

Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\  
 Data File : BF094204.D  
 Acq On : 5 Apr 2017 22:06  
 Operator : SJ/MA  
 Sample : I2522-17  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 BASING-9(0-2)

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-BF032217.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         | 1.922       | 3             | 6           | 13           | rVB      | 49130          | 61508         | 1.42%           | 0.141%        |
| 2         | 4.010       | 356           | 361         | 370          | rBV      | 438095         | 601390        | 13.88%          | 1.383%        |
| 3         | 4.404       | 422           | 428         | 431          | rBV      | 3876792        | 3865249       | 89.23%          | 8.887%        |
| 4         | 5.475       | 603           | 610         | 613          | rBV      | 3401915        | 4026595       | 92.96%          | 9.258%        |
| 5         | 5.610       | 628           | 633         | 636          | rBV      | 3854587        | 4183173       | 96.57%          | 9.618%        |
| 6         | 5.863       | 672           | 676         | 680          | rBV      | 1274160        | 1089847       | 25.16%          | 2.506%        |
| 7         | 6.045       | 702           | 707         | 710          | rBV      | 3166388        | 3206145       | 74.02%          | 7.372%        |
| 8         | 6.516       | 780           | 787         | 790          | rBV      | 2457083        | 2705917       | 62.47%          | 6.222%        |
