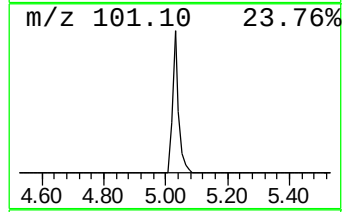
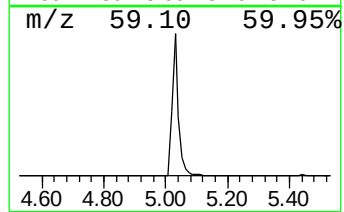
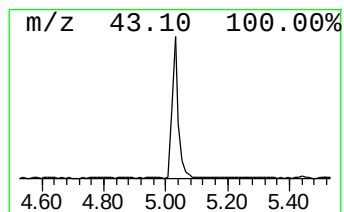
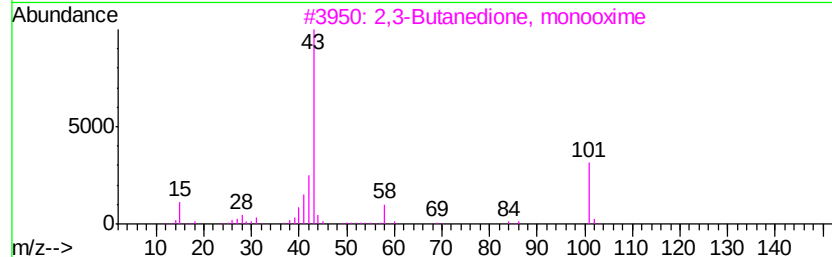
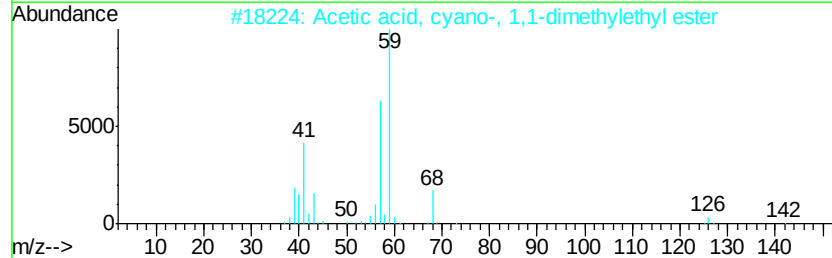
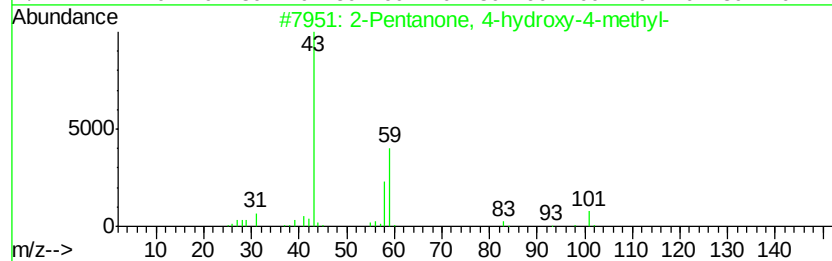
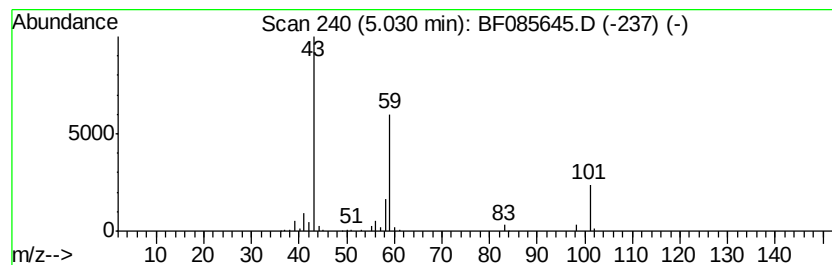
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.03 | 9.44 ng | 322826 | 1,4-Dichlorobenzene-d4 | 6.84 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 72   |
| 2    |    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 3    |    | 2,3-Butanedione, monooxime          | 101 | C4H7NO2  | 000057-71-6 | 17   |
| 4    |    | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |
| 5    |    | 5-Hexen-2-one                       | 98  | C6H10O   | 000109-49-9 | 9    |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

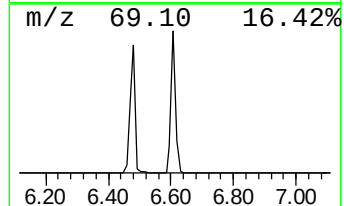
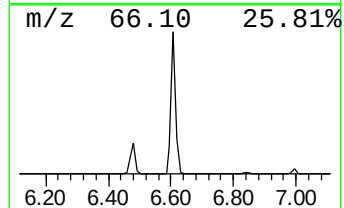
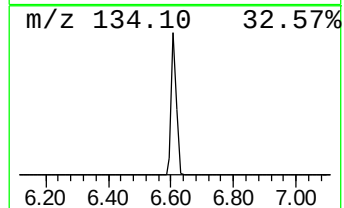
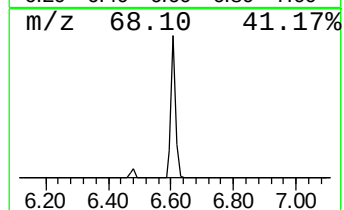
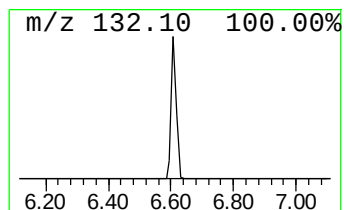
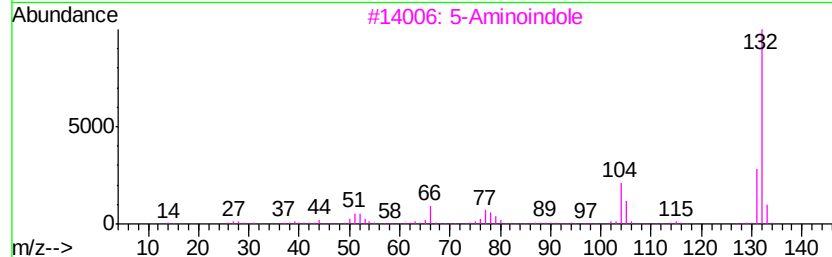
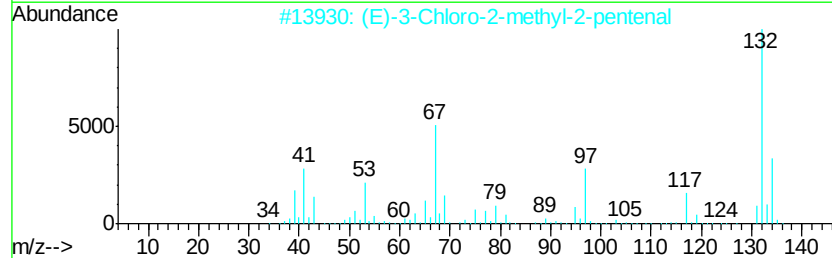
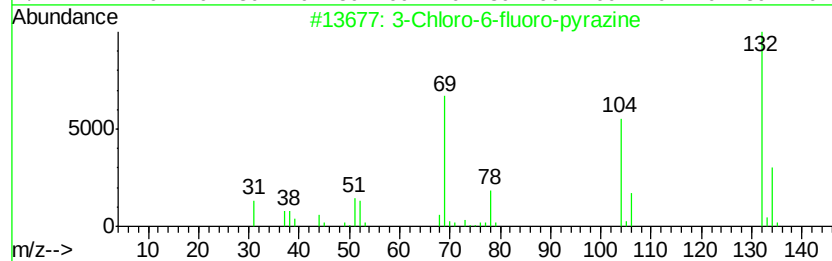
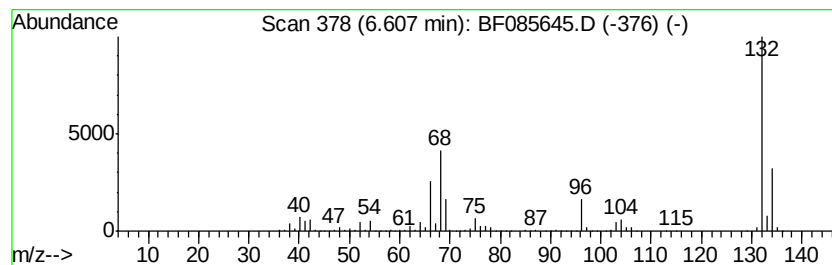
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown6.61 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.61 | 75.97 ng | 2598740 | 1,4-Dichlorobenzene-d4 | 6.84 |

| 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  | 17   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | 4-Chloro-2-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-00-5  | 9    |
| 5    |    | 1-Methylpyrrolo[1,2-a]pyrazine   | 132 | C8H8N2    | 064608-59-9  | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PJ-SB-11-1.5-2.0

