

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052516\  
 Data File : BF087371.D  
 Acq On : 25 May 2016 14:29  
 Operator : UM/SJ  
 Sample : H3218-01  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 E-3(6.5-7)

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-BF052316.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 1.644       | 4             | 5           | 57           | rVB      | 5022711        | 10083639      | 100.00%         | 26.668%       |
| 2         | 4.776       | 277           | 279         | 295          | rBV      | 43204          | 143868        | 1.43%           | 0.380%        |
| 3         | 5.404       | 332           | 334         | 345          | rBV      | 1055373        | 1994386       | 19.78%          | 5.274%        |
| 4         | 7.016       | 473           | 475         | 477          | rBV      | 1192067        | 1604786       | 15.91%          | 4.244%        |
| 5         | 7.050       | 477           | 478         | 486          | rVV      | 1574980        | 2575997       | 25.55%          | 6.813%        |
| 6         | 7.165       | 486           | 488         | 494          | rVV      | 1618262        | 2430439       | 24.10%          | 6.428%        |
| 7         | 7.313       | 498           | 501         | 506          | rVB      | 78851          | 114117        | 1.13%           | 0.302%        |
| 8         | 7.519       | 517           | 519         | 524          | rBV      | 506910         | 707338        | 7.01%           | 1.871%        |
| 9         | 7.748       | 537           | 539         | 555          | rVB      | 1025677        | 1762154       | 17.48%          | 4.660%        |
| 10        | 8.433       | 597           | 599         | 607          | rBV      | 1234347        | 1802841       | 17.88%          | 4.768%        |
| 11        | 8.548       | 607           | 609         | 615          | rVB      | 70987          | 103618        | 1.03%           | 0.274%        |
| 12        | 8.650       | 615           | 618         | 621          | rBV      | 108260         | 171876        | 1.70%           | 0.455%        |
| 13        | 9.176       | 657           | 664         | 668          | rBV4     | 40826          | 123925        | 1.23%           | 0.328%        |
| 14        | 9.553       | 694           | 697         | 699          | rBV      | 686410         | 927216        | 9.20%           | 2.452%        |
| 15        | 9.633       | 702           | 704         | 712          | rVB      | 111808         | 196491        | 1.95%           | 0.520%        |
| 16        | 10.616      | 787           | 790         | 794          | rVB      | 163444         | 204182        | 2.02%           | 0.540%        |
| 17        | 10.936      | 815           | 818         | 826          | rBV2     | 49026          | 137358        | 1.36%           | 0.363%        |
| 18        | 11.325      | 849           | 852         | 855          | rBV      | 2134530        | 2525663       | 25.05%          | 6.680%        |
| 19        | 11.542      | 868           | 871         | 875          | rVV      | 85037          | 133386        | 1.32%           | 0.353%        |
| 20        | 12.022      | 911           | 913         | 915          | rBV      | 293848         | 376827        | 3.74%           | 0.997%        |
| 21        | 12.377      | 941           | 944         | 950          | rBV      | 770669         | 1152888       | 11.43%          | 3.049%        |
| 22        | 13.222      | 1015          | 1018        | 1019         | rBV      | 82824          | 126127        | 1.25%           | 0.334%        |
| 23        | 13.668      | 1054          | 1057        | 1063         | rBV      | 878108         | 1462488       | 14.50%          | 3.868%        |
| 24        | 14.777      | 1151          | 1154        | 1156         | rBV      | 755869         | 978055        | 9.70%           | 2.587%        |
| 25        | 14.823      | 1156          | 1158        | 1162         | rVB      | 100141         | 179541        | 1.78%           | 0.475%        |
| 26        | 15.074      | 1178          | 1180        | 1187         | rVB3     | 67428          | 159550        | 1.58%           | 0.422%        |
| 27        | 15.577      | 1222          | 1224        | 1226         | rBV      | 118794         | 176389        | 1.75%           | 0.466%        |
| 28        | 15.611      | 1226          | 1227        | 1233         | rVV      | 163138         | 257481        | 2.55%           | 0.681%        |
| 29        | 15.725      | 1236          | 1237        | 1239         | rVV2     | 73793          | 118128        | 1.17%           | 0.312%        |
| 30        | 15.760      | 1239          | 1240        | 1244         | rVB2     | 75722          | 116539        | 1.16%           | 0.308%        |
| 31        | 16.023      | 1258          | 1263        | 1268         | rVB5     | 74970          | 190670        | 1.89%           | 0.504%        |
| 32        | 16.274      | 1283          | 1285        | 1287         | rBV      | 116807         | 125341        | 1.24%           | 0.331%        |
| 33        | 16.377      | 1291          | 1294        | 1296         | rBV      | 128111         | 173873        | 1.72%           | 0.460%        |
| 34        | 16.960      | 1342          | 1345        | 1347         | rBV2     | 53647          | 103652        | 1.03%           | 0.274%        |

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ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

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-BF052316.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 17.120 | 1357 | 1359 | 1362 | rVV  | 1500223 | 1808462 | 17.93% | 4.783% |
| 36 | 17.223 | 1364 | 1368 | 1370 | rBV  | 56515   | 112511  | 1.12%  | 0.298% |
| 37 | 17.920 | 1424 | 1429 | 1431 | rBV2 | 95379   | 235834  | 2.34%  | 0.624% |
| 38 | 18.046 | 1437 | 1440 | 1443 | rVB  | 51491   | 102035  | 1.01%  | 0.270% |
| 39 | 18.411 | 1467 | 1472 | 1474 | rBV2 | 77799   | 212171  | 2.10%  | 0.561% |
| 40 | 18.469 | 1474 | 1477 | 1479 | rVV  | 686851  | 914347  | 9.07%  | 2.418% |
| 41 | 18.514 | 1479 | 1481 | 1484 | rVV2 | 100970  | 221173  | 2.19%  | 0.585% |
| 42 | 18.983 | 1516 | 1522 | 1524 | rBV  | 86786   | 176958  | 1.75%  | 0.468% |
| 43 | 20.137 | 1621 | 1623 | 1628 | rVV  | 434952  | 587550  | 5.83%  | 1.554% |

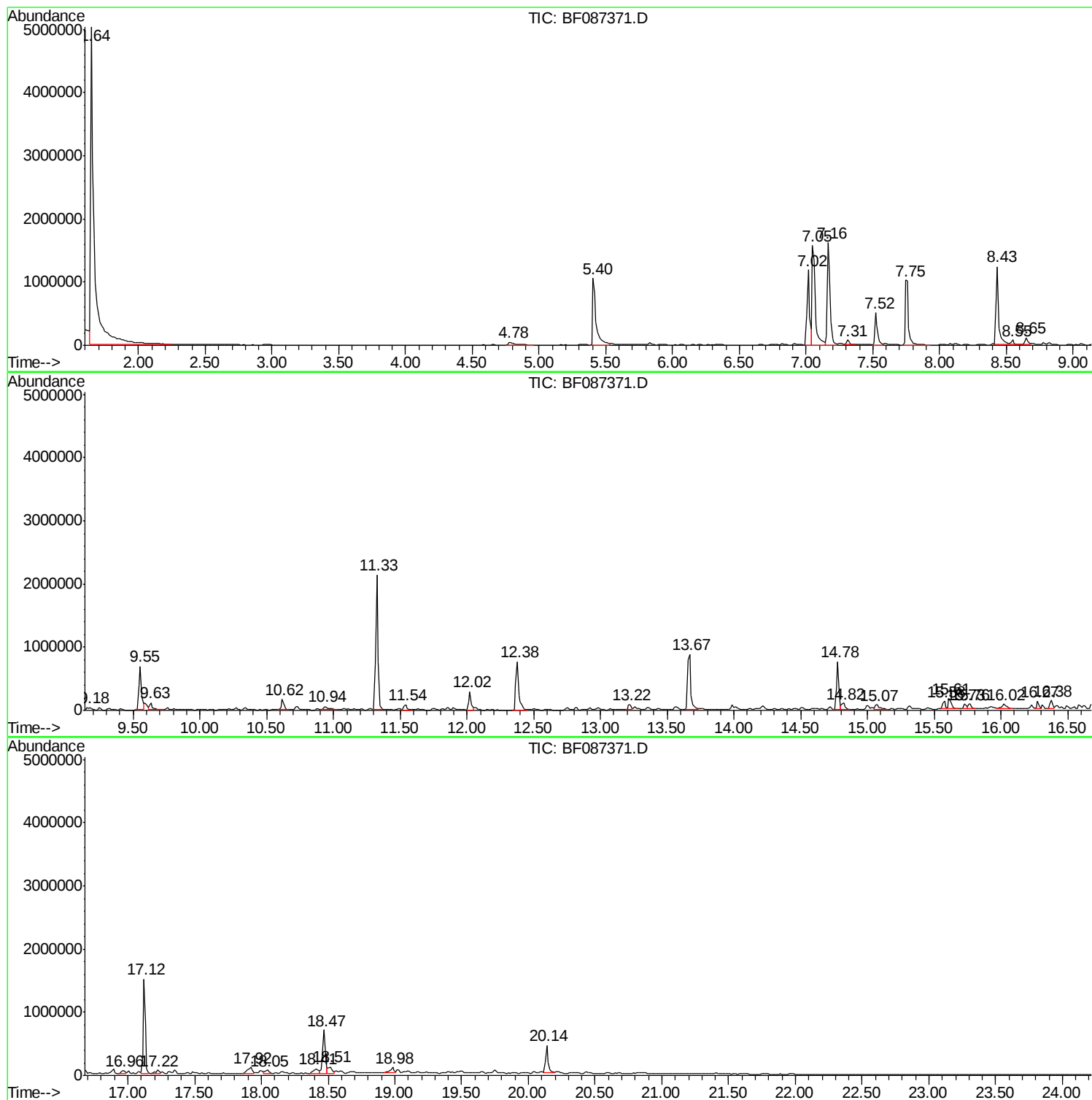
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Instrument :  
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ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