| 9         | 7.363       | 926           | 931         | 935          | rBV      | 1392736        | 1467467       | 33.88%          | 3.374%        |
| 10        | 8.692       | 1150          | 1157        | 1160         | rVB      | 4052061        | 4331631       | 100.00%         | 9.960%        |
| 11        | 8.757       | 1165          | 1168        | 1170         | rBV2     | 55302          | 60074         | 1.39%           | 0.138%        |
| 12        | 8.845       | 1178          | 1183        | 1188         | rVB3     | 51711          | 55474         | 1.28%           | 0.128%        |
| 13        | 8.998       | 1203          | 1209        | 1215         | rVB2     | 32056          | 50755         | 1.17%           | 0.117%        |
| 14        | 9.098       | 1219          | 1226        | 1229         | rVV2     | 141302         | 175829        | 4.06%           | 0.404%        |
| 15        | 9.157       | 1233          | 1236        | 1238         | rVV      | 111079         | 117525        | 2.71%           | 0.270%        |
| 16        | 9.186       | 1238          | 1241        | 1244         | rVV      | 342903         | 330805        | 7.64%           | 0.761%        |
| 17        | 9.233       | 1246          | 1249        | 1256         | rVB3     | 30700          | 49232         | 1.14%           | 0.113%        |
| 18        | 9.333       | 1262          | 1266        | 1270         | rBV2     | 70214          | 80851         | 1.87%           | 0.186%        |
| 19        | 9.422       | 1276          | 1281        | 1284         | rBV      | 148112         | 153781        | 3.55%           | 0.354%        |