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

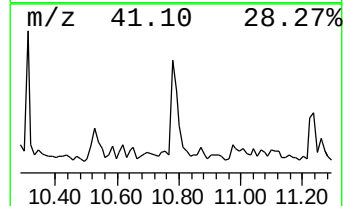
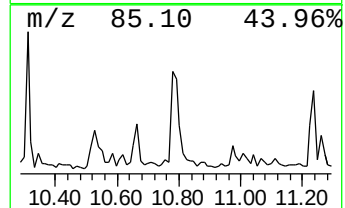
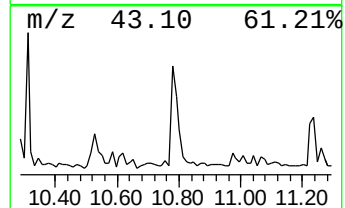
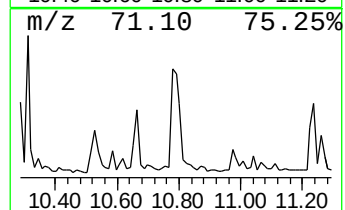
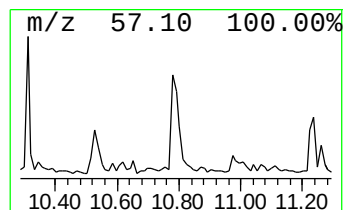
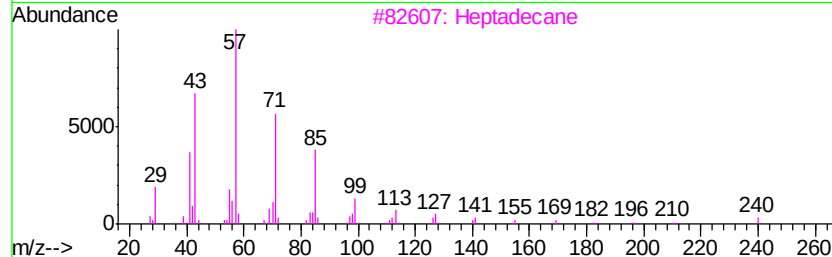
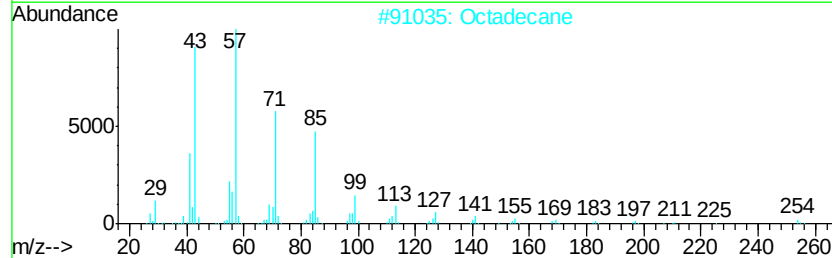
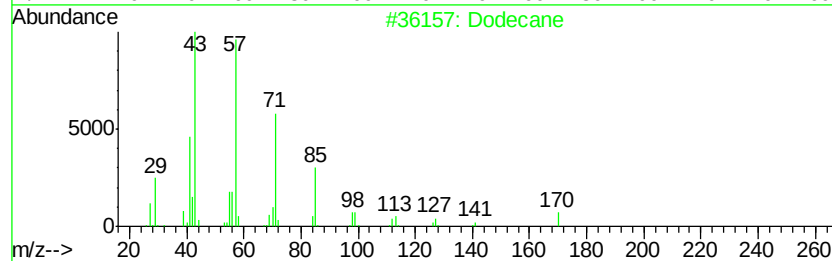
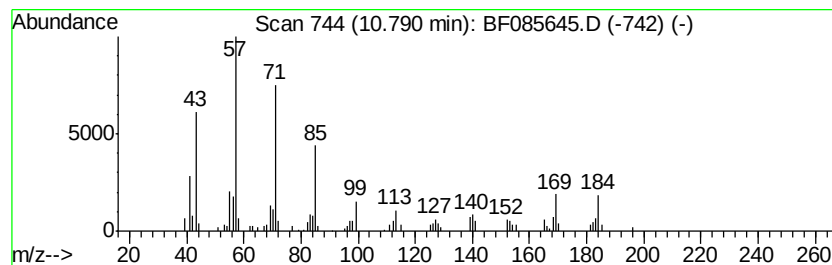
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 Dodecane Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.79 | 4.57 ng | 174041 | Phenanthrene-d10 | 11.36 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane     | 170 | C12H26  | 000112-40-3 | 87   |
| 2    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 83   |
| 3    |    | Heptadecane  | 240 | C17H36  | 000629-78-7 | 83   |
| 4    |    | Octacosane   | 394 | C28H58  | 000630-02-4 | 80   |
| 5    |    | Pentadecane  | 212 | C15H32  | 000629-62-9 | 80   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

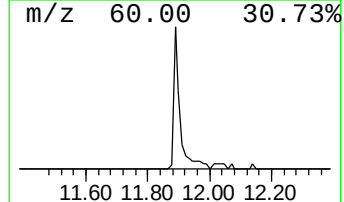
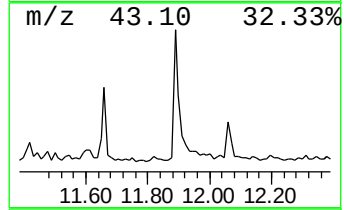
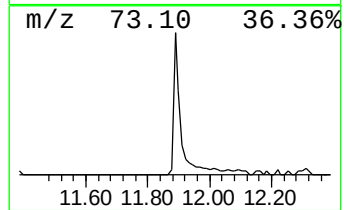
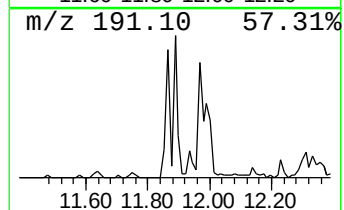
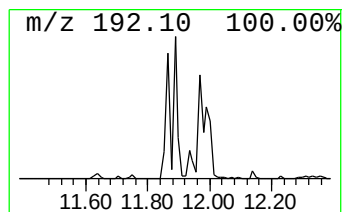
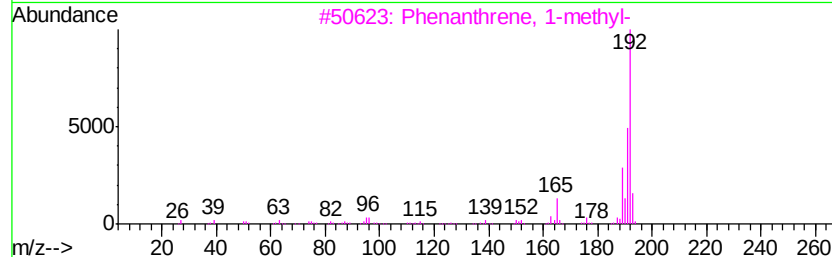
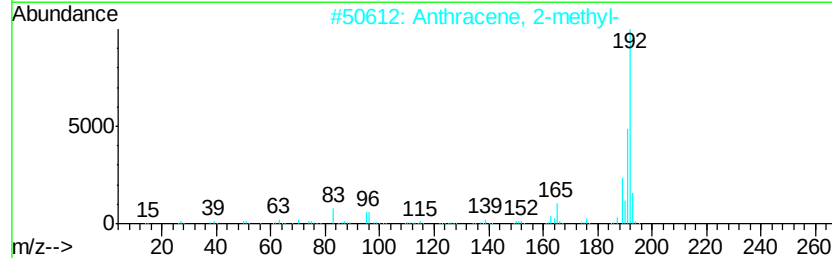
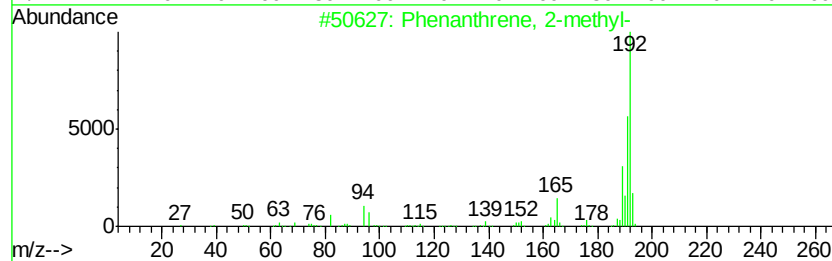
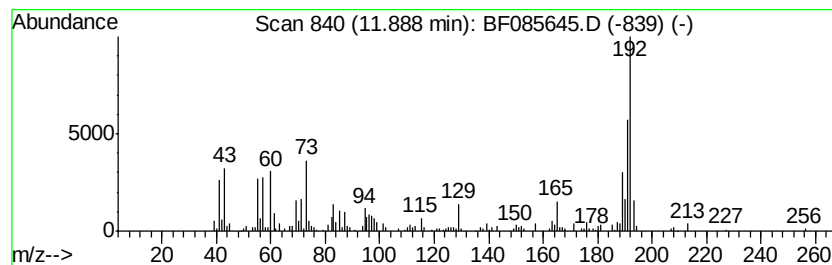
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ClientSampleId :  
PJ-SB-11-1.5-2.0

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.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 Phenanthrene, 2-methyl- Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.89     | 7.03 ng                             | 267543 | Phenanthrene-d10 | 11.36       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 98   |
| 2         | Anthracene, 2-methyl-               | 192    | C15H12           | 000613-12-7 | 97   |
| 3         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 96   |
| 4         | 1H-Cyclopropa[l]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 96   |
| 5         | 1H-Indene, 2-phenyl-                | 192    | C15H12           | 004505-48-0 | 93   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PJ-SB-11-1.5-2.0