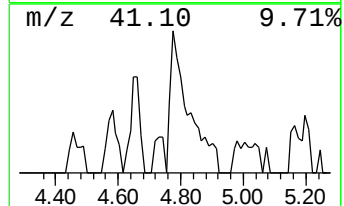
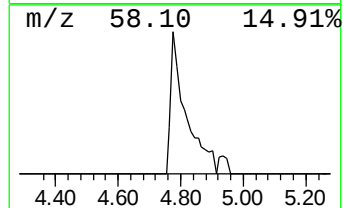
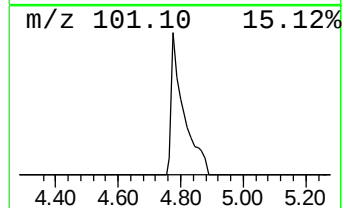
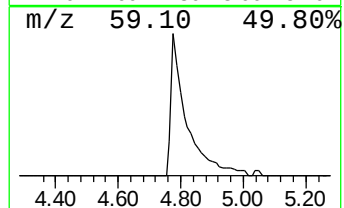
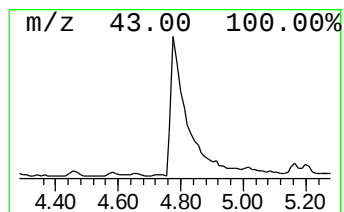
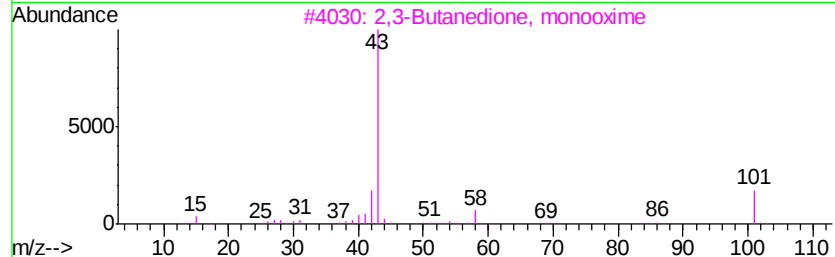
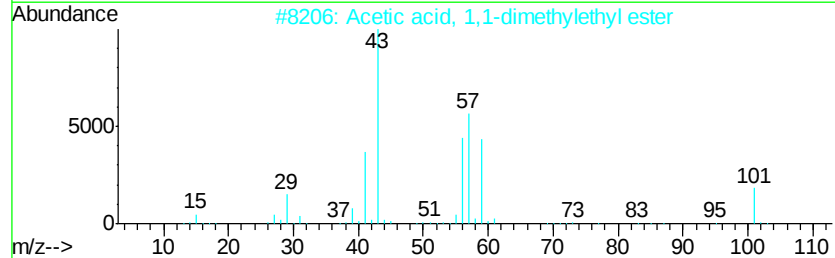
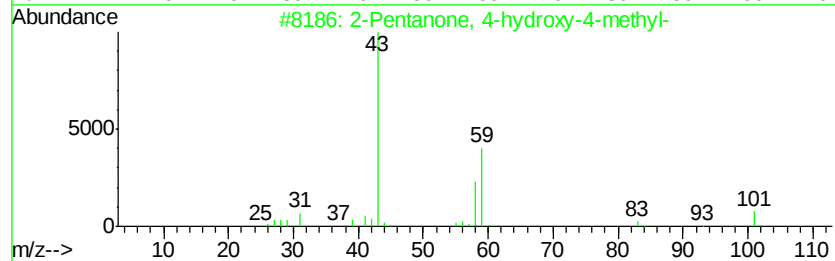
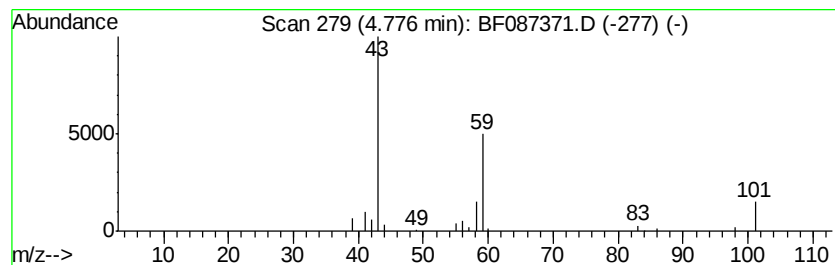
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.78 | 4.07 ng | 143868 | 1,4-Dichlorobenzene-d4 | 7.52 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2 | 000123-42-2 | 50   |
| 2    |    | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5 | 39   |
| 3    |    | 2,3-Butanedione, monooxime          | 101 | C4H7NO2 | 000057-71-6 | 9    |
| 4    |    | 2-Propyn-1-ol, acetate              | 98  | C5H6O2  | 000627-09-8 | 9    |
| 5    |    | 4-Penten-2-one, 4-methyl-           | 98  | C6H10O  | 003744-02-3 | 9    |



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Instrument :  
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E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

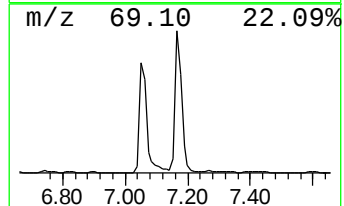
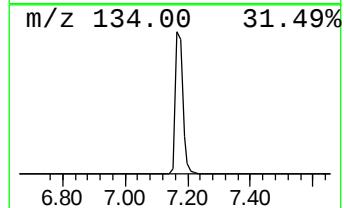
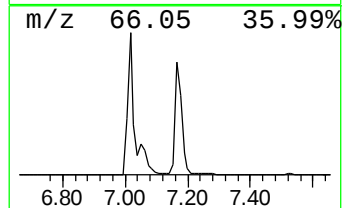
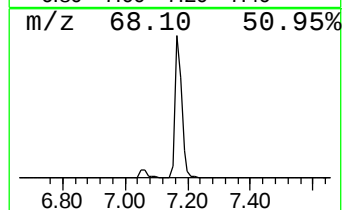
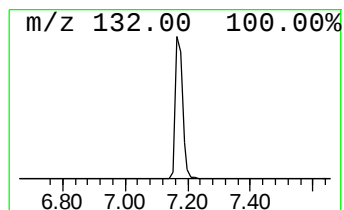
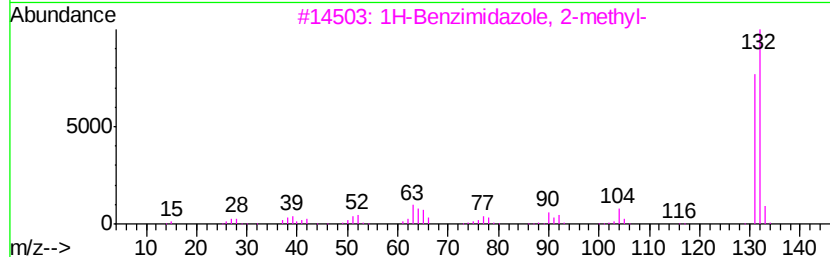
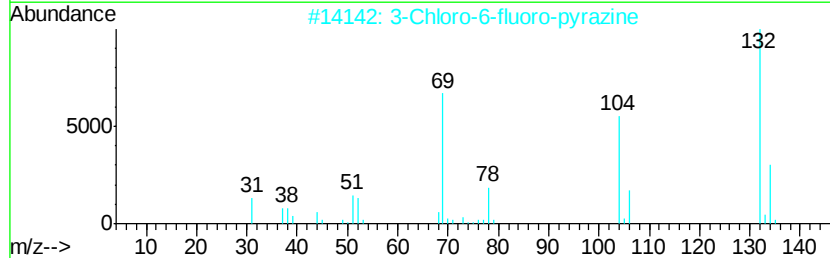
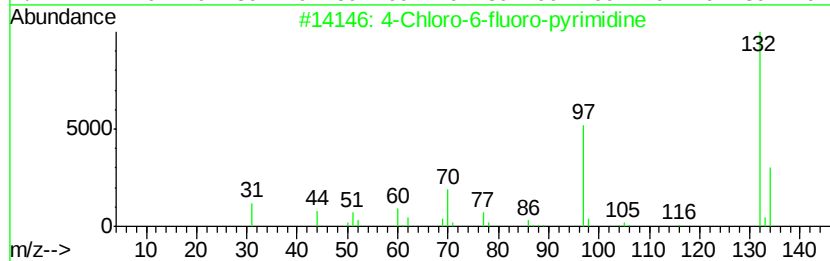
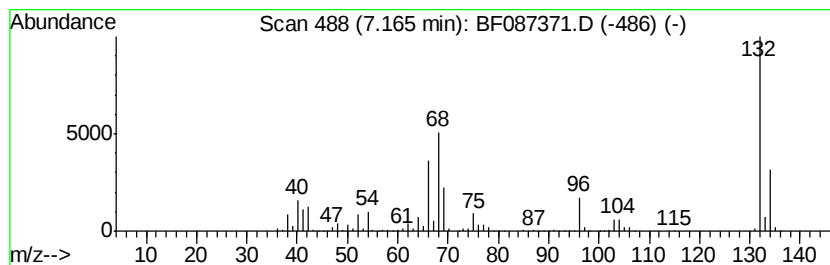
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown7.16 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.16 | 68.72 ng | 2430440 | 1,4-Dichlorobenzene-d4 | 7.52 |

| Hit# | of | Tentative ID                 | MW  | MolForm   | CAS#         | Qual |
|------|----|------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 4-Chloro-6-fluoro-pyrimidine | 132 | C4H2ClFN2 | 051422-01-6  | 27   |
| 2    |    | 3-Chloro-6-fluoro-pyrazine   | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 3    |    | 1H-Benzimidazole, 2-methyl-  | 132 | C8H8N2    | 000615-15-6  | 22   |
| 4    |    | Tranylcypromine-propionyl    | 189 | C12H15NO  | 1000123-86-3 | 22   |
| 5    |    | 1H-Indazole, 4-methyl-       | 132 | C8H8N2    | 1000316-00-9 | 12   |



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ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

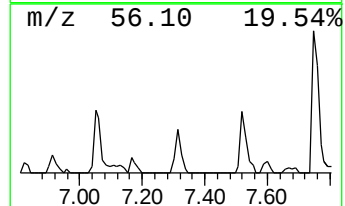
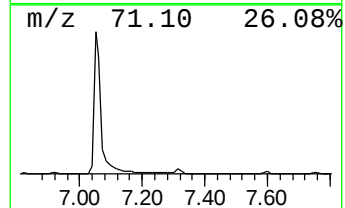
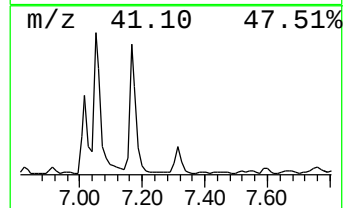
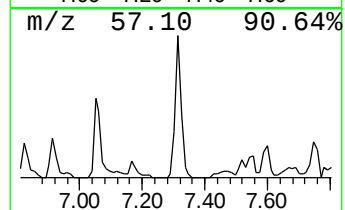
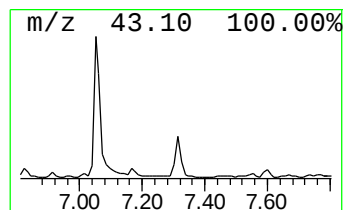
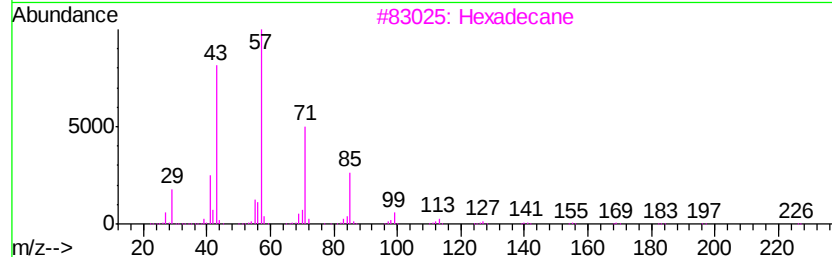
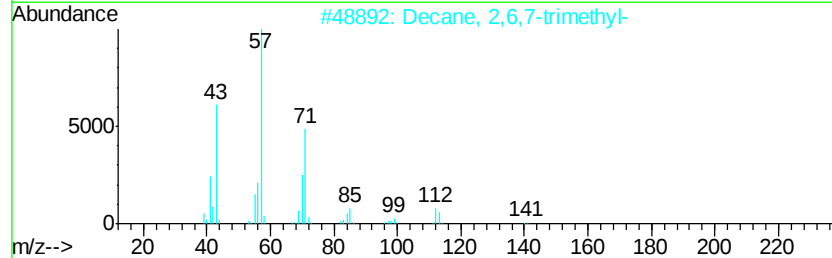
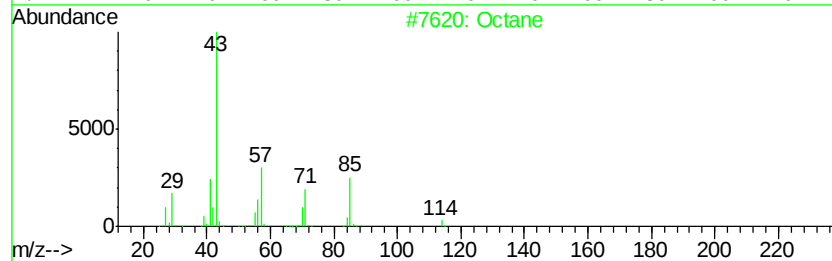
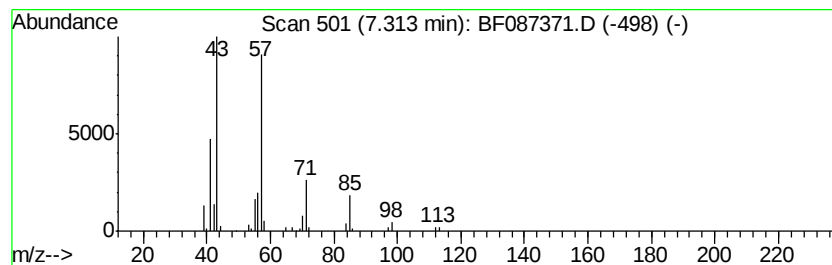
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 7.31 | 3.23 ng | 114117 | 1,4-Dichlorobenzene-d4 | 7.52 |

| Hit# | of | Tentative ID                      | MW  | MolForm  | CAS#         | Qual |
|------|----|-----------------------------------|-----|----------|--------------|------|
| 1    | 5  | Octane                            | 114 | C8H18    | 000111-65-9  | 64   |
| 2    |    | Decane, 2,6,7-trimethyl-          | 184 | C13H28   | 062108-25-2  | 64   |
| 3    |    | Hexadecane                        | 226 | C16H34   | 000544-76-3  | 59   |
| 4    |    | Decane                            | 142 | C10H22   | 000124-18-5  | 53   |
| 5    |    | Oxalic acid, isobutyl octyl ester | 258 | C14H26O4 | 1000309-37-3 | 53   |