| 20        | 9.504       | 1288          | 1295        | 1299         | rVV2     | 1549462        | 1646996       | 38.02%          | 3.787%        |
| 21        | 9.545       | 1299          | 1302        | 1304         | rVV      | 178505         | 196877        | 4.55%           | 0.453%        |
| 22        | 9.569       | 1304          | 1306        | 1309         | rVB      | 238274         | 196449        | 4.54%           | 0.452%        |
| 23        | 9.686       | 1321          | 1326        | 1329         | rBV2     | 143497         | 181092        | 4.18%           | 0.416%        |
| 24        | 9.727       | 1329          | 1333        | 1335         | rVV2     | 102090         | 125141        | 2.89%           | 0.288%        |
| 25        | 9.757       | 1335          | 1338        | 1343         | rVV      | 267455         | 276792        | 6.39%           | 0.636%        |
| 26        | 9.816       | 1346          | 1348        | 1353         | rVV3     | 72405          | 74649         | 1.72%           | 0.172%        |
| 27        | 9.857       | 1353          | 1355        | 1361         | rVB2     | 35576          | 52055         | 1.20%           | 0.120%        |
| 28        | 10.098      | 1391          | 1396        | 1401         | rBV      | 146284         | 161082        | 3.72%           | 0.370%        |
| 29        | 10.151      | 1401          | 1405        | 1407         | rVV4     | 52869          | 77608         | 1.79%           | 0.178%        |
| 30        | 10.174      | 1407          | 1409        | 1413         | rVB      | 222800         | 218871        | 5.05%           | 0.503%        |