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

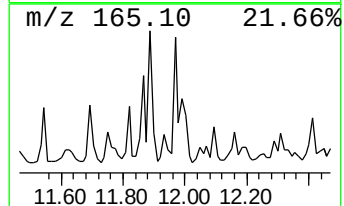
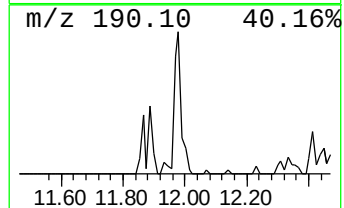
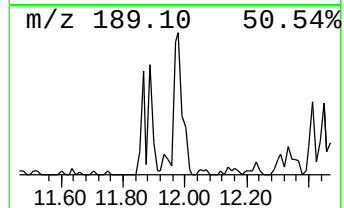
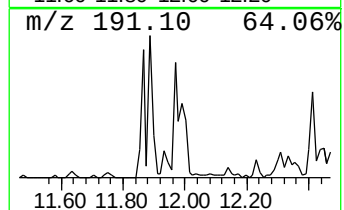
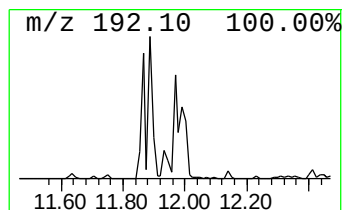
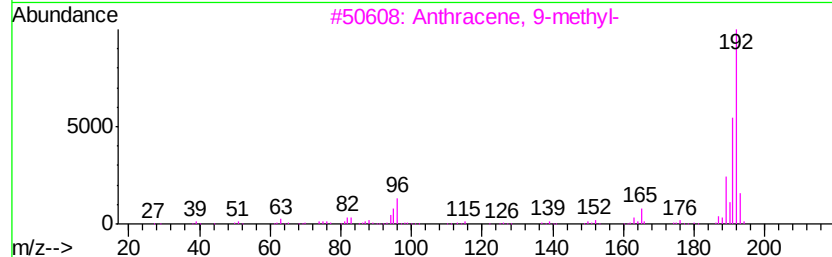
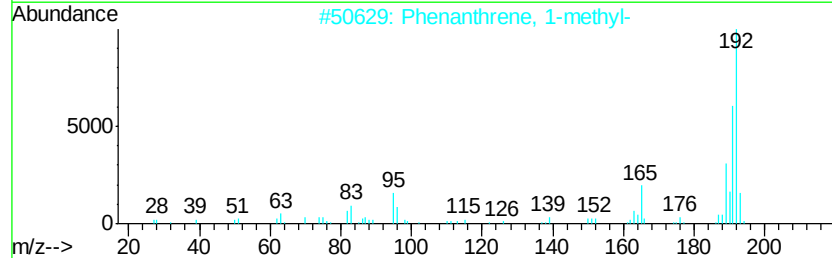
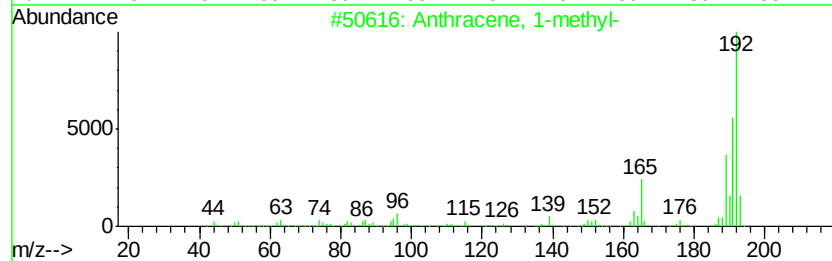
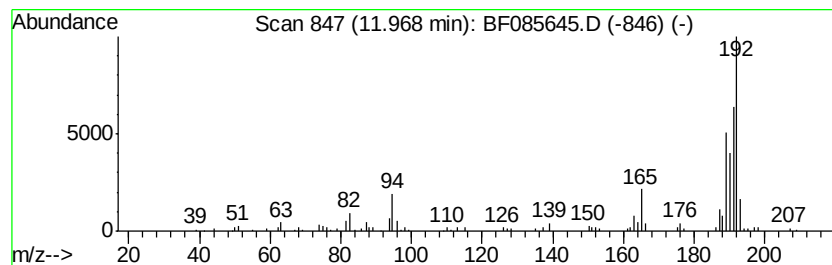
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Anthracene, 1-methyl- Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.97 | 5.45 ng | 207709 | Phenanthrene-d10 | 11.36 |

| Hit# of | 5 | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|---------|---|---------------------------------------|-----|---------|-------------|------|
| 1       |   | Anthracene, 1-methyl-                 | 192 | C15H12  | 000610-48-0 | 90   |
| 2       |   | Phenanthrene, 1-methyl-               | 192 | C15H12  | 000832-69-9 | 78   |
| 3       |   | Anthracene, 9-methyl-                 | 192 | C15H12  | 000779-02-2 | 78   |
| 4       |   | Phenanthrene, 2-methyl-               | 192 | C15H12  | 002531-84-2 | 60   |
| 5       |   | 1H-Cyclopropa[l]phenanthrene, 1a, ... | 192 | C15H12  | 000949-41-7 | 60   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PJ-SB-11-1.5-2.0

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

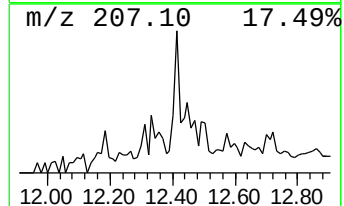
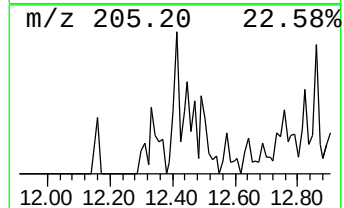
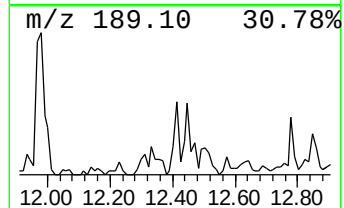
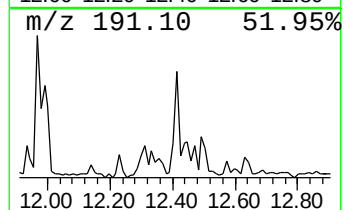
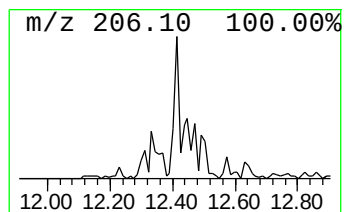
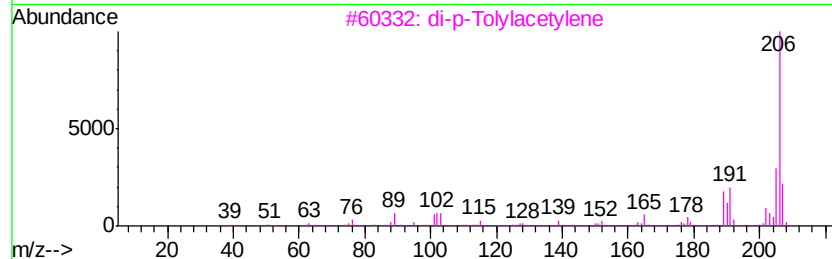
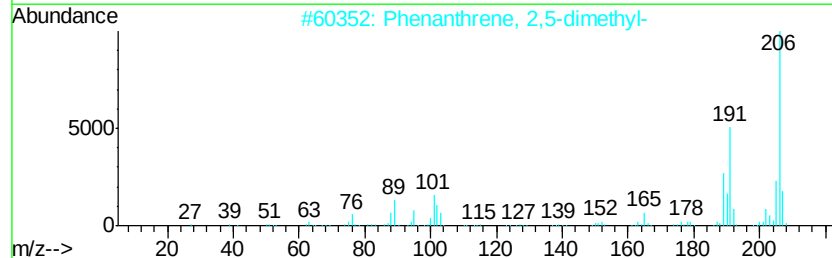
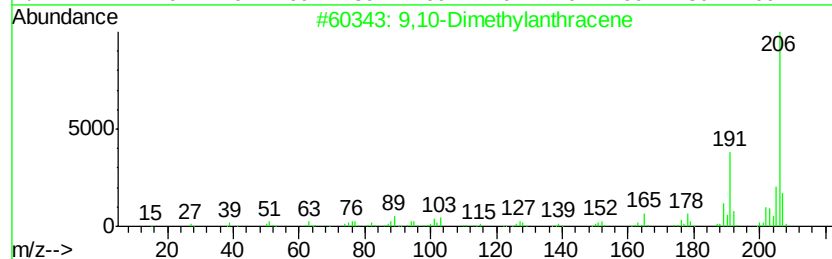
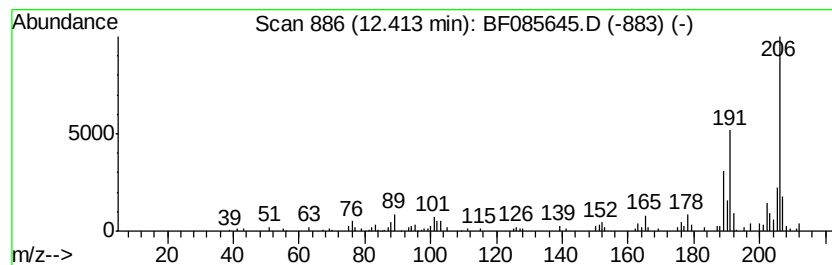
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 9,10-Dimethylantracene Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.41 | 3.13 ng | 119045 | Phenanthrene-d10 | 11.36 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | 9,10-Dimethylantracene        | 206 | C16H14  | 000781-43-1 | 97   |
| 2    |    |   | Phenanthrene, 2,5-dimethyl-   | 206 | C16H14  | 003674-66-6 | 95   |
| 3    |    |   | di-p-Tolylacetylene           | 206 | C16H14  | 002789-88-0 | 93   |
| 4    |    |   | Phenanthrene, 2,3-dimethyl-   | 206 | C16H14  | 003674-65-5 | 93   |
| 5    |    |   | 1H-Indene, 2-methyl-3-phenyl- | 206 | C16H14  | 035099-60-6 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PJ-SB-11-1.5-2.0