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Instrument :  
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ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

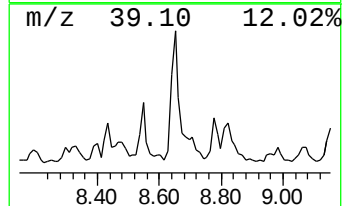
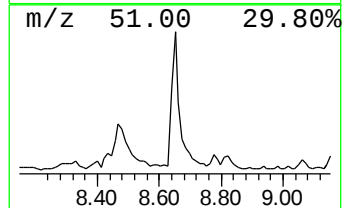
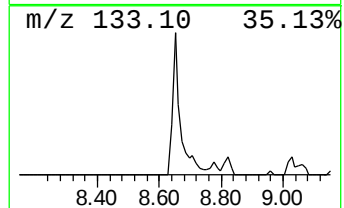
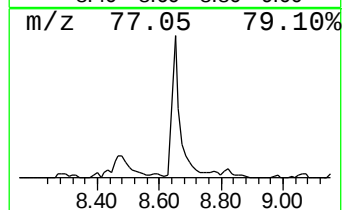
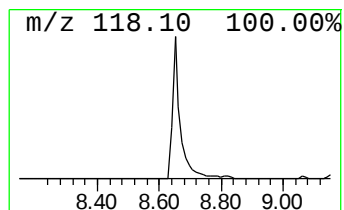
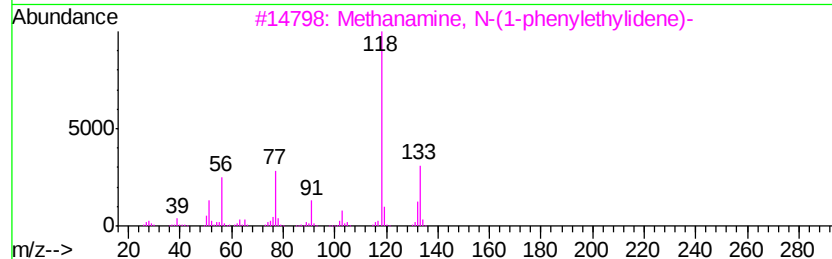
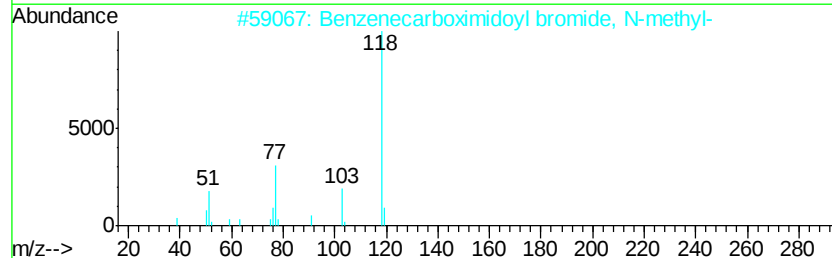
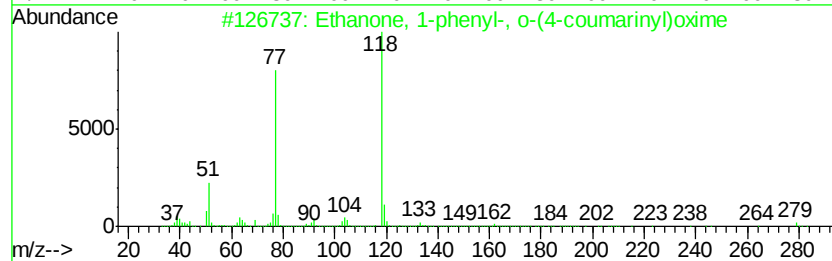
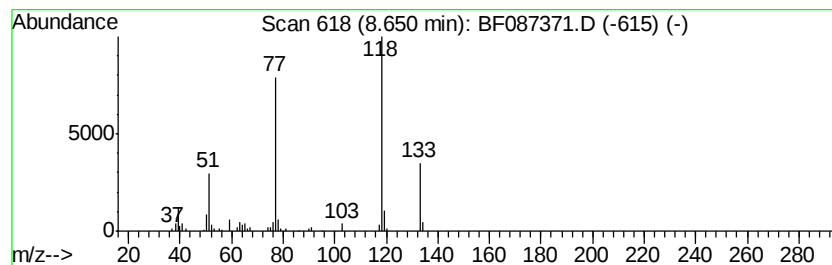
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Ethanone, 1-phenyl-, o-(4-c... Concentration Rank 11

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 8.65 | 3.71 ng | 171876 | Naphthalene-d8   | 9.55 |

| Hit# | of | Tentative ID                          | MW  | MolForm   | CAS#        | Qual |
|------|----|---------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Ethanone, 1-phenyl-, o-(4-coumar...   | 279 | C17H13NO3 | 034605-11-3 | 64   |
| 2    |    | Benzenecarboximidoyl bromide, N-...   | 197 | C8H8BrN   | 041182-85-8 | 40   |
| 3    |    | Methanamine, N-(1-phenylethylidene... | 133 | C9H11N    | 006907-71-7 | 40   |
| 4    |    | 2-Methylindoline                      | 133 | C9H11N    | 006872-06-6 | 38   |
| 5    |    | Cyanoic acid, phenyl ester            | 119 | C7H5NO    | 001122-85-6 | 30   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

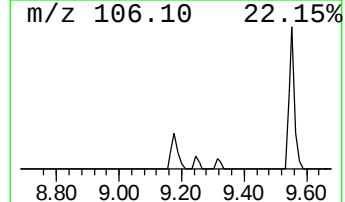
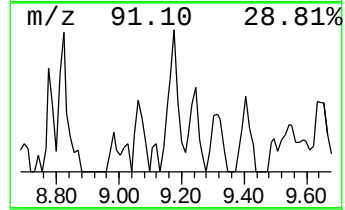
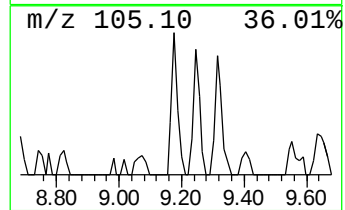
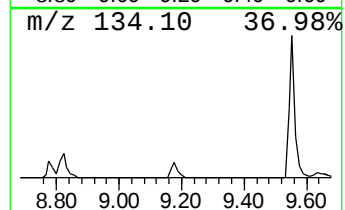
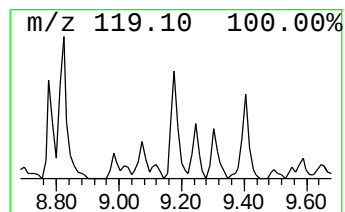
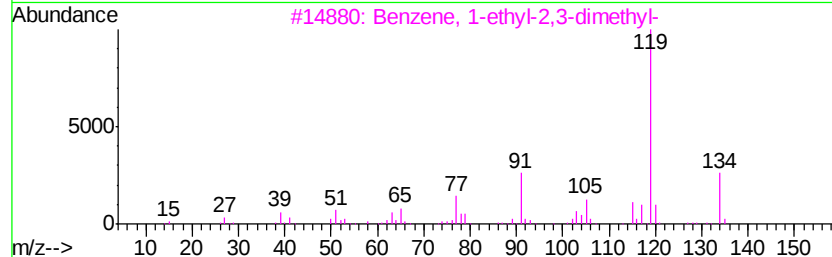
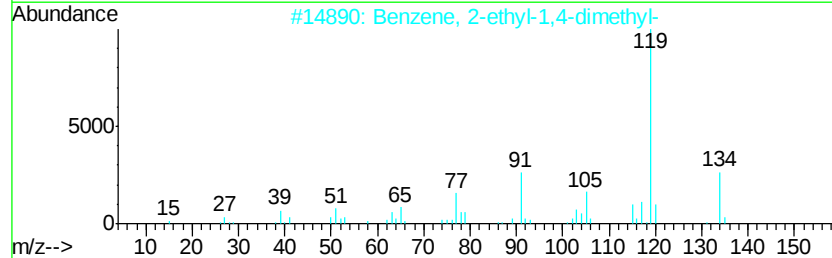
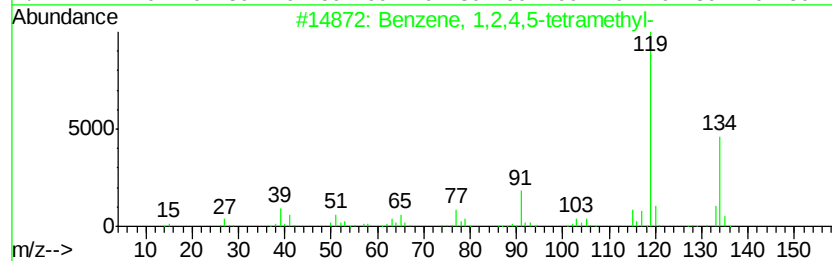
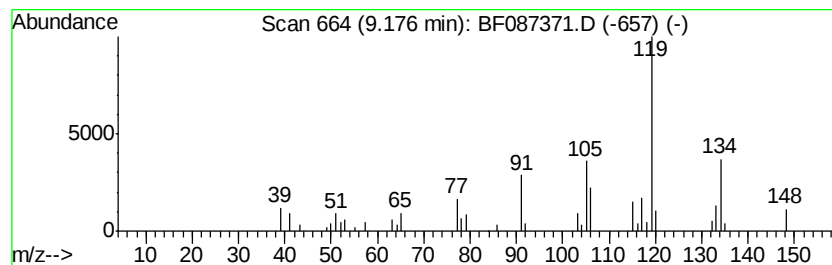
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 17

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.18 | 2.67 ng | 123925 | Naphthalene-d8   | 9.55 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 76   |
| 2    |    | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 76   |
| 3    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 76   |
| 4    |    | 1,4-Cyclohexadiene, 3-ethenyl-1,... | 134 | C10H14  | 062338-57-2 | 74   |
| 5    |    | p-Cymene                            | 134 | C10H14  | 000099-87-6 | 70   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