| 31        | 10.380      | 1436          | 1444        | 1451         | rVB4     | 100648         | 203291        | 4.69%           | 0.467%        |
| 32        | 10.480      | 1455          | 1461        | 1465         | rVV      | 2117386        | 2367726       | 54.66%          | 5.444%        |
| 33        | 10.545      | 1469          | 1472        | 1477         | rVB4     | 93282          | 95915         | 2.21%           | 0.221%        |
| 34        | 10.645      | 1486          | 1489        | 1492         | rBV      | 91608          | 91071         | 2.10%           | 0.209%        |

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 BASING-9(0-2)

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 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032217.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 10.680 | 1492 | 1495 | 1506 | rVB  | 144361  | 203708  | 4.70%  | 0.468% |
| 36 | 11.086 | 1560 | 1564 | 1567 | rBV  | 116038  | 108367  | 2.50%  | 0.249% |
| 37 | 11.198 | 1580 | 1583 | 1586 | rVB  | 65549   | 57312   | 1.32%  | 0.132% |
| 38 | 11.239 | 1586 | 1590 | 1593 | rVB  | 87580   | 89369   | 2.06%  | 0.205% |
| 39 | 11.327 | 1600 | 1605 | 1607 | rBV  | 1383219 | 1357862 | 31.35% | 3.122% |
| 40 | 11.357 | 1607 | 1610 | 1615 | rVV  | 984830  | 929029  | 21.45% | 2.136% |
| 41 | 11.416 | 1617 | 1620 | 1625 | rVB  | 81344   | 78974   | 1.82%  | 0.182% |
| 42 | 11.745 | 1672 | 1676 | 1678 | rBV3 | 32571   | 43426   | 1.00%  | 0.100% |
| 43 | 11.951 | 1707 | 1711 | 1714 | rBV  | 89434   | 92053   | 2.13%  | 0.212% |