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

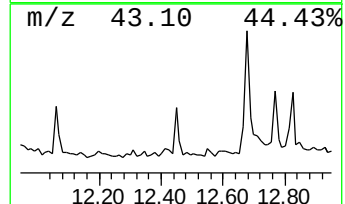
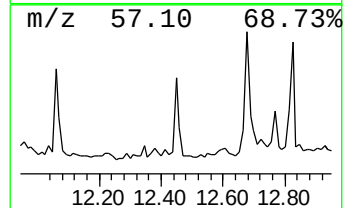
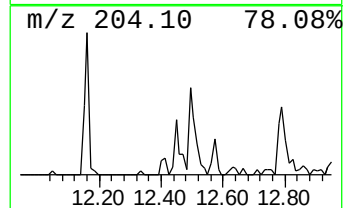
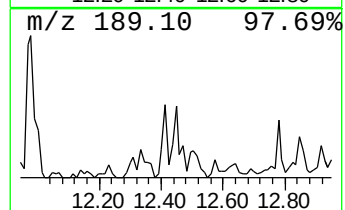
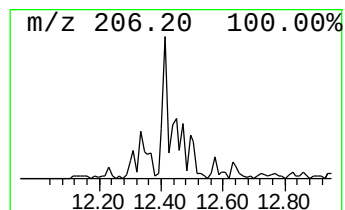
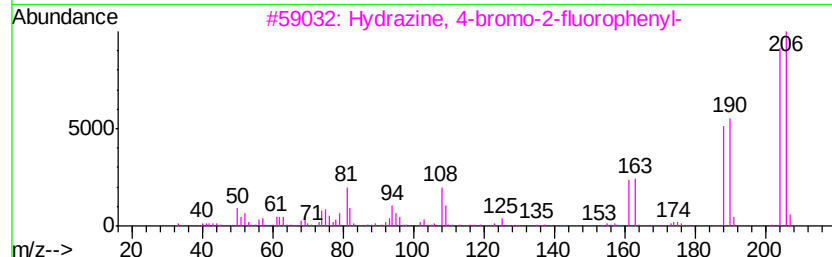
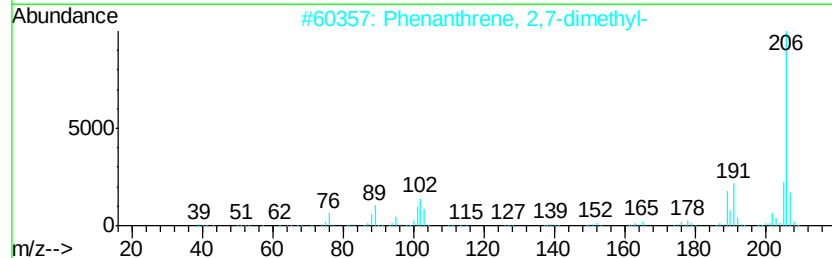
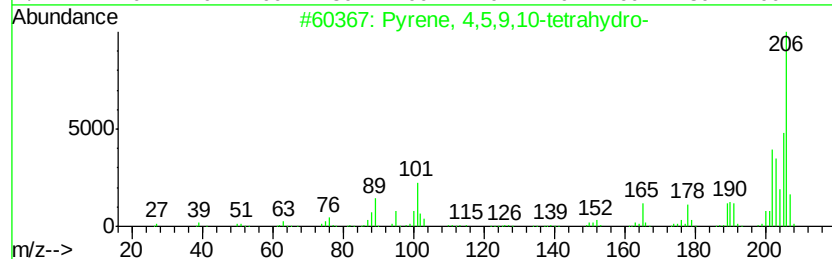
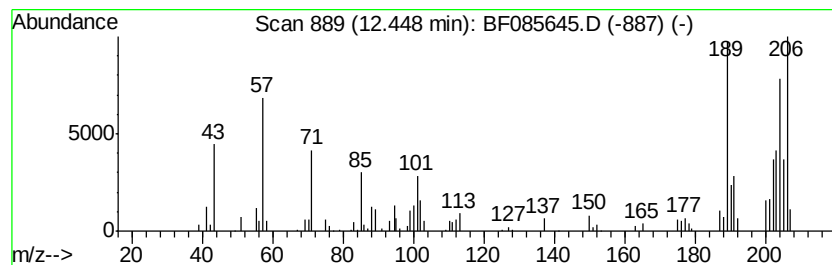
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 unknown12.45 Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.45 | 2.80 ng | 106557 | Phenanthrene-d10 | 11.36 |

| Hit# | of | Tentative ID                       | MW  | MolForm   | CAS#        | Qual |
|------|----|------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Pyrene, 4,5,9,10-tetrahydro-       | 206 | C16H14    | 000781-17-9 | 38   |
| 2    |    | Phenanthrene, 2,7-dimethyl-        | 206 | C16H14    | 001576-69-8 | 38   |
| 3    |    | Hydrazine, 4-bromo-2-fluorophenyl- | 204 | C6H6BrFN2 | 299440-17-8 | 35   |
| 4    |    | 4-(Trifluoromethoxy)benzoic acid   | 206 | C8H5F3O3  | 000330-12-1 | 30   |
| 5    |    | Phenanthrene, 3,6-dimethyl-        | 206 | C16H14    | 001576-67-6 | 27   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PJ-SB-11-1.5-2.0

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

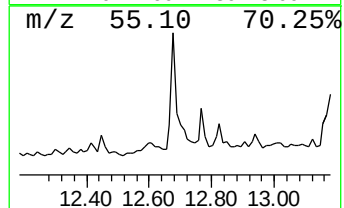
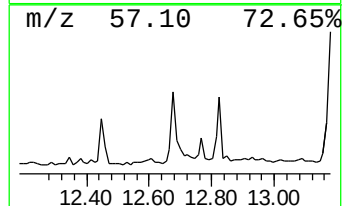
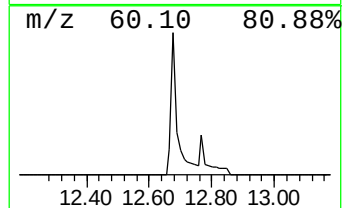
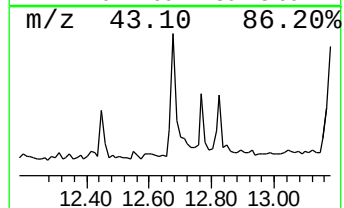
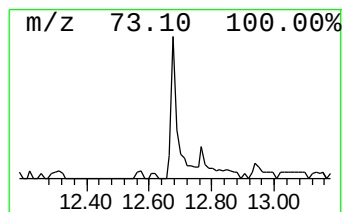
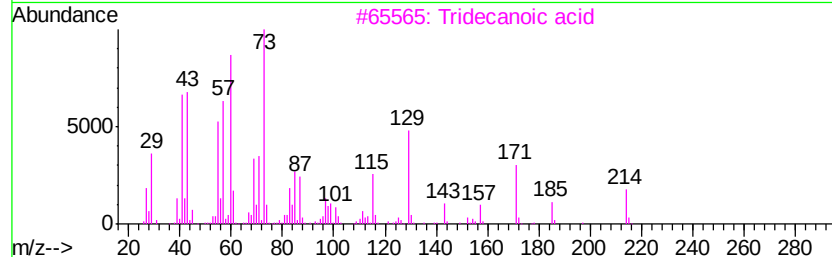
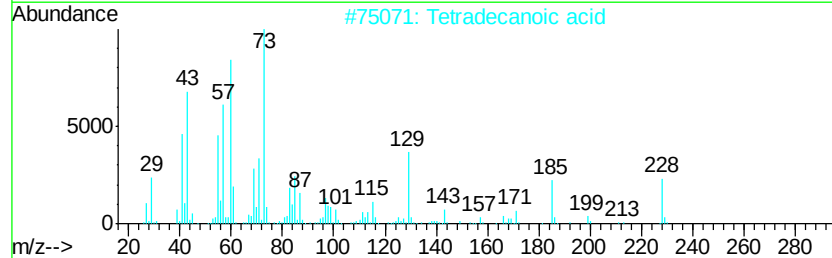
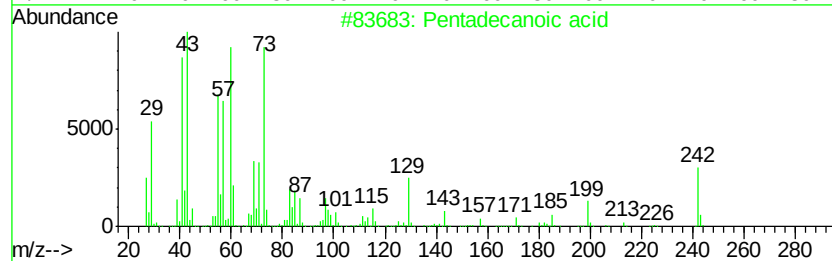
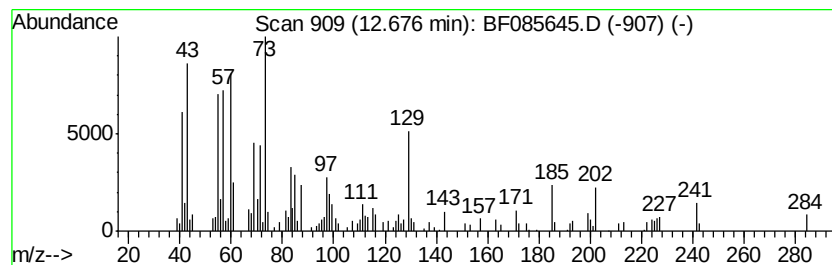
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Pentadecanoic acid Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.68 | 5.72 ng | 217929 | Phenanthrene-d10 | 11.36 |

| Hit# | of | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------|-----|----------|-------------|------|
| 1    | 5  | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 93   |
| 2    |    | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 74   |
| 3    |    | Tridecanoic acid   | 214 | C13H26O2 | 000638-53-9 | 72   |
| 4    |    | Dodecanoic acid    | 200 | C12H24O2 | 000143-07-7 | 70   |
| 5    |    | n-Decanoic acid    | 172 | C10H20O2 | 000334-48-5 | 60   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