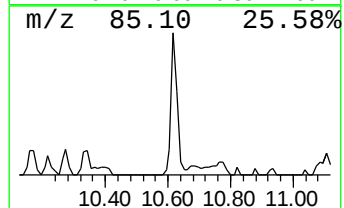
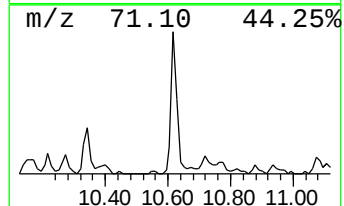
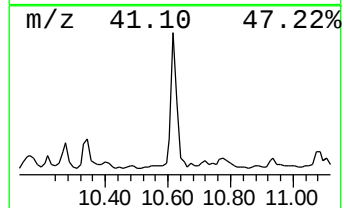
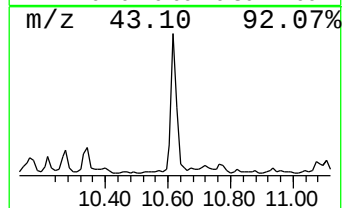
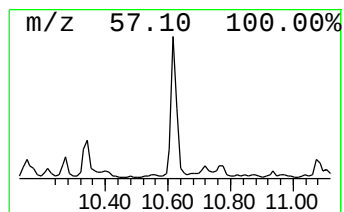
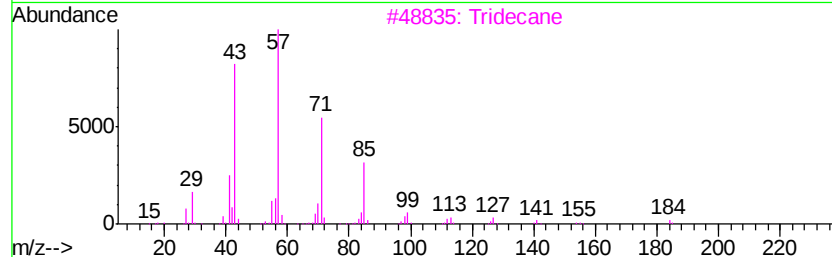
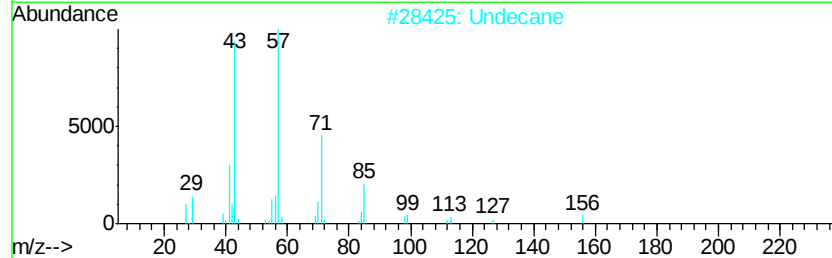
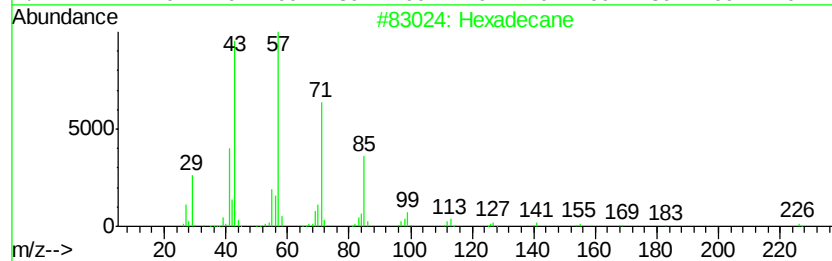
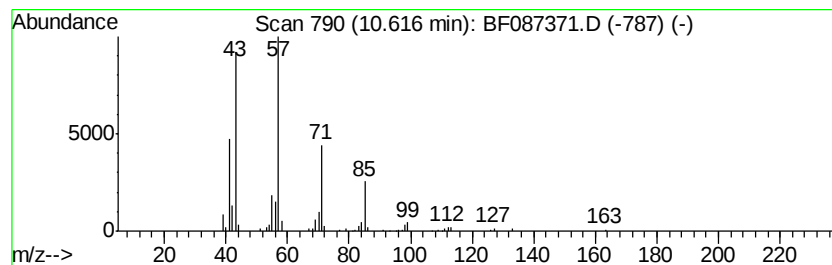
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Hexadecane Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T. |
|-------|---------|--------|------------------|------|
| 10.62 | 4.40 ng | 204182 | Naphthalene-d8   | 9.55 |

| Hit# of | 5                                    | Tentative ID | MW  | MolForm   | CAS#         | Qual |
|---------|--------------------------------------|--------------|-----|-----------|--------------|------|
| 1       | Hexadecane                           |              | 226 | C16H34    | 000544-76-3  | 80   |
| 2       | Undecane                             |              | 156 | C11H24    | 001120-21-4  | 78   |
| 3       | Tridecane                            |              | 184 | C13H28    | 000629-50-5  | 78   |
| 4       | Sulfurous acid, 2-propyl tetradec... |              | 320 | C17H36O3S | 1000309-12-5 | 78   |
| 5       | Tetradecane                          |              | 198 | C14H30    | 000629-59-4  | 72   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

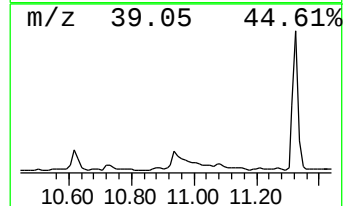
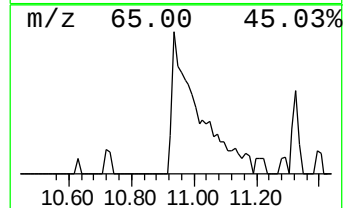
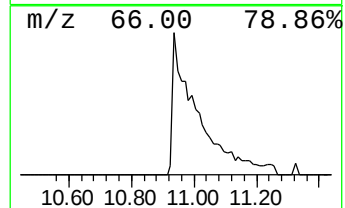
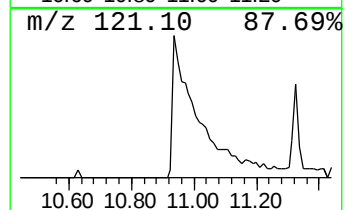
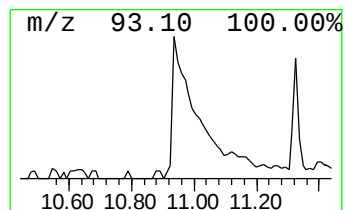
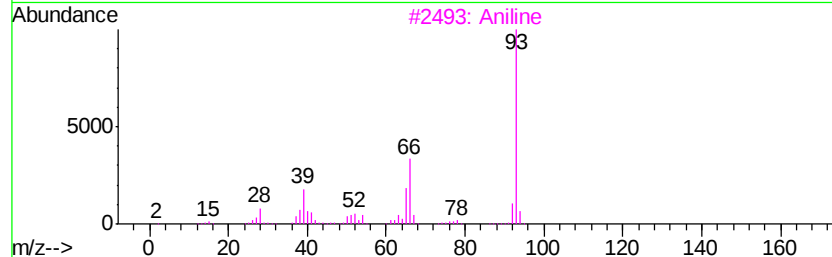
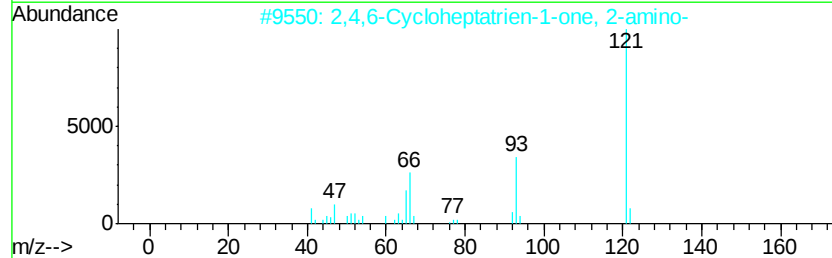
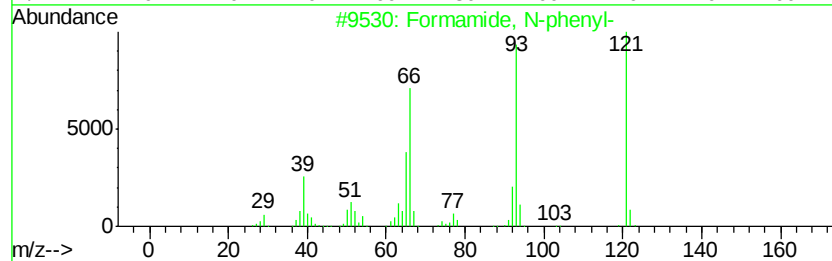
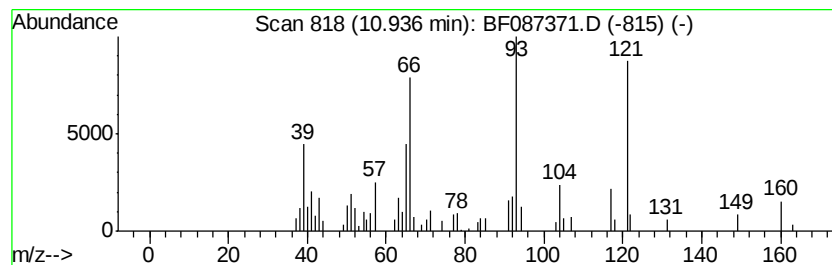
Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 Formamide, N-phenyl- Concentration Rank 16

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.           |
|---------|---------|-------------------------------------|------------------|----------------|
| 10.94   | 2.96 ng | 137358                              | Napthalene-d8    | 9.55           |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |         | Formamide, N-phenyl-                | 121 C7H7NO       | 000103-70-8 94 |
| 2       |         | 2,4,6-Cycloheptatrien-1-one, 2-a... | 121 C7H7NO       | 006264-93-3 93 |
| 3       |         | Aniline                             | 93 C6H7N         | 000062-53-3 81 |
| 4       |         | Pyridine, 4-methyl-                 | 93 C6H7N         | 000108-89-4 81 |
| 5       |         | Pyridine, 3-methyl-                 | 93 C6H7N         | 000108-99-6 81 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

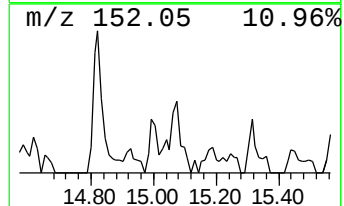
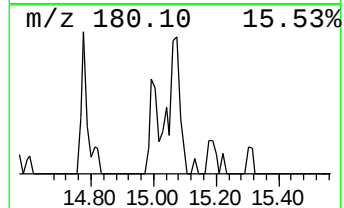
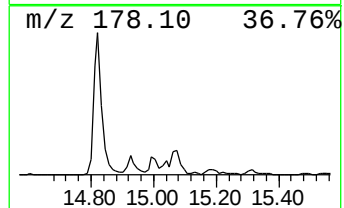
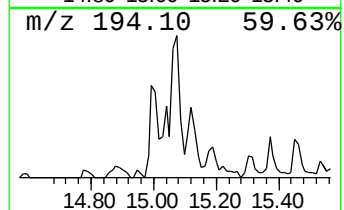
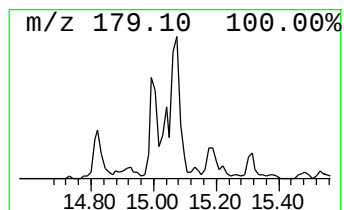
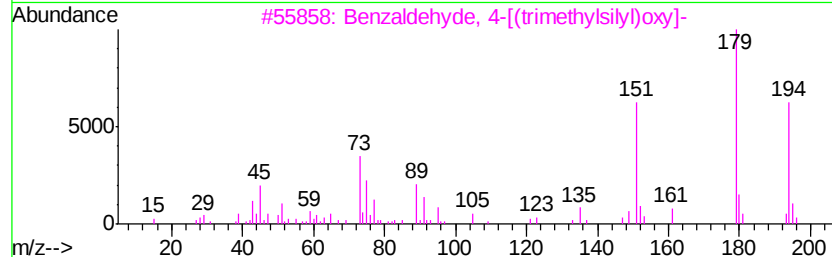
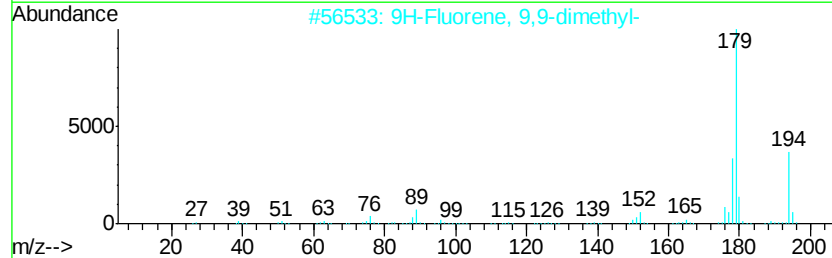
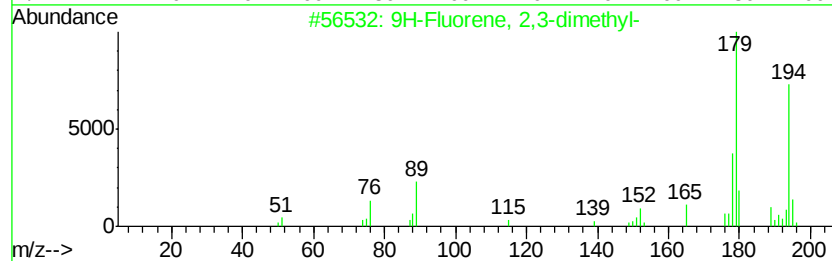
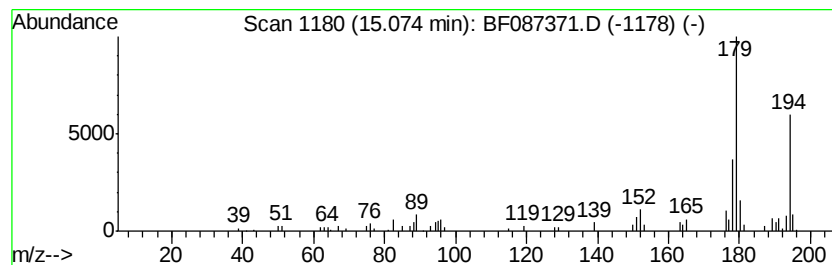
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 9H-Fluorene, 2,3-dimethyl- Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.07 | 3.26 ng | 159550 | Phenanthrene-d10 | 14.78 |

| Hit# | of | Tentative ID                           | MW  | MolForm    | CAS#        | Qual |
|------|----|----------------------------------------|-----|------------|-------------|------|
| 1    | 5  | 9H-Fluorene, 2,3-dimethyl-             | 194 | C15H14     | 004612-63-9 | 91   |
| 2    |    | 9H-Fluorene, 9,9-dimethyl-             | 194 | C15H14     | 004569-45-3 | 87   |
| 3    |    | Benzaldehyde, 4-[(trimethylsilyl)oxy]- | 194 | C10H14O2Si | 001012-12-0 | 72   |
| 4    |    | Anthracene, 9,10-dihydro-2-methyl-     | 194 | C15H14     | 000948-67-4 | 64   |
| 5    |    | 9H-Fluorene, 1,9-dimethyl-             | 194 | C15H14     | 017057-98-6 | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