| 44 | 11.986 | 1714 | 1717 | 1723 | rVB  | 84394   | 102121  | 2.36%  | 0.235% |
| 45 | 12.080 | 1730 | 1733 | 1737 | rBV2 | 98079   | 102946  | 2.38%  | 0.237% |
| 46 | 12.116 | 1737 | 1739 | 1751 | rVB2 | 33461   | 47555   | 1.10%  | 0.109% |
| 47 | 12.321 | 1771 | 1774 | 1780 | rBV2 | 45396   | 65994   | 1.52%  | 0.152% |
| 48 | 12.815 | 1850 | 1858 | 1862 | rBV  | 437117  | 446362  | 10.30% | 1.026% |
| 49 | 13.092 | 1900 | 1905 | 1909 | rVB2 | 307513  | 334899  | 7.73%  | 0.770% |
| 50 | 13.304 | 1936 | 1941 | 1945 | rBV  | 2372039 | 2723432 | 62.87% | 6.262% |
| 51 | 13.451 | 1962 | 1966 | 1969 | rBV  | 65160   | 68233   | 1.58%  | 0.157% |
| 52 | 13.639 | 1994 | 1998 | 2006 | rVB6 | 27116   | 53271   | 1.23%  | 0.122% |
| 53 | 14.474 | 2135 | 2140 | 2150 | rBV3 | 113266  | 224618  | 5.19%  | 0.516% |
| 54 | 14.574 | 2152 | 2157 | 2161 | rBV  | 958509  | 1068896 | 24.68% | 2.458% |
| 55 | 15.257 | 2269 | 2273 | 2281 | rBV3 | 85888   | 173171  | 4.00%  | 0.398% |
| 56 | 15.615 | 2331 | 2334 | 2338 | rBV  | 79469   | 90920   | 2.10%  | 0.209% |
| 57 | 15.968 | 2391 | 2394 | 2398 | rBV  | 142569  | 169046  | 3.90%  | 0.389% |
| 58 | 16.203 | 2429 | 2434 | 2442 | rVB  | 781036  | 897342  | 20.72% | 2.063% |
| 59 | 16.327 | 2451 | 2455 | 2459 | rBV  | 219299  | 257898  | 5.95%  | 0.593% |
| 60 | 16.727 | 2518 | 2523 | 2528 | rVB  | 207081  | 302429  | 6.98%  | 0.695% |
| 61 | 17.180 | 2595 | 2600 | 2607 | rBV  | 186867  | 328421  | 7.58%  | 0.755% |
| 62 | 17.709 | 2685 | 2690 | 2698 | rVB2 | 124002  | 245715  | 5.67%  | 0.565% |