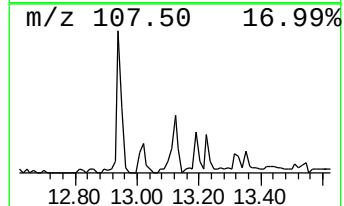
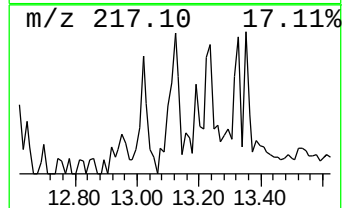
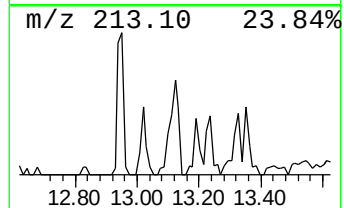
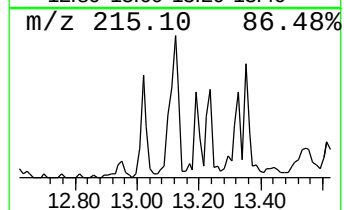
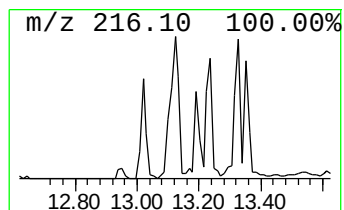
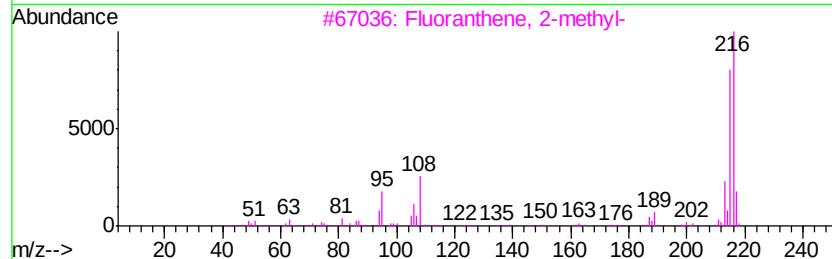
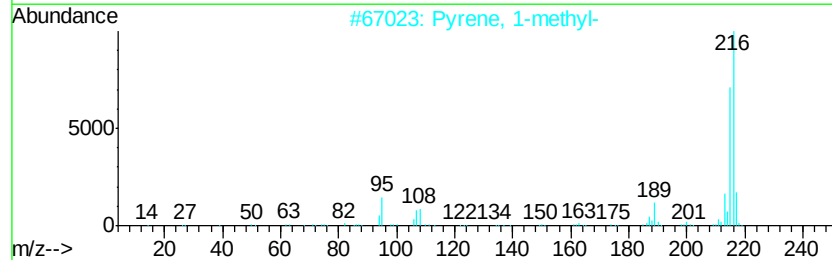
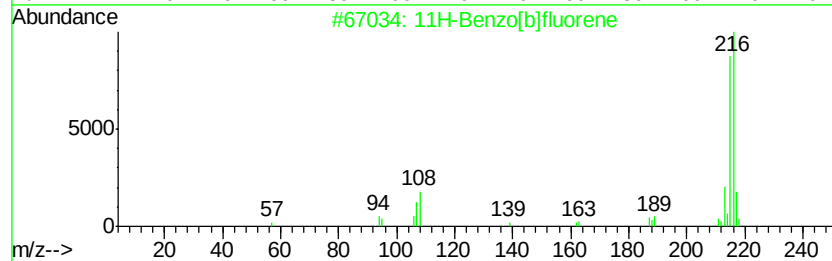
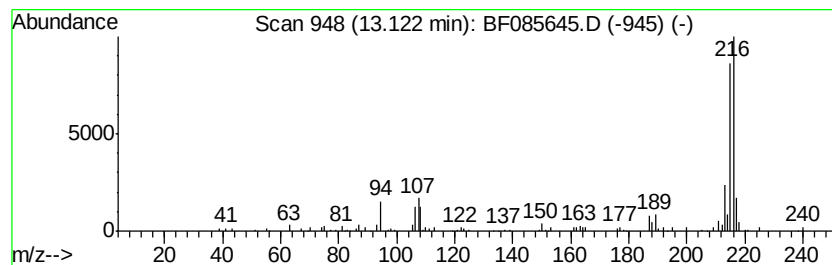
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 20

| R.T.      | EstConc                 | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------|--------|------------------|-------------|------|
| 13.12     | 2.54 ng                 | 146945 | Chrysene-d12     | 13.99       |      |
| Hit# of 5 | Tentative ID            | MW     | MolForm          | CAS#        | Qual |
| 1         | 11H-Benzo[b]fluorene    | 216    | C17H12           | 000243-17-4 | 94   |
| 2         | Pyrene, 1-methyl-       | 216    | C17H12           | 002381-21-7 | 94   |
| 3         | Fluoranthene, 2-methyl- | 216    | C17H12           | 033543-31-6 | 91   |
| 4         | 11H-Benzo[a]fluorene    | 216    | C17H12           | 000238-84-6 | 76   |
| 5         | Pyrene, 2-methyl-       | 216    | C17H12           | 003442-78-2 | 76   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