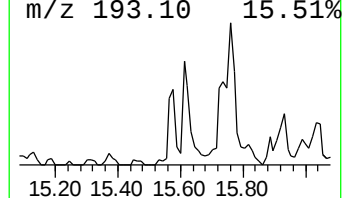
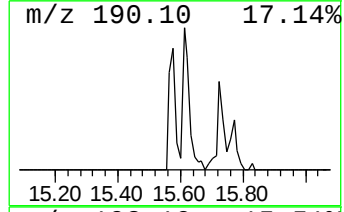
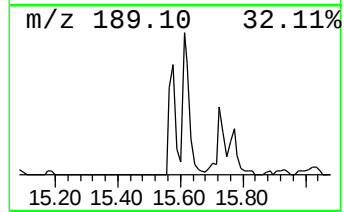
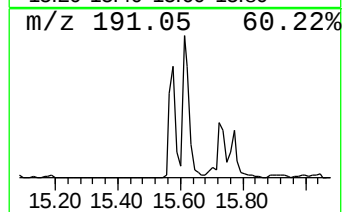
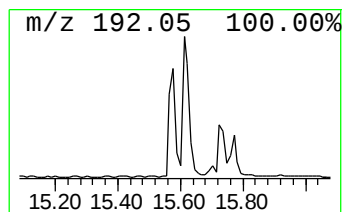
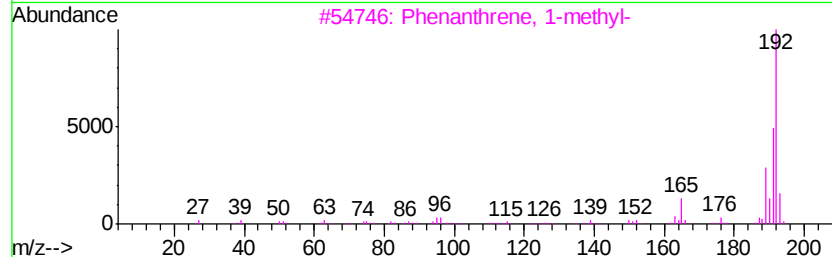
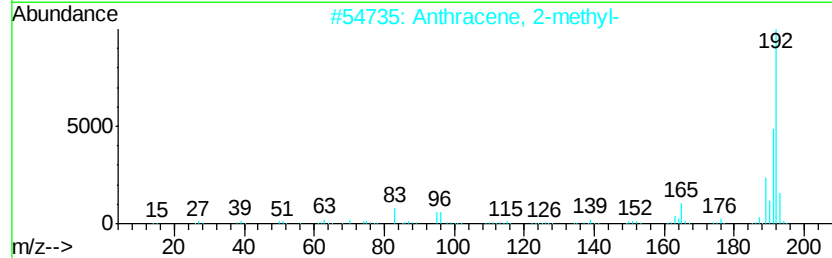
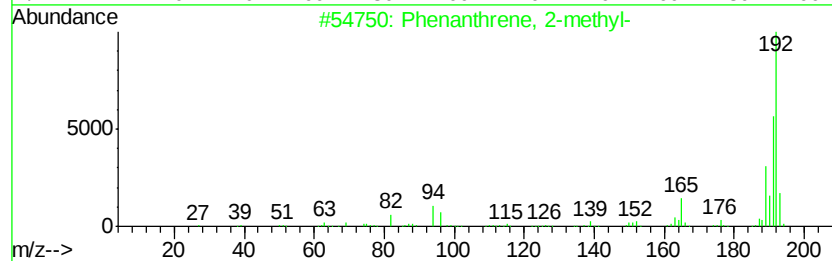
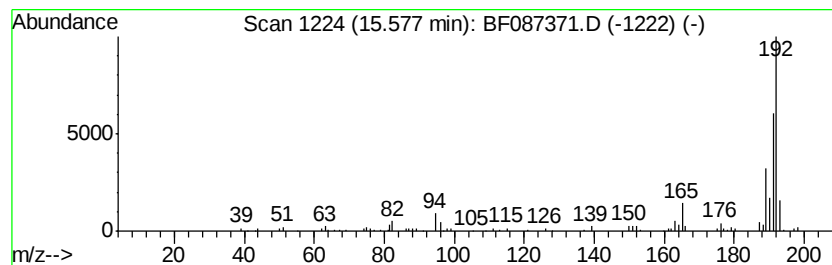
Instrument :  
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ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Phenanthrene, 2-methyl- Concentration Rank 12

| R.T.      | EstConc                 | Area   | Relative to ISTD |             | R.T.  |
|-----------|-------------------------|--------|------------------|-------------|-------|
| 15.58     | 3.61 ng                 | 176389 | Phenanthrene-d10 |             | 14.78 |
| Hit# of 5 | Tentative ID            | MW     | MolForm          | CAS#        | Qual  |
| 1         | Phenanthrene, 2-methyl- | 192    | C15H12           | 002531-84-2 | 96    |
| 2         | Anthracene, 2-methyl-   | 192    | C15H12           | 000613-12-7 | 94    |
| 3         | Phenanthrene, 1-methyl- | 192    | C15H12           | 000832-69-9 | 93    |
| 4         | Anthracene, 9-methyl-   | 192    | C15H12           | 000779-02-2 | 93    |
| 5         | Anthracene, 1-methyl-   | 192    | C15H12           | 000610-48-0 | 91    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF052516\  
Data File : BF087371.D  
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Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

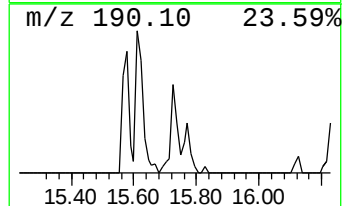
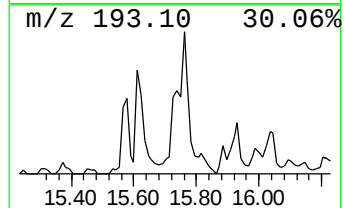
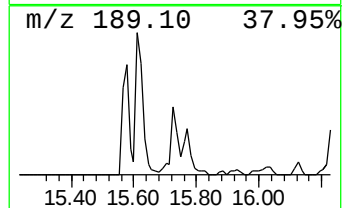
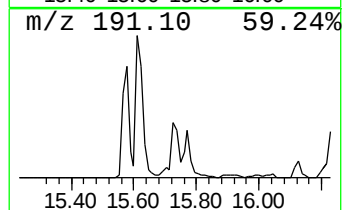
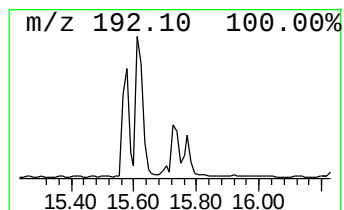
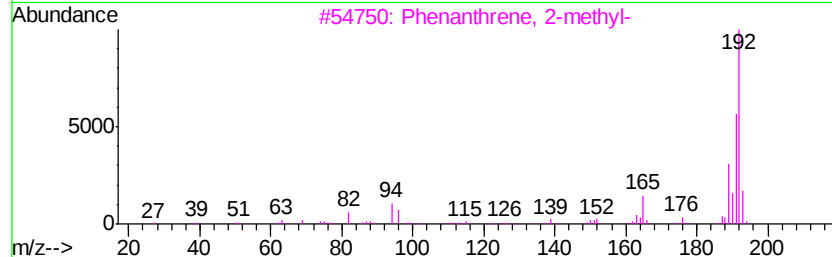
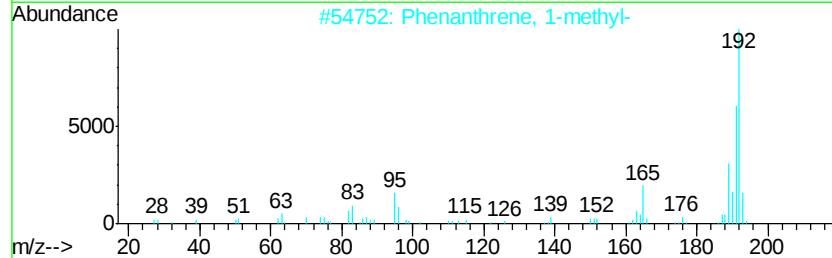
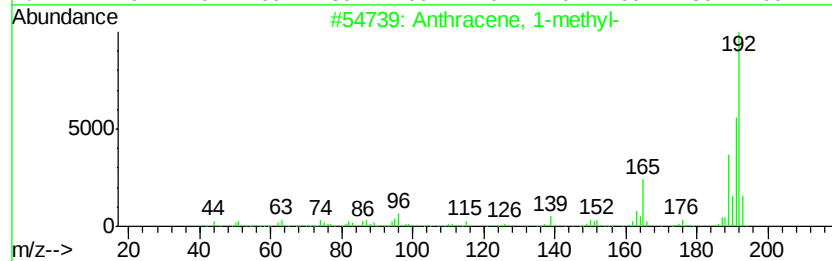
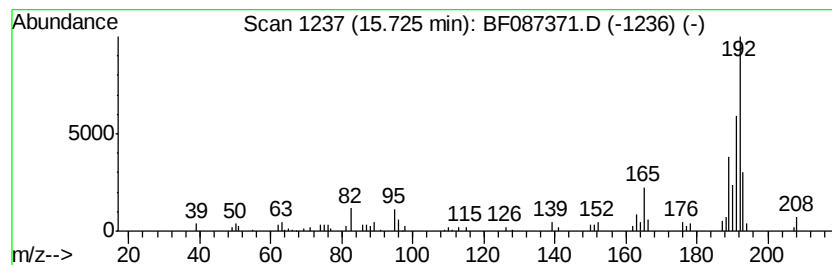
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TIC Integration Parameters: LSCINT.P

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Peak Number 13 Anthracene, 1-methyl- Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.73 | 2.42 ng | 118128 | Phenanthrene-d10 | 14.78 |

| Hit# | of | 5 | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 1-methyl-                 | 192 | C15H12  | 000610-48-0 | 92   |
| 2    |    |   | Phenanthrene, 1-methyl-               | 192 | C15H12  | 000832-69-9 | 89   |
| 3    |    |   | Phenanthrene, 2-methyl-               | 192 | C15H12  | 002531-84-2 | 76   |
| 4    |    |   | Anthracene, 9-methyl-                 | 192 | C15H12  | 000779-02-2 | 76   |
| 5    |    |   | 1H-Cyclopropa[1]phenanthrene, 1a, ... | 192 | C15H12  | 000949-41-7 | 76   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