| 63 | 18.327 | 2789 | 2795 | 2806 | rVB3 | 93374   | 248961  | 5.75%  | 0.572% |

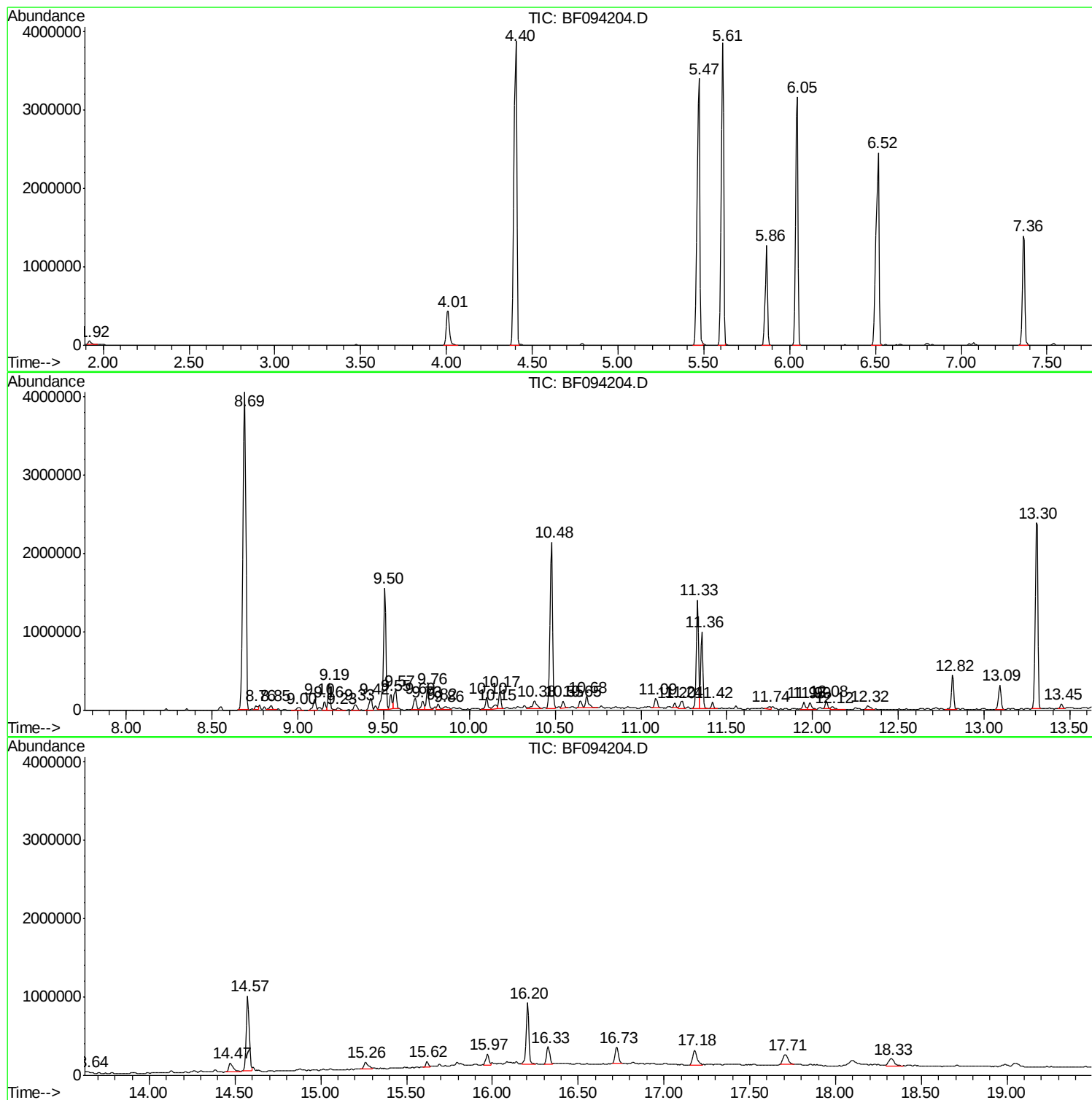
Sum of corrected areas: 43491193

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BNA\_F  
ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.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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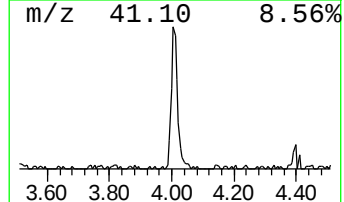
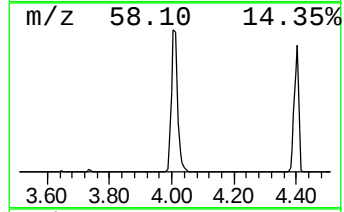
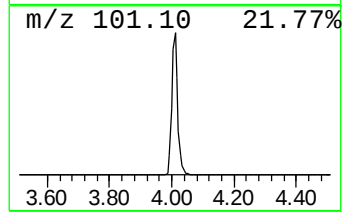
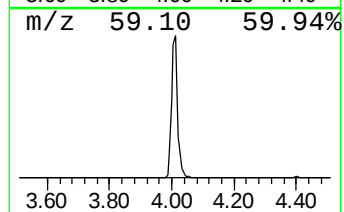
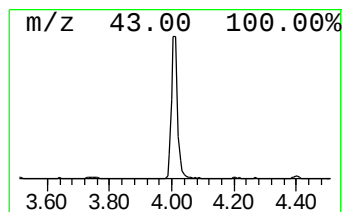
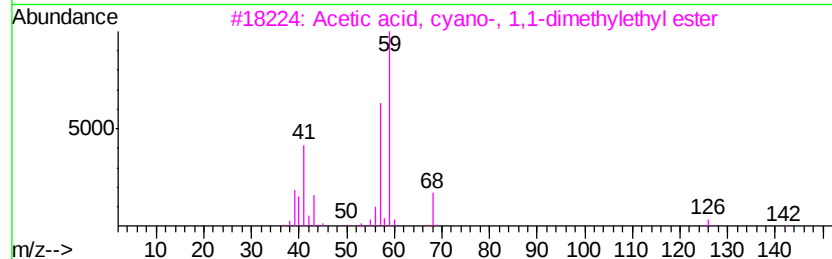
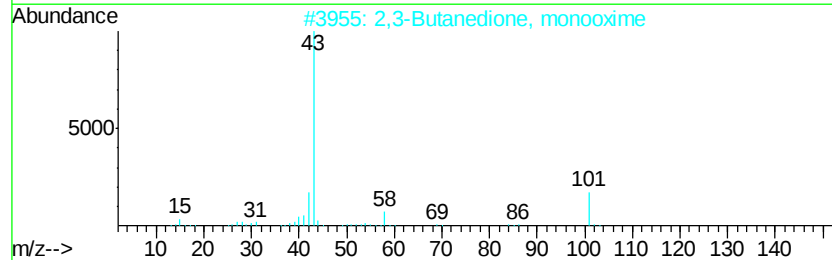
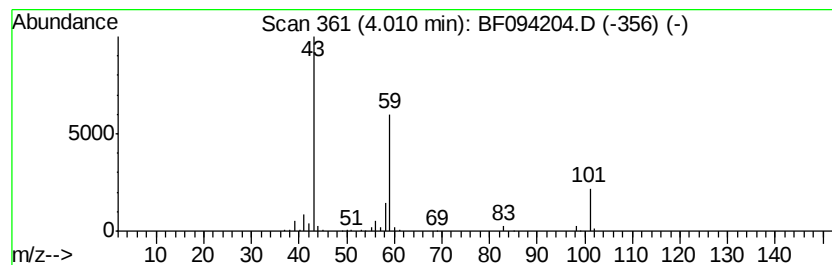
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 4.01 | 11.04 ng | 601390 | 1,4-Dichlorobenzene-d4 | 5.86 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 72   |
| 2    |    |   | 2,3-Butanedione, monooxime           | 101 | C4H7NO2  | 000057-71-6 | 35   |
| 3    |    |   | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 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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BNA\_F  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

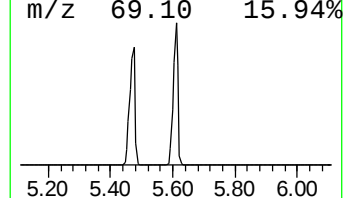
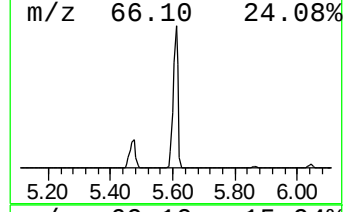
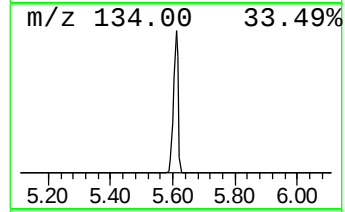
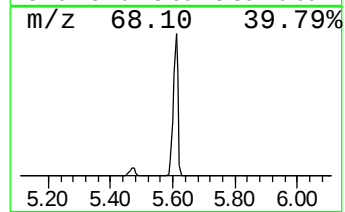
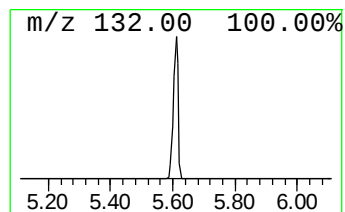
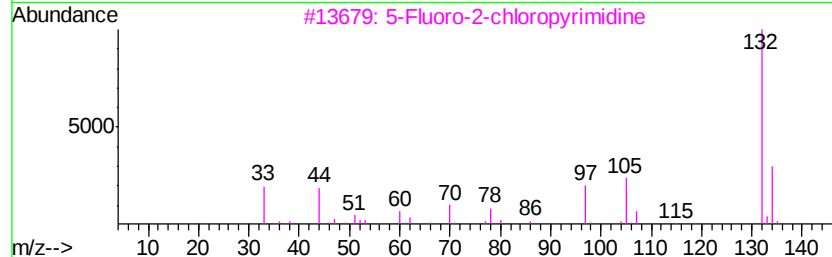
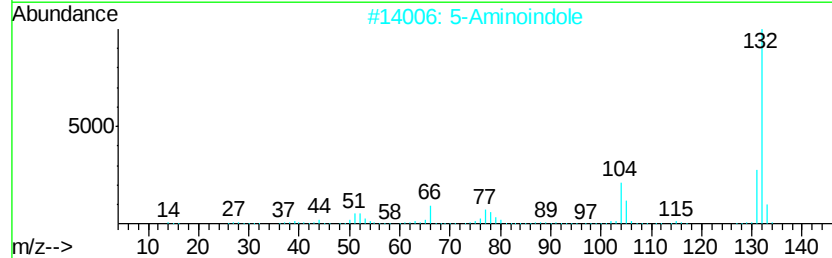
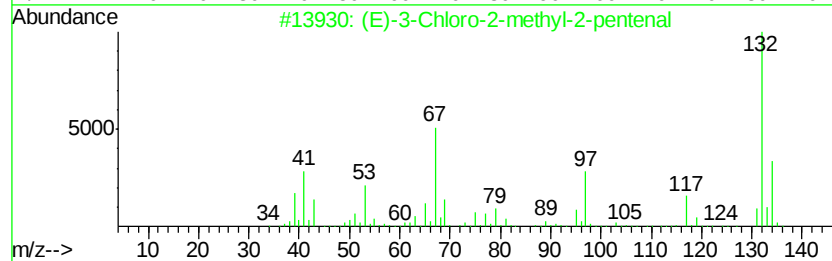
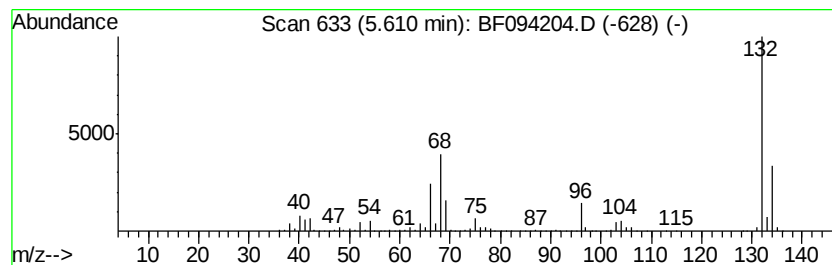