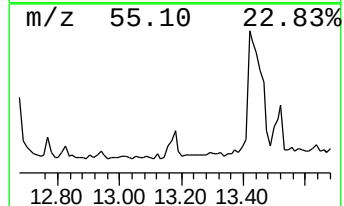
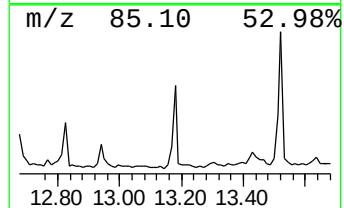
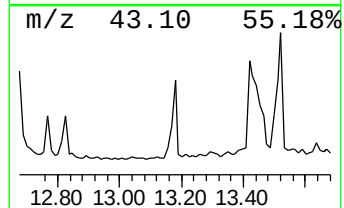
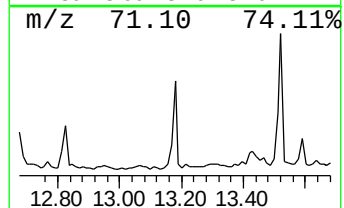
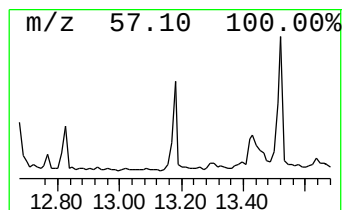
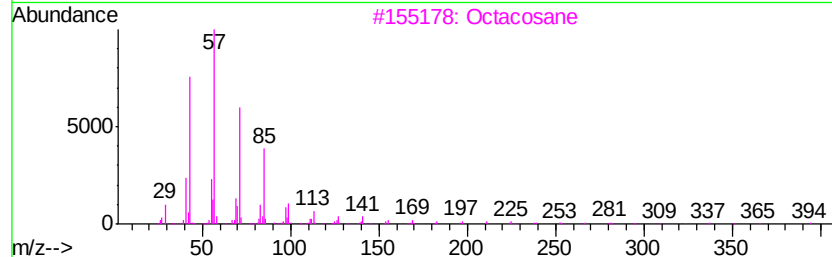
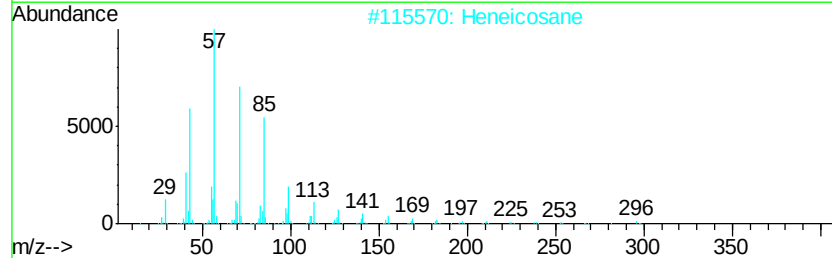
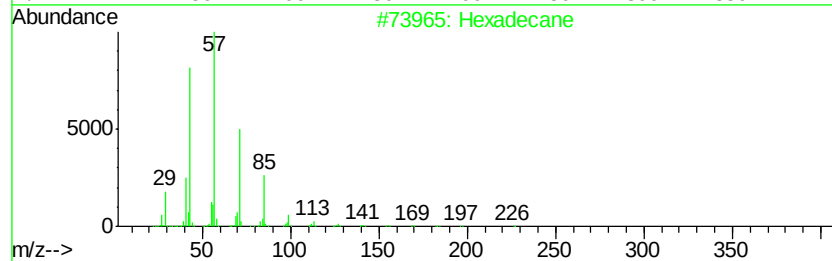
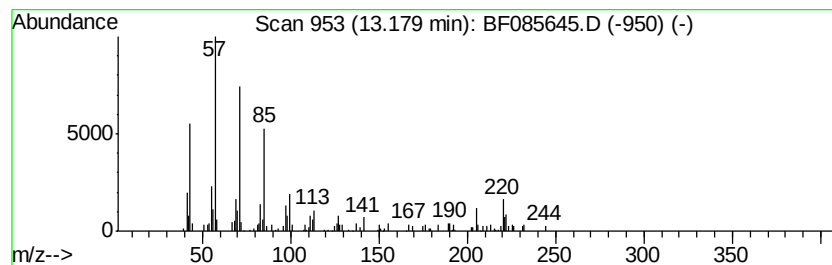
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 Hexadecane Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.18 | 2.77 ng | 160379 | Chrysene-d12     | 13.99 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane             | 226 | C16H34  | 000544-76-3 | 94   |
| 2    |    | Heneicosane            | 296 | C21H44  | 000629-94-7 | 87   |
| 3    |    | Octacosane             | 394 | C28H58  | 000630-02-4 | 87   |
| 4    |    | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 83   |
| 5    |    | Triacontane            | 422 | C30H62  | 000638-68-6 | 83   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
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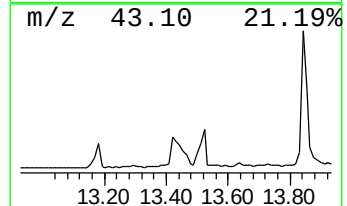
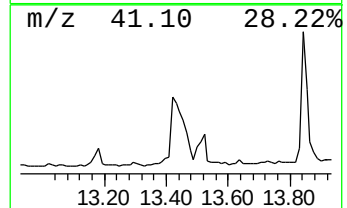
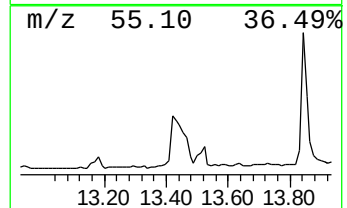
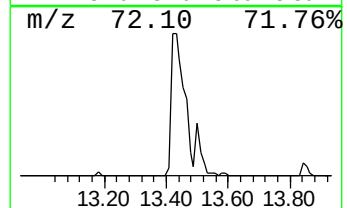
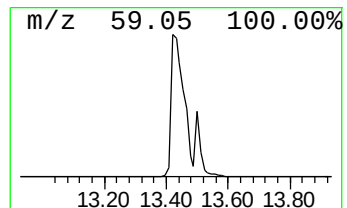
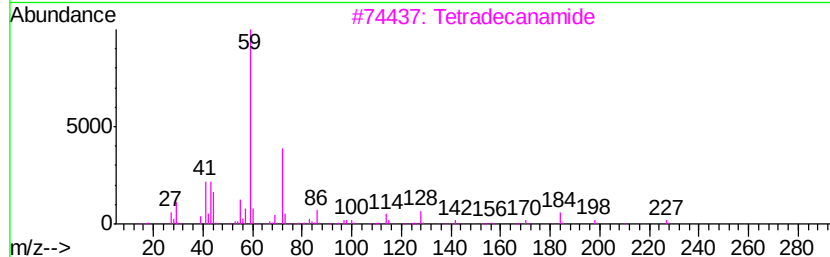
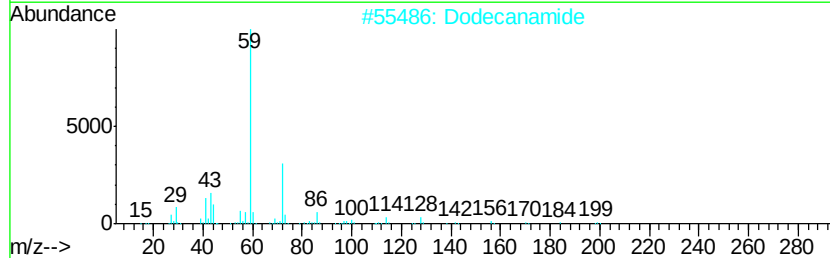
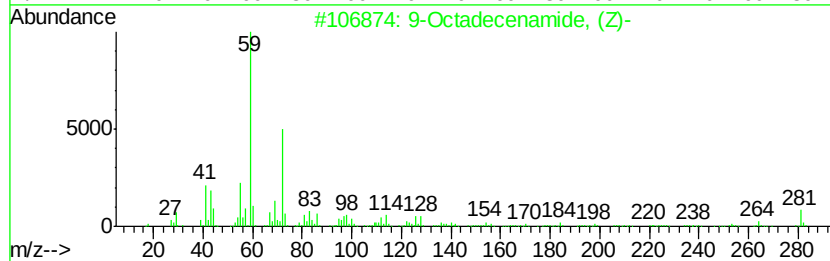
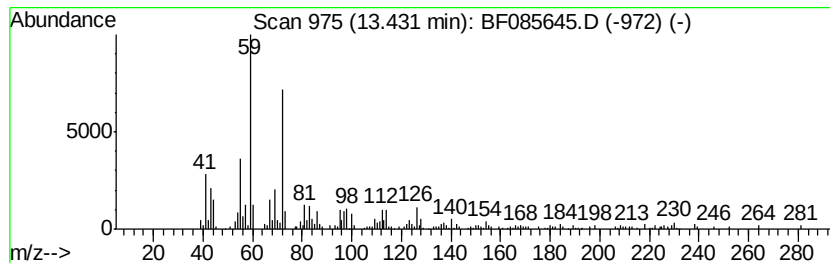
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 13 9-Octadecenamide, (Z)- Concentration Rank 5

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 13.43     | 8.67 ng                             | 501481 | Chrysene-d12     | 13.99        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 9-Octadecenamide, (Z)-              | 281    | C18H35NO         | 000301-02-0  | 91   |
| 2         | Dodecanamide                        | 199    | C12H25NO         | 001120-16-7  | 53   |
| 3         | Tetradecanamide                     | 227    | C14H29NO         | 000638-58-4  | 49   |
| 4         | Benzeneethanamine, 2-fluoro-.bet... | 229    | C11H16FN03       | 061338-98-5  | 47   |
| 5         | 2-Methyl-tridecane-2,12-diol        | 230    | C14H30O2         | 1000187-03-5 | 38   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PJ-SB-11-1.5-2.0

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

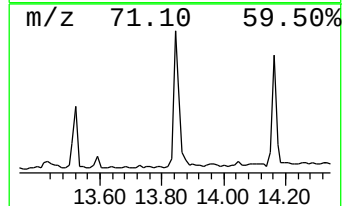
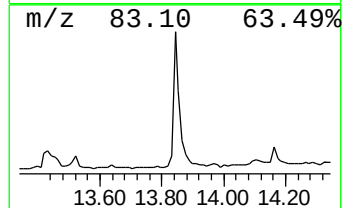
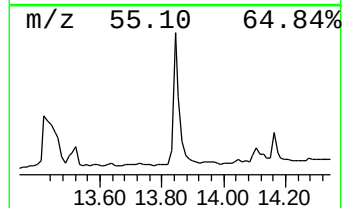
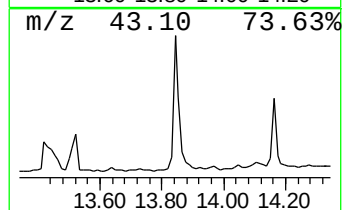
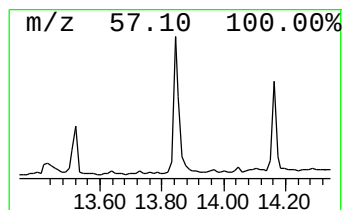
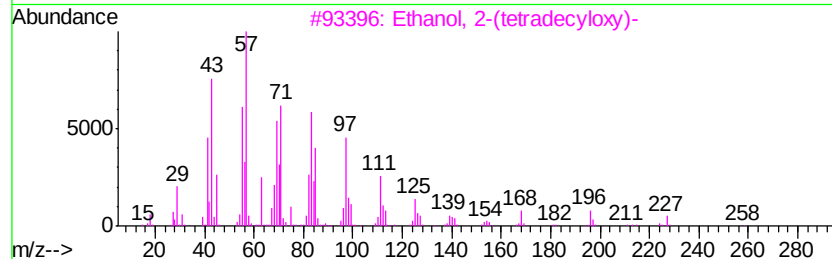
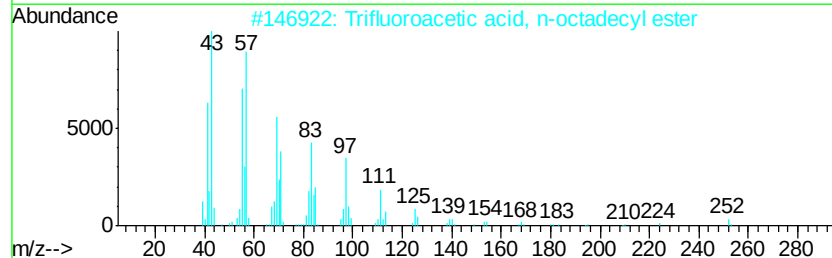
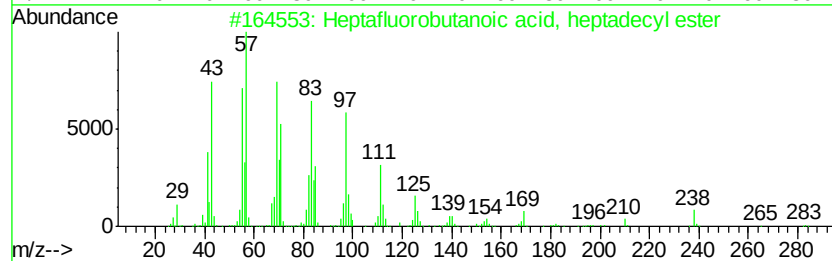
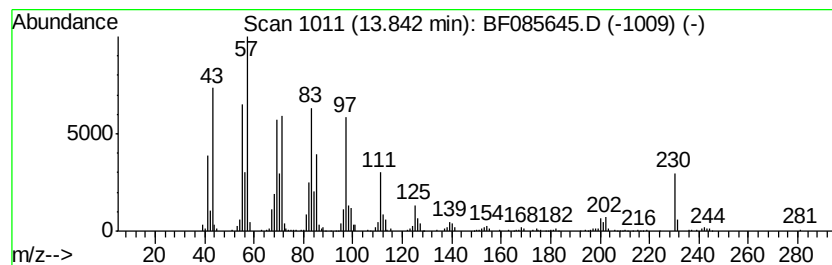
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Heptafluorobutanoic acid, h... Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.84 | 10.14 ng | 586809 | Chrysene-d12     | 13.99 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Heptafluorobutanoic acid, heptad... | 452 | C21H35F7O2  | 1000282-97-3 | 86   |
| 2    |    | Trifluoroacetic acid, n-octadecy... | 366 | C20H37F3O2  | 079392-43-1  | 81   |
| 3    |    | Ethanol, 2-(tetradecyloxy)-         | 258 | C16H34O2    | 002136-70-1  | 81   |
| 4    |    | Pentafluoropropionic acid, hepta... | 402 | C20H35F5O2  | 1000283-04-2 | 81   |
| 5    |    | 1-Hexadecanesulfonyl chloride       | 324 | C16H33ClO2S | 038775-38-1  | 74   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PJ-SB-11-1.5-2.0