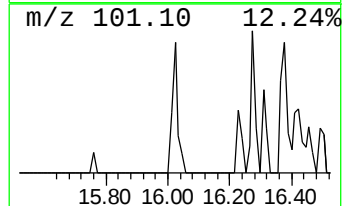
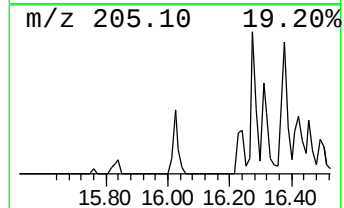
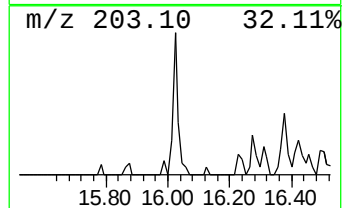
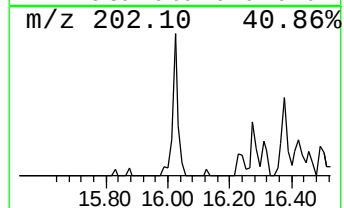
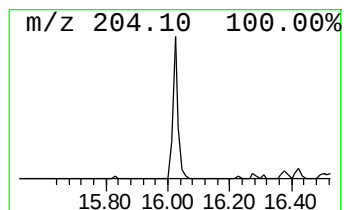
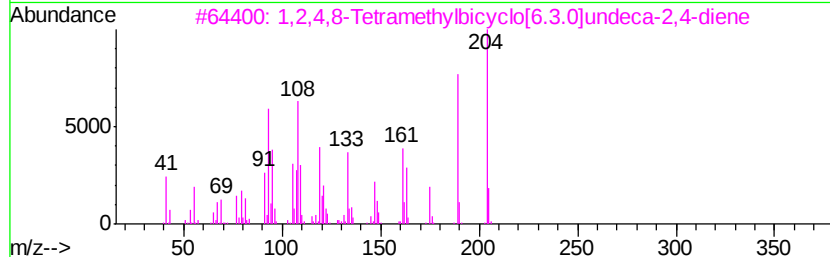
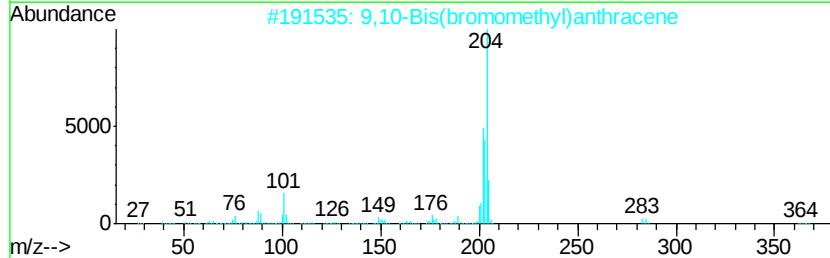
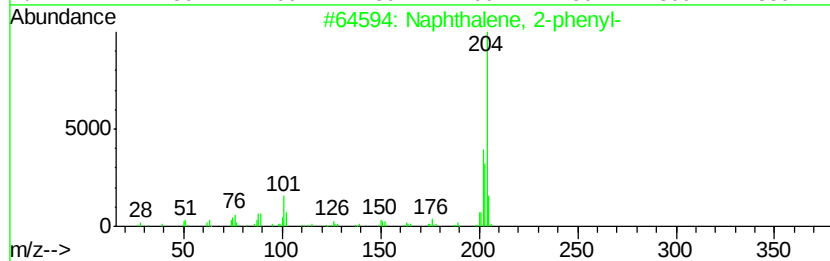
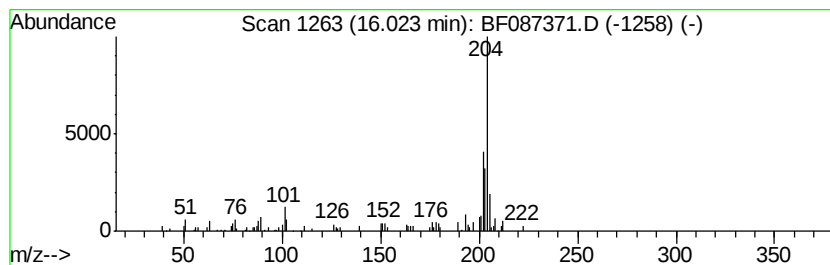
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 Naphthalene, 2-phenyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.02 | 3.90 ng | 190670 | Phenanthrene-d10 | 14.78 |

| Hit# | of | Tentative ID                                            | MW  | MolForm   | CAS#        | Qual |
|------|----|---------------------------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Naphthalene, 2-phenyl-                                  | 204 | C16H12    | 000612-94-2 | 91   |
| 2    |    | 9,10-Bis(bromomethyl)anthracene                         | 362 | C16H12Br2 | 034373-96-1 | 90   |
| 3    |    | 1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene       | 204 | C15H24    | 137235-51-9 | 86   |
| 4    |    | Tricyclo[8.2.2.2(4,7)]hexadeca-2,4,6,8-tetraene         | 204 | C16H12    | 006572-60-7 | 83   |
| 5    |    | 5,16[1',2']:8,13[1'',2'']-Dibenz[1,2-b:4,5-b']pentalene | 408 | C32H24    | 005672-97-9 | 80   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

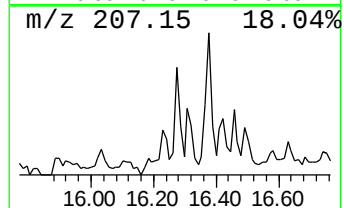
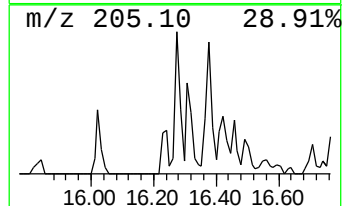
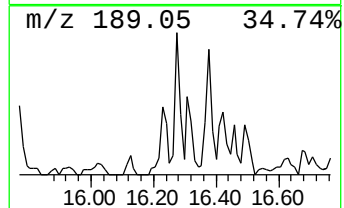
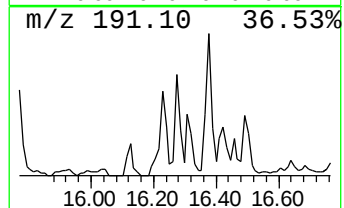
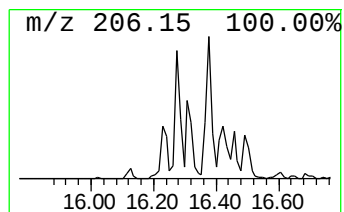
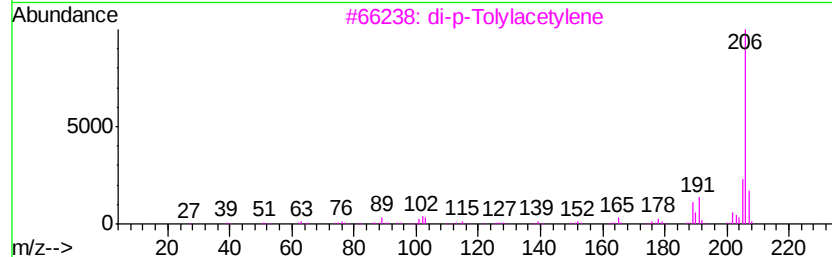
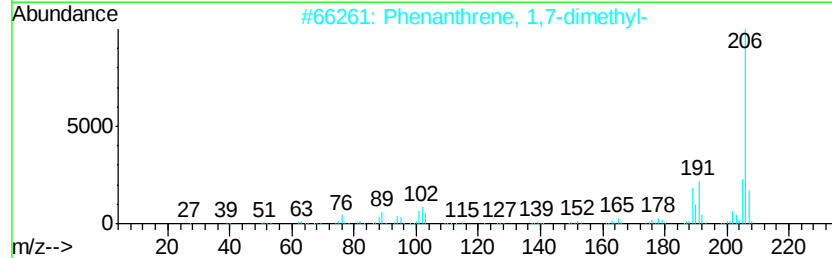
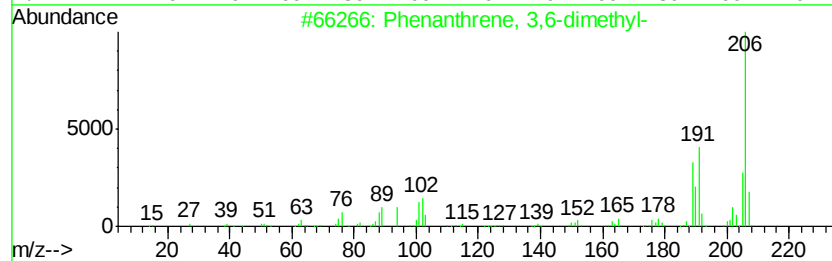
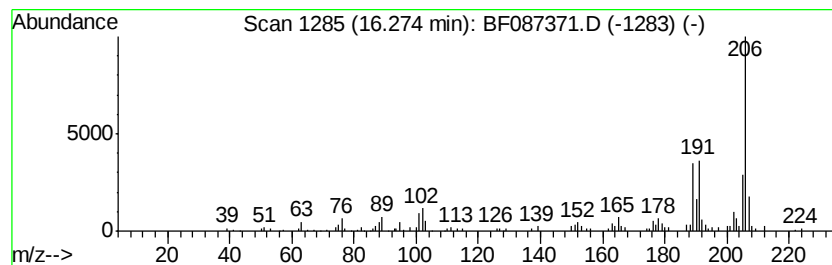
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TIC Integration Parameters: LSCINT.P

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Peak Number 15 Phenanthrene, 3,6-dimethyl- Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.27 | 2.56 ng | 125341 | Phenanthrene-d10 | 14.78 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 97   |
| 2    |    |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 94   |
| 3    |    |   | di-p-Tolylacetvlene         | 206 | C16H14  | 002789-88-0 | 94   |
| 4    |    |   | Phenanthrene, 2,7-dimethyl- | 206 | C16H14  | 001576-69-8 | 93   |
| 5    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

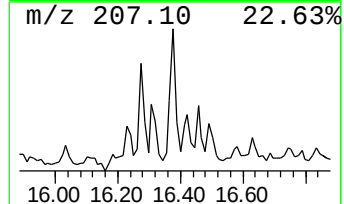
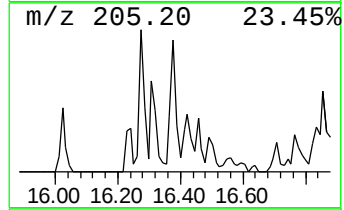
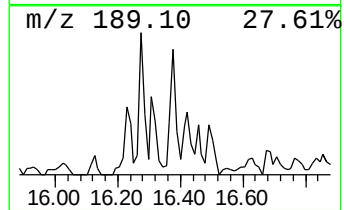
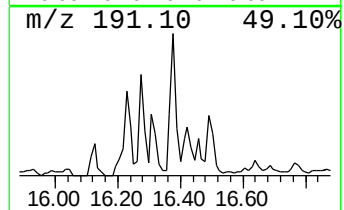
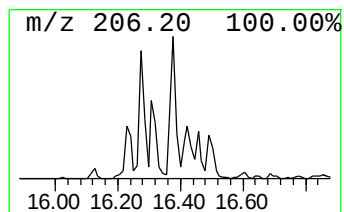
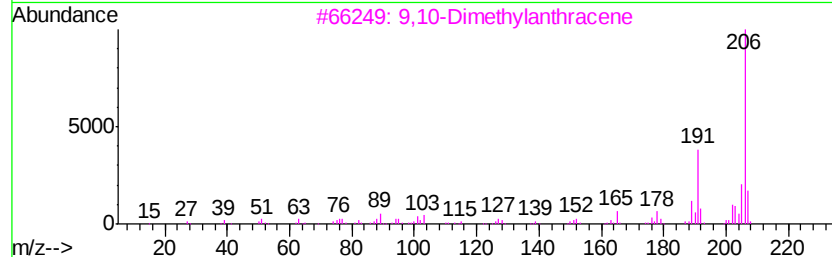
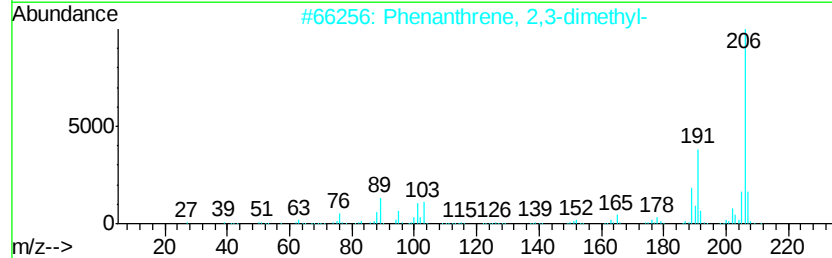
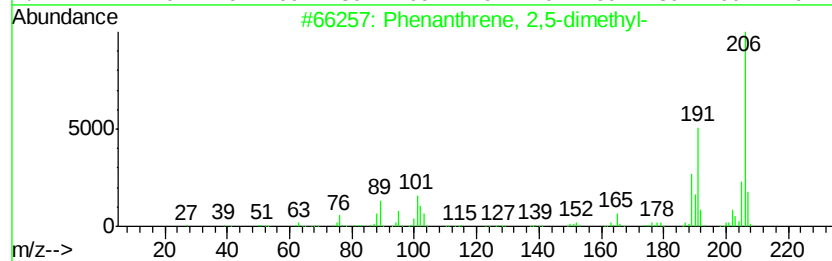
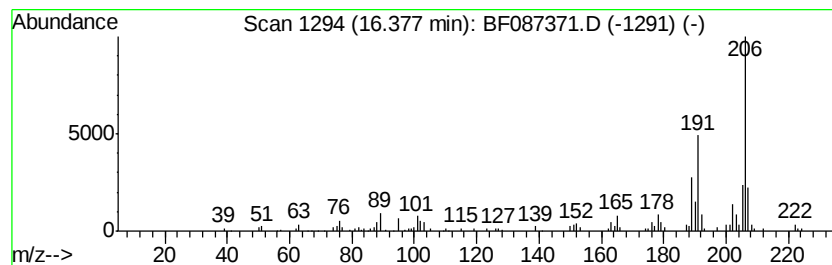
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Phenanthrene, 2,5-dimethyl- Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.38 | 3.56 ng | 173873 | Phenanthrene-d10 | 14.78 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 95   |
| 2    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 93   |
| 3    |    |   | 9,10-Dimethylanthracene     | 206 | C16H14  | 000781-43-1 | 91   |
| 4    |    |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 90   |
| 5    |    |   | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 89   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