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 unknown5.61 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.61 | 76.77 ng | 4183170 | 1,4-Dichlorobenzene-d4 | 5.86 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#        | Qual |
|------|----|----------------------------------|-----|-----------|-------------|------|
| 1    | 5  | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3 | 27   |
| 2    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0 | 12   |
| 3    |    | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0 | 9    |
| 4    |    | 1-Methylpyrrolo[1,2-a]pyrazine   | 132 | C8H8N2    | 064608-59-9 | 9    |
| 5    |    | 4-Methylpyrrolo[1,2-a]pyrazine   | 132 | C8H8N2    | 064608-60-2 | 9    |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

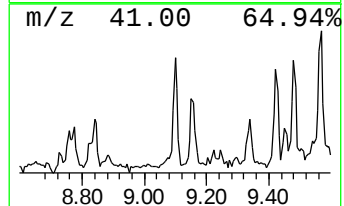
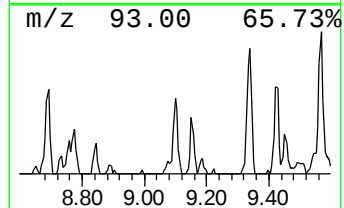
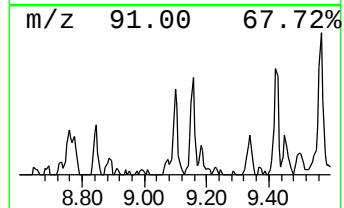
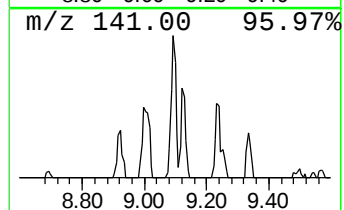
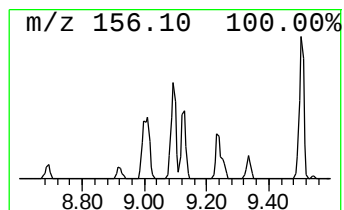
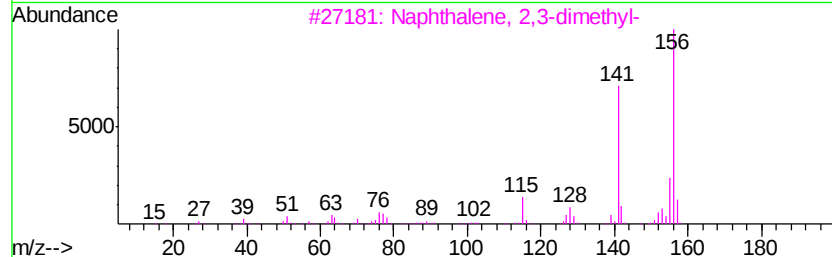
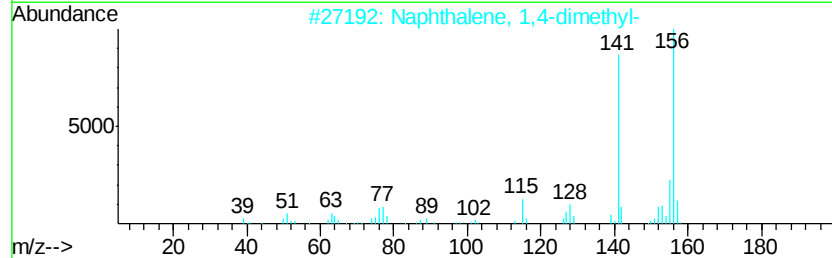
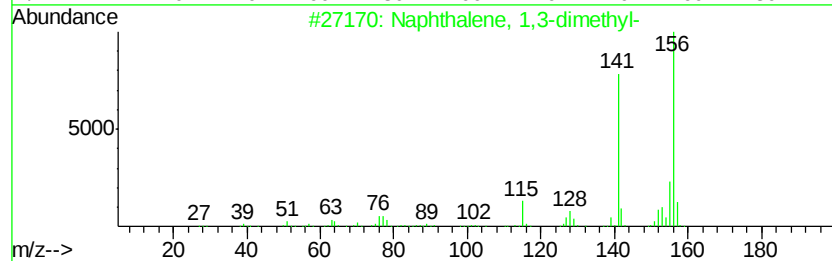
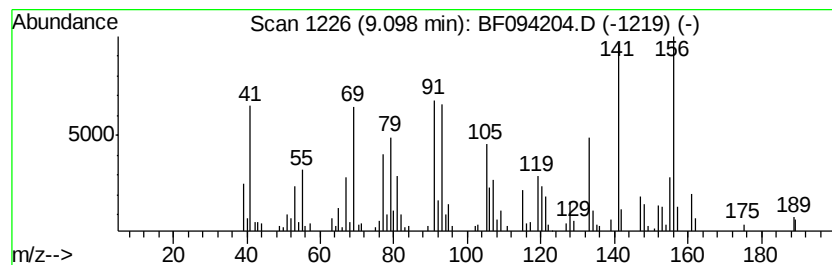
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Naphthalene, 1,3-dimethyl- Concentration Rank 16

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.10 | 2.14 ng | 175829 | Acenaphthene-d10 | 9.50 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 95   |
| 2    |    | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 90   |
| 3    |    | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 90   |
| 4    |    | Naphthalene, 1,8-dimethyl- | 156 | C12H12  | 000569-41-5 | 86   |