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

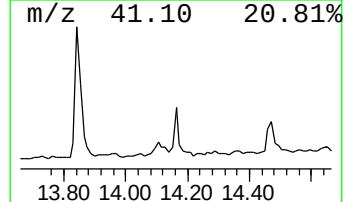
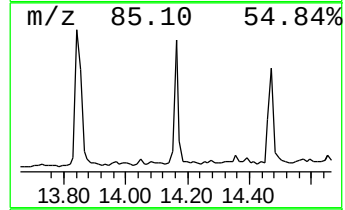
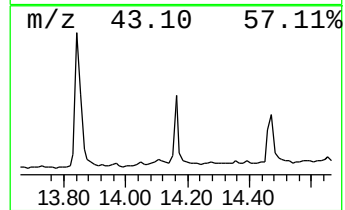
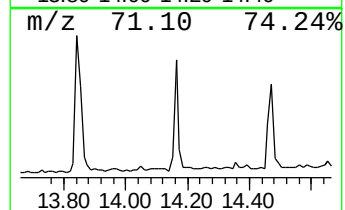
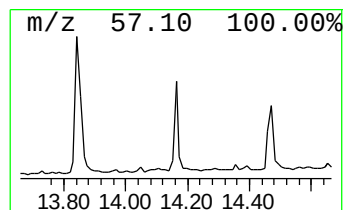
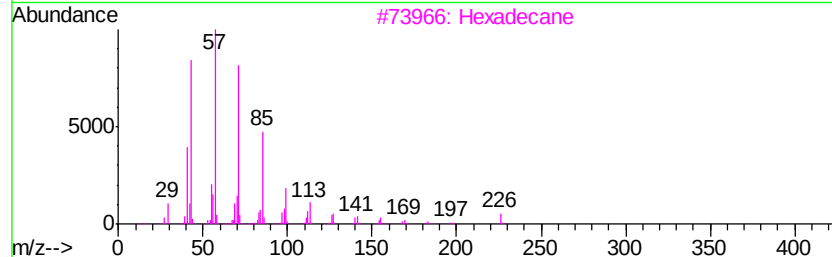
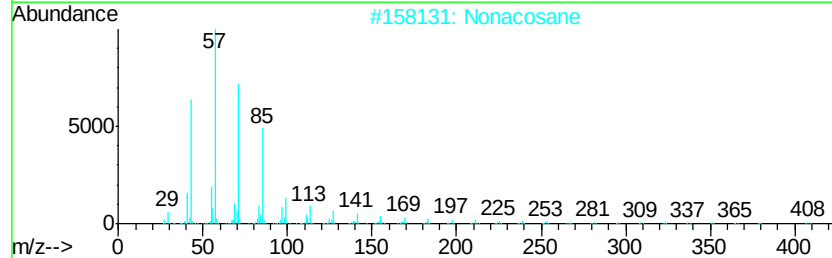
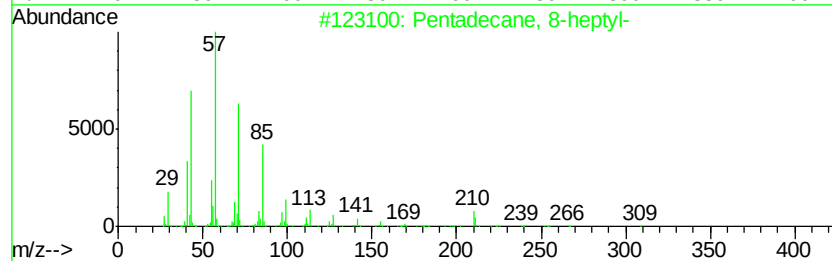
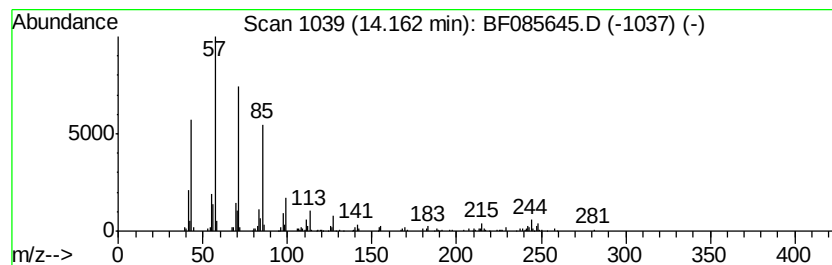
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 Pentadecane, 8-heptyl- Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.16 | 3.11 ng | 179691 | Chrysene-d12     | 13.99 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 87   |
| 2    |    |   | Nonacosane             | 408 | C29H60  | 000630-03-5 | 86   |
| 3    |    |   | Hexadecane             | 226 | C16H34  | 000544-76-3 | 86   |
| 4    |    |   | Octadecane, 1-iodo-    | 380 | C18H37I | 000629-93-6 | 80   |
| 5    |    |   | Heptacosane            | 380 | C27H56  | 000593-49-7 | 74   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
Acq On : 22 Mar 2016 4:21  
Operator : UM/SJ  
Sample : H1840-03  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PJ-SB-11-1.5-2.0

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

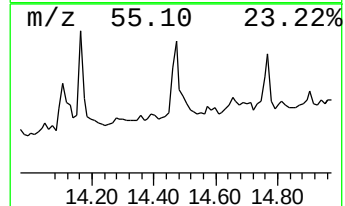
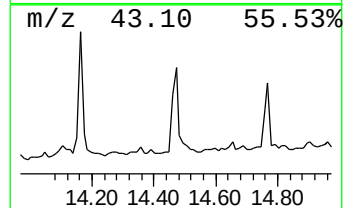
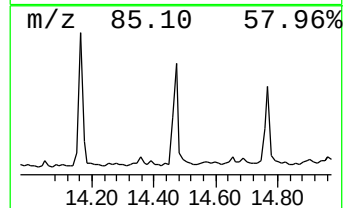
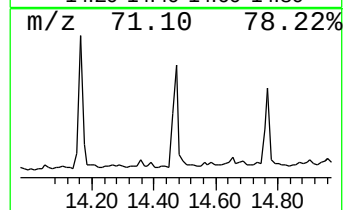
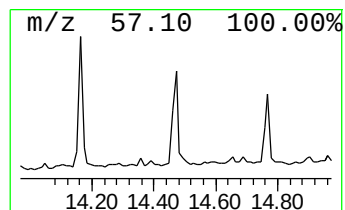
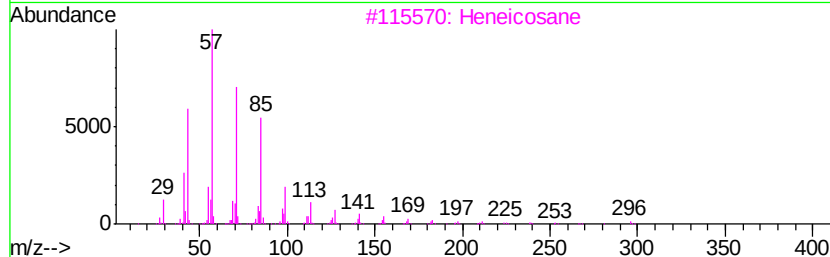
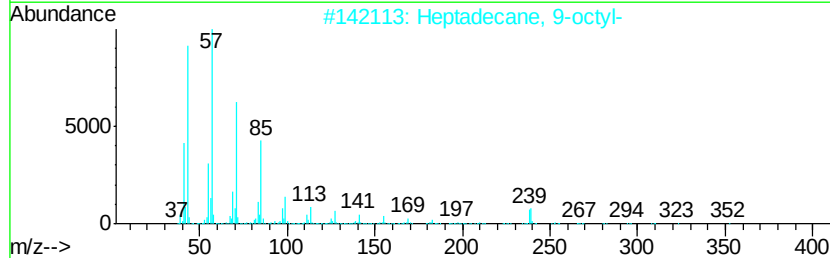
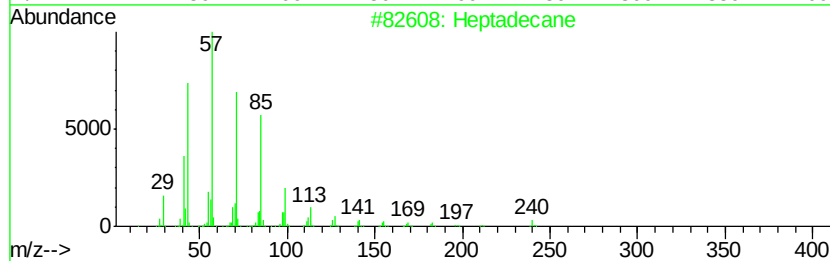
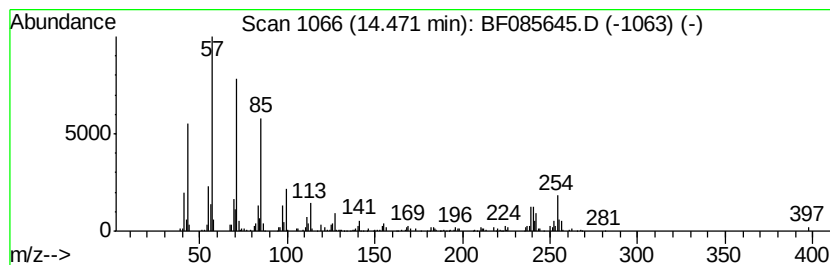
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Heptadecane Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.47 | 4.02 ng | 232839 | Chrysene-d12     | 13.99 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane           | 240 | C17H36  | 000629-78-7 | 90   |
| 2    |    |   | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 81   |
| 3    |    |   | Heneicosane           | 296 | C21H44  | 000629-94-7 | 76   |
| 4    |    |   | Nonadecane, 2-methyl- | 282 | C20H42  | 001560-86-7 | 74   |
| 5    |    |   | Eicosane              | 282 | C20H42  | 000112-95-8 | 70   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
Data File : BF085645.D  
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Operator : UM/SJ  
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ALS Vial : 28 Sample Multiplier: 1