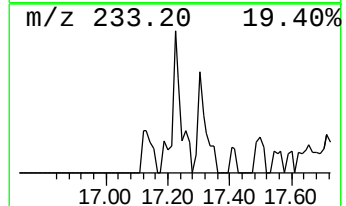
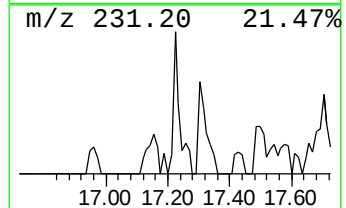
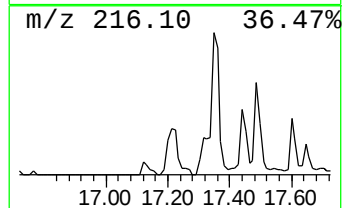
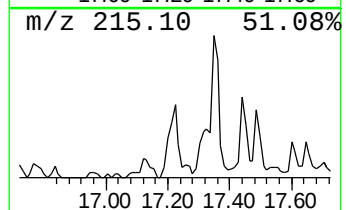
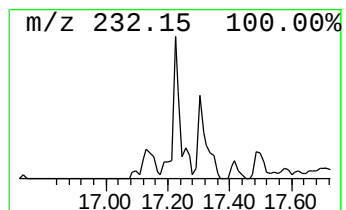
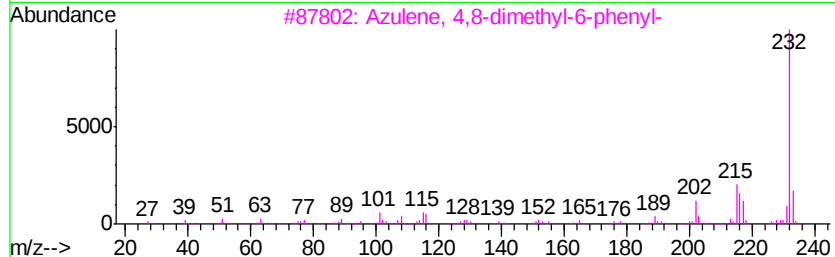
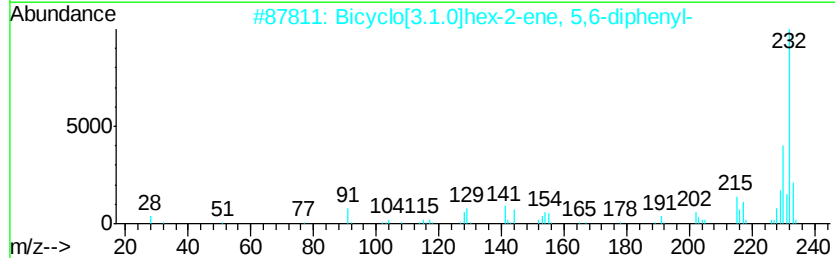
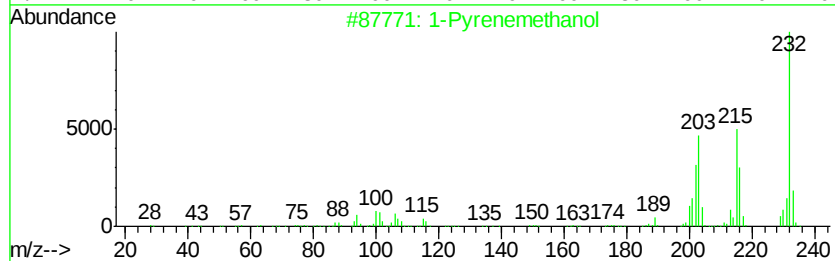
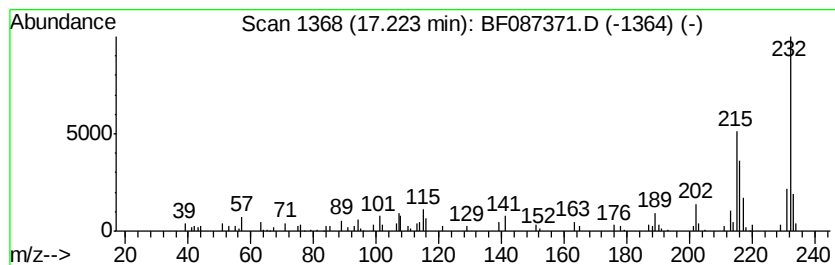
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 1-Pyrenemethanol Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.22 | 2.46 ng | 112511 | Chrysene-d12     | 18.47 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 1-Pyrenemethanol                    | 232 | C17H12O | 024463-15-8  | 91   |
| 2    |    | Bicyclo[3.1.0]hex-2-ene, 5,6-dip... | 232 | C18H16  | 056143-24-9  | 59   |
| 3    |    | Azulene, 4,8-dimethyl-6-phenyl-     | 232 | C18H16  | 042758-88-3  | 58   |
| 4    |    | 1,3-dimethyl-2-phenyl-naphthalene   | 232 | C18H16  | 1000379-93-4 | 58   |
| 5    |    | 1-benzyl-3-methylnaphthalene        | 232 | C18H16  | 1000379-93-5 | 56   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

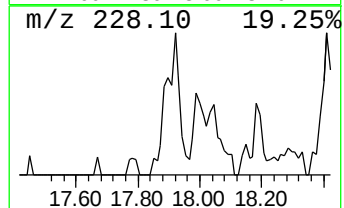
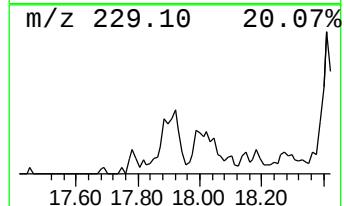
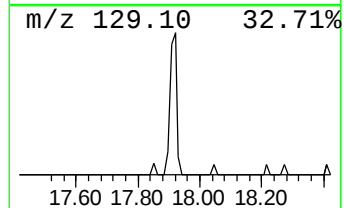
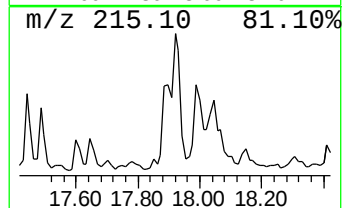
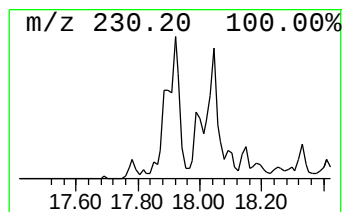
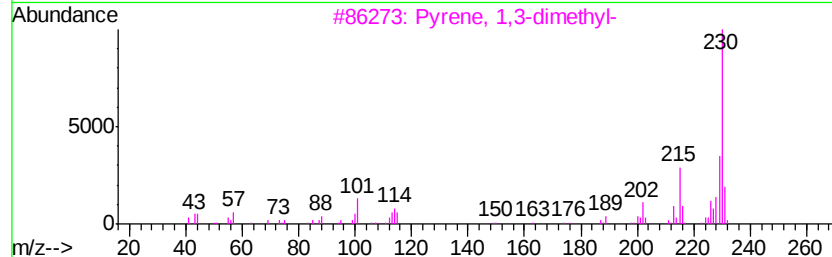
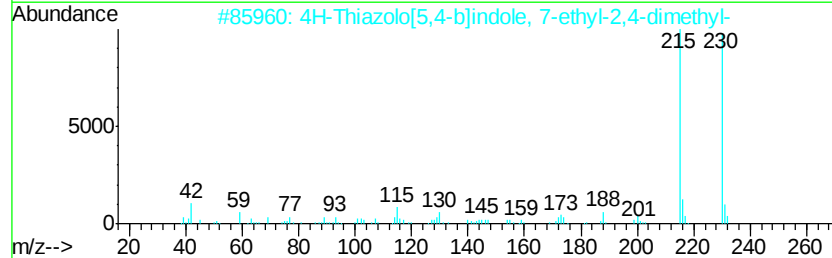
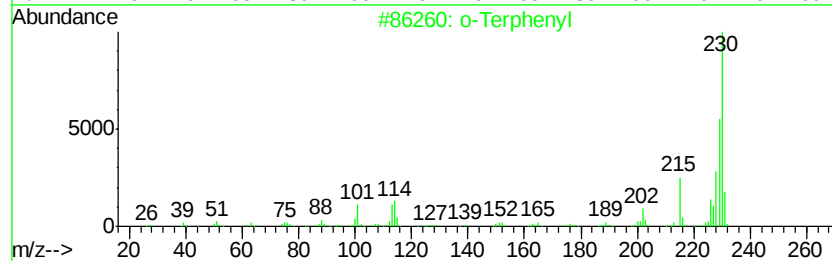
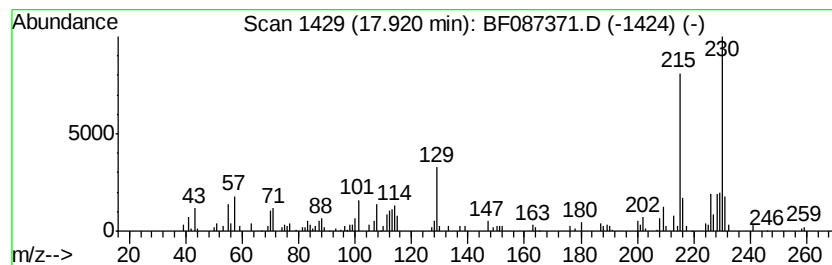
Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 18 o-Terphenyl Concentration Rank 5

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 17.92     | 5.16 ng                             | 235834 | Chrysene-d12     | 18.47        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | o-Terphenyl                         | 230    | C18H14           | 000084-15-1  | 64   |
| 2         | 4H-Thiazolo[5,4-b]indole, 7-ethy... | 230    | C13H14N2S        | 1000304-41-6 | 52   |
| 3         | Pvrene, 1,3-dimethyl-               | 230    | C18H14           | 064401-21-4  | 50   |
| 4         | 6,6-Diphenylfulvene                 | 230    | C18H14           | 002175-90-8  | 49   |
| 5         | Benzo[g]pteridine, 2,4-diamino-6... | 230    | C11H14N6         | 063630-26-2  | 43   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