| 5    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 80   |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

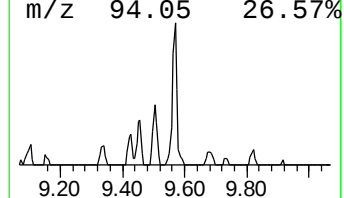
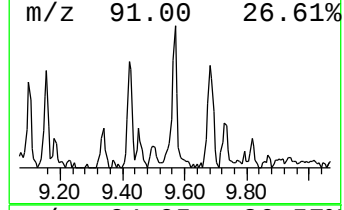
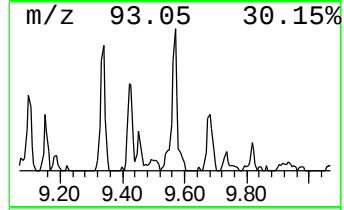
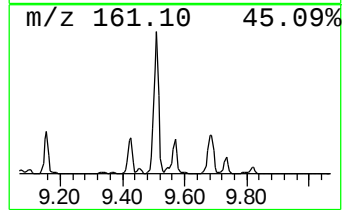
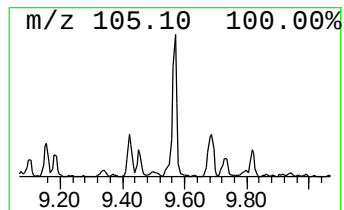
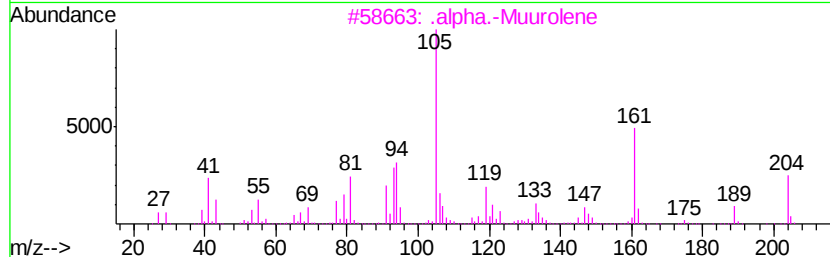
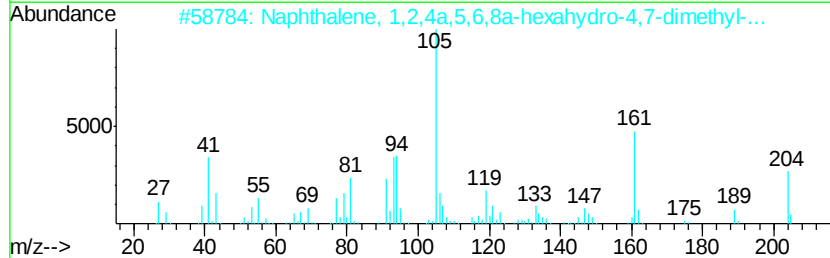
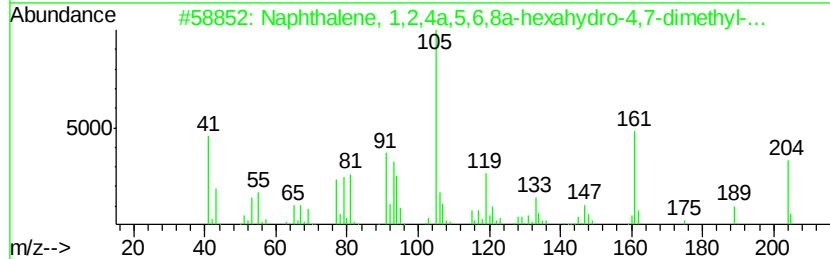
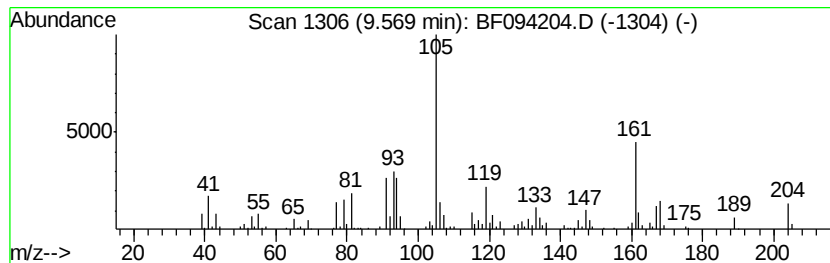
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 Naphthalene, 1,2,4a,5,6,8a-... Concentration Rank 14

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.57 | 2.39 ng | 196449 | Acenaphthene-d10 | 9.50 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,2,4a,5,6,8a-hexah... | 204 | C15H24  | 031983-22-9 | 97   |
| 2    |    | Naphthalene, 1,2,4a,5,6,8a-hexah... | 204 | C15H24  | 000483-75-0 | 96   |
| 3    |    | .alpha.-Muurolene                   | 204 | C15H24  | 010208-80-7 | 92   |
| 4    |    | Naphthalene, 1,2,4a,5,6,8a-hexah... | 204 | C15H24  | 017627-24-6 | 74   |
| 5    |    | Naphthalene, 1,2,4a,5,6,8a-hexah... | 204 | C15H24  | 024406-05-1 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\Data\BF040517\  
Data File : BF094204.D  
Acq On : 5 Apr 2017 22:06  
Operator : SJ/MA  
Sample : I2522-17  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

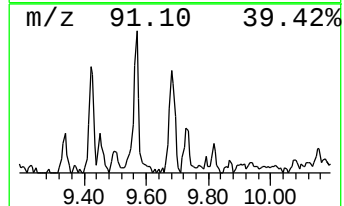
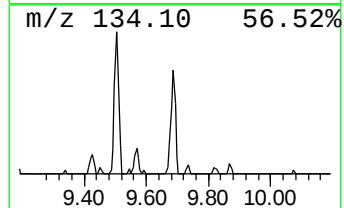
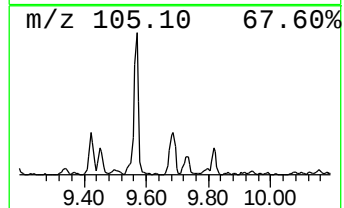
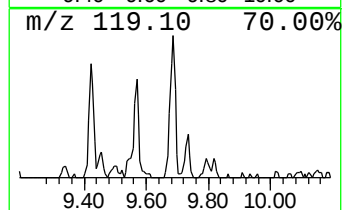
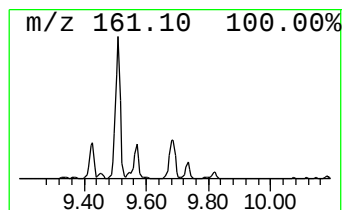
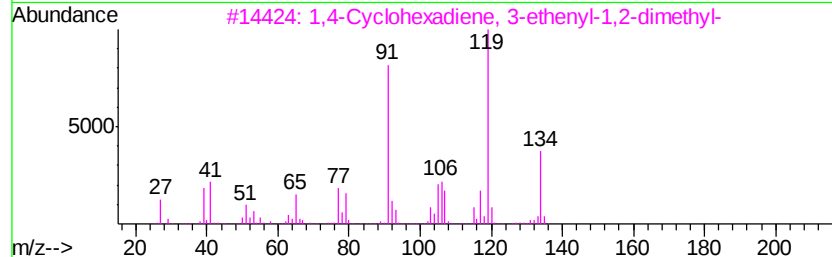