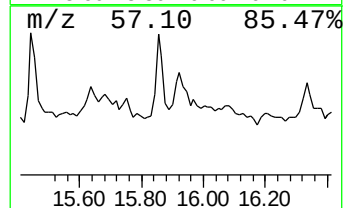
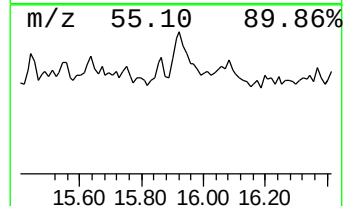
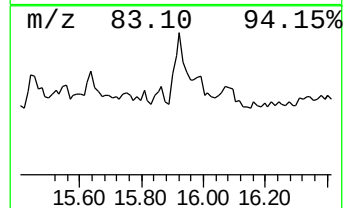
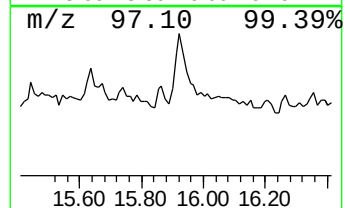
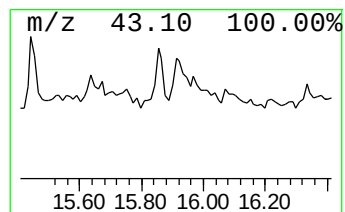
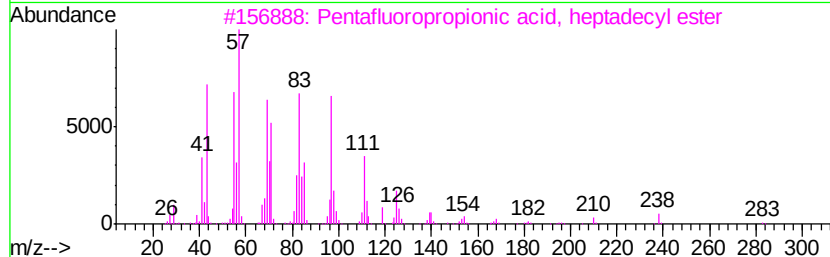
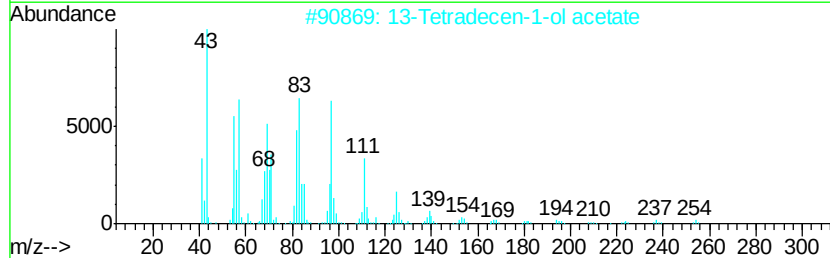
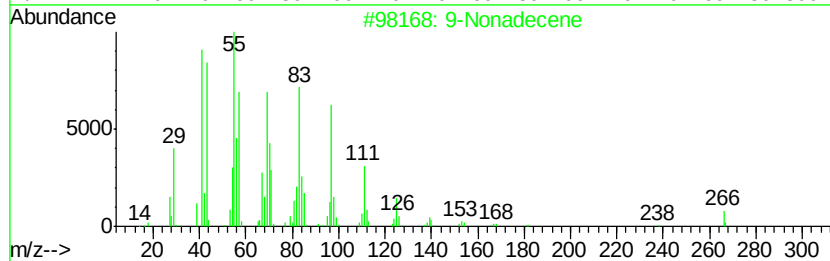
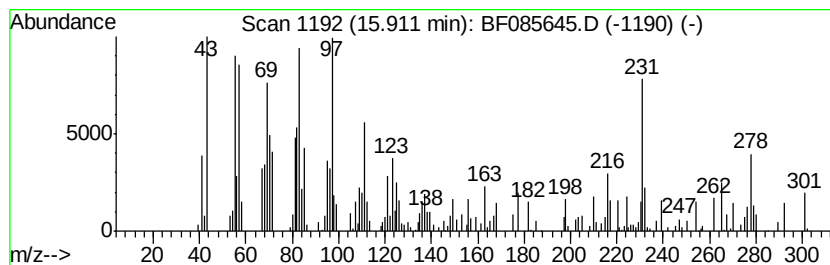
Instrument :  
BNA\_F  
ClientSampled :  
PJ-SB-11-1.5-2.0

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 20 9-Nonadecene Concentration Rank 13

| R.T.    | EstConc                             | Area           | Relative to ISTD | R.T.      |
|---------|-------------------------------------|----------------|------------------|-----------|
| 15.91   | 3.21 ng                             | 114891         | Perylene-d12     | 15.42     |
| Hit# of | 5                                   | Tentative ID   | MW MolForm       | CAS# Qual |
| 1       | 9-Nonadecene                        | 266 C19H38     | 031035-07-1      | 58        |
| 2       | 13-Tetradecen-1-ol acetate          | 254 C16H30O2   | 056221-91-1      | 50        |
| 3       | Pentafluoropropionic acid, hepta... | 402 C20H35F5O2 | 1000283-04-2     | 43        |
| 4       | Bromoacetic acid, pentadecyl ester  | 348 C17H33BrO2 | 131143-01-6      | 43        |
| 5       | 1-Eicosene                          | 280 C20H40     | 003452-07-1      | 43        |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032116\  
 Data File : BF085645.D  
 Acq On : 22 Mar 2016 4:21  
 Operator : UM/SJ  
 Sample : H1840-03  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PJ-SB-11-1.5-2.0

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.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 |
| Tetrachloroethylene   | 4.50  | 4.4 ng  |       | 150536   | 1                     | 6.84  | 684185  | 20.0 |
| 2-Pentanone, 4-hy...  | 5.03  | 9.4 ng  |       | 322826   | 1                     | 6.84  | 684185  | 20.0 |
| unknown6.61           | 6.61  | 76.0 ng |       | 2598740  | 1                     | 6.84  | 684185  | 20.0 |
| Dodecane              | 10.79 | 4.6 ng  |       | 174041   | 4                     | 11.36 | 761674  | 20.0 |
| Phenanthrene, 2-m...  | 11.89 | 7.0 ng  |       | 267543   | 4                     | 11.36 | 761674  | 20.0 |
| Anthracene, 1-met...  | 11.97 | 5.5 ng  |       | 207709   | 4                     | 11.36 | 761674  | 20.0 |
| 9,10-Dimethylanthe... | 12.41 | 3.1 ng  |       | 119045   | 4                     | 11.36 | 761674  | 20.0 |
| unknown12.45          | 12.45 | 2.8 ng  |       | 106557   | 4                     | 11.36 | 761674  | 20.0 |
| Pentadecanoic acid    | 12.68 | 5.7 ng  |       | 217929   | 4                     | 11.36 | 761674  | 20.0 |
| 11H-Benzo[b]fluorene  | 13.12 | 2.5 ng  |       | 146945   | 5                     | 13.99 | 1157070 | 20.0 |
| Hexadecane            | 13.18 | 2.8 ng  |       | 160379   | 5                     | 13.99 | 1157070 | 20.0 |
| 9-Octadecenamide, ... | 13.43 | 8.7 ng  |       | 501481   | 5                     | 13.99 | 1157070 | 20.0 |
| Heptafluorobutano...  | 13.84 | 10.1 ng |       | 586809   | 5                     | 13.99 | 1157070 | 20.0 |
| Pentadecane, 8-he...  | 14.16 | 3.1 ng  |       | 179691   | 5                     | 13.99 | 1157070 | 20.0 |
| Heptadecane           | 14.47 | 4.0 ng  |       | 232839   | 5                     | 13.99 | 1157070 | 20.0 |
| 9-Nonadecene          | 15.91 | 3.2 ng  |       | 114891   | 6                     | 15.42 | 716056  | 20.0 |