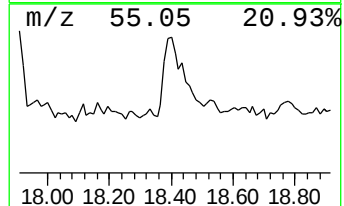
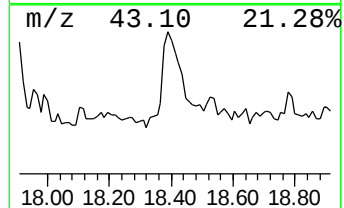
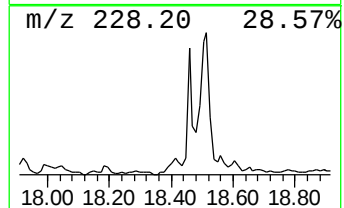
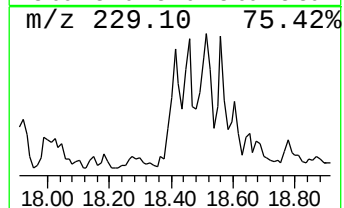
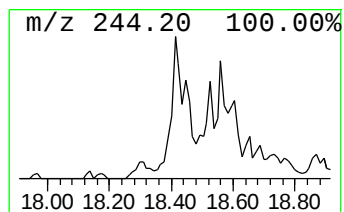
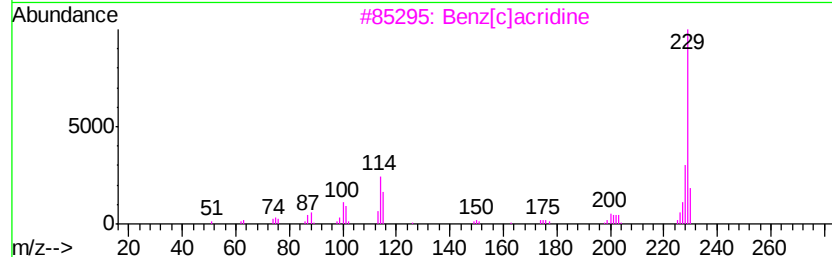
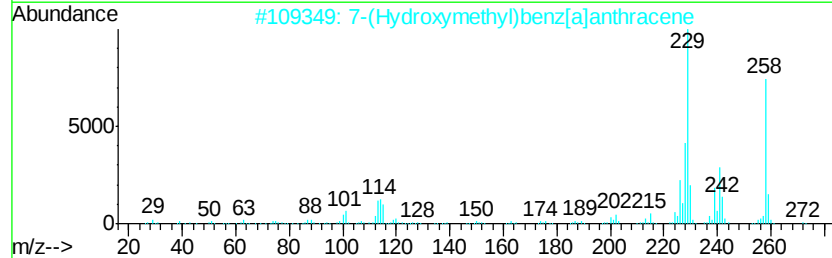
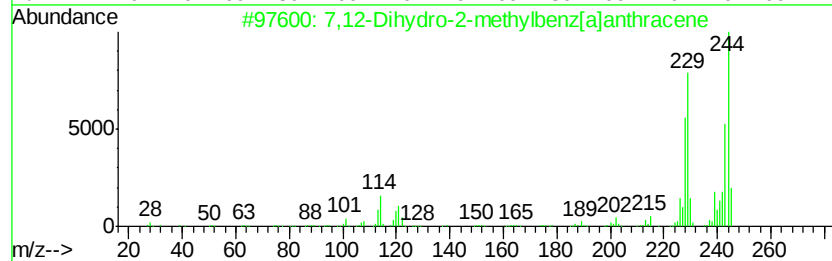
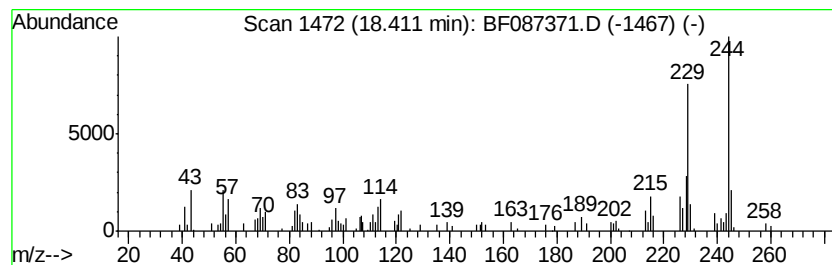
Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 7,12-Dihydro-2-methylbenz[a... Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 18.41     | 4.64 ng                             | 212171 | Chrysene-d12     | 18.47        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 7,12-Dihydro-2-methylbenz[a]anth... | 244    | C19H16           | 035187-44-1  | 81   |
| 2         | 7-(Hydroxymethyl)benz[a]anthracene  | 258    | C19H14O          | 1000326-13-2 | 70   |
| 3         | Benz[c]acridine                     | 229    | C17H11N          | 000225-51-4  | 60   |
| 4         | 4-Benzylbiphenyl                    | 244    | C19H16           | 000613-42-3  | 60   |
| 5         | 6-Acetyl-5-hydroxy-2,7-dimethyl-... | 244    | C14H12O4         | 001086-07-3  | 59   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF052516\  
Data File : BF087371.D  
Acq On : 25 May 2016 14:29  
Operator : UM/SJ  
Sample : H3218-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

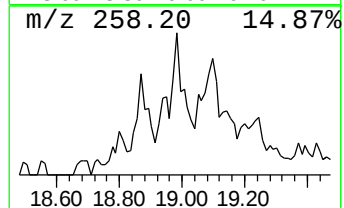
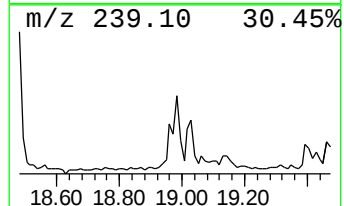
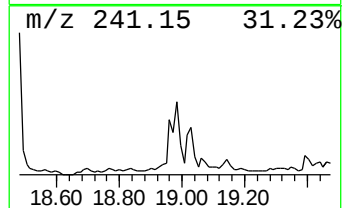
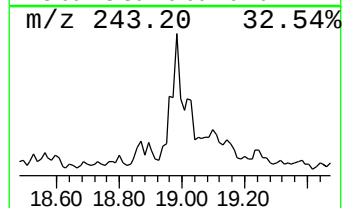
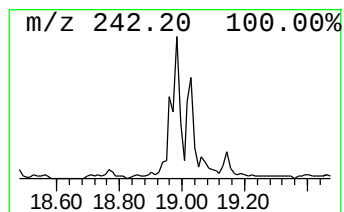
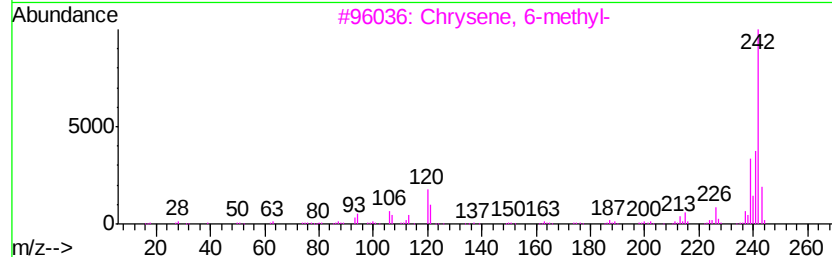
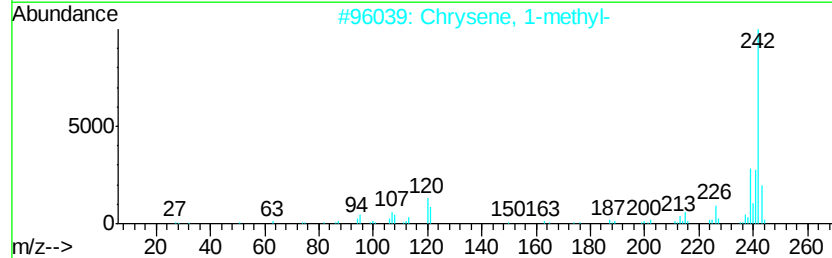
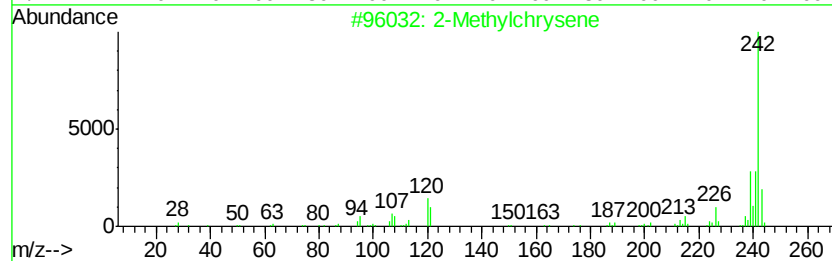
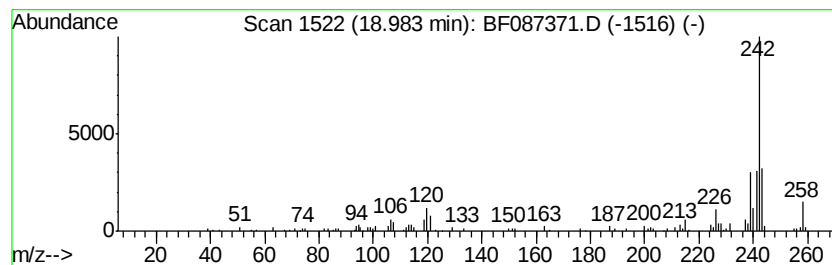
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 2-Methylchrysene Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.98 | 3.87 ng | 176958 | Chrysene-d12     | 18.47 |

| Hit# | of | 5 | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Methylchrysene                      | 242 | C19H14  | 003351-32-4 | 93   |
| 2    |    |   | Chrysene, 1-methyl-                   | 242 | C19H14  | 003351-28-8 | 93   |
| 3    |    |   | Chrysene, 6-methyl-                   | 242 | C19H14  | 001705-85-7 | 93   |
| 4    |    |   | Benz[ <i>a</i> ]anthracene, 1-methyl- | 242 | C19H14  | 002498-77-3 | 92   |
| 5    |    |   | Chrysene, 5-methyl-                   | 242 | C19H14  | 003697-24-3 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF052516\  
 Data File : BF087371.D  
 Acq On : 25 May 2016 14:29  
 Operator : UM/SJ  
 Sample : H3218-01  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 E-3(6.5-7)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF052316.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| 2-Pentanone, 4-hy... | 4.78  | 4.1 ng  |       | 143868   | 1                     | 7.52  | 707338 | 20.0 |
| unknown7.16          | 7.16  | 68.7 ng |       | 2430440  | 1                     | 7.52  | 707338 | 20.0 |
| Octane               | 7.31  | 3.2 ng  |       | 114117   | 1                     | 7.52  | 707338 | 20.0 |
| Ethanone, 1-pheny... | 8.65  | 3.7 ng  |       | 171876   | 2                     | 9.55  | 927216 | 20.0 |
| Benzene, 1,2,4,5-... | 9.18  | 2.7 ng  |       | 123925   | 2                     | 9.55  | 927216 | 20.0 |
| Hexadecane           | 10.62 | 4.4 ng  |       | 204182   | 2                     | 9.55  | 927216 | 20.0 |
| Formamide, N-phenyl- | 10.94 | 3.0 ng  |       | 137358   | 2                     | 9.55  | 927216 | 20.0 |
| 9H-Fluorene, 2,3-... | 15.07 | 3.3 ng  |       | 159550   | 4                     | 14.78 | 978055 | 20.0 |
| Phenanthrene, 2-m... | 15.58 | 3.6 ng  |       | 176389   | 4                     | 14.78 | 978055 | 20.0 |
| Anthracene, 1-met... | 15.73 | 2.4 ng  |       | 118128   | 4                     | 14.78 | 978055 | 20.0 |
| Naphthalene, 2-ph... | 16.02 | 3.9 ng  |       | 190670   | 4                     | 14.78 | 978055 | 20.0 |
| Phenanthrene, 3,6... | 16.27 | 2.6 ng  |       | 125341   | 4                     | 14.78 | 978055 | 20.0 |
| Phenanthrene, 2,5... | 16.38 | 3.6 ng  |       | 173873   | 4                     | 14.78 | 978055 | 20.0 |
| 1-Pyrenemethanol     | 17.22 | 2.5 ng  |       | 112511   | 5                     | 18.47 | 914347 | 20.0 |
| o-Terphenyl          | 17.92 | 5.2 ng  |       | 235834   | 5                     | 18.47 | 914347 | 20.0 |
| 7,12-Dihydro-2-me... | 18.41 | 4.6 ng  |       | 212171   | 5                     | 18.47 | 914347 | 20.0 |
| 2-Methylchrysene     | 18.98 | 3.9 ng  |       | 176958   | 5                     | 18.47 | 914347 | 20.0 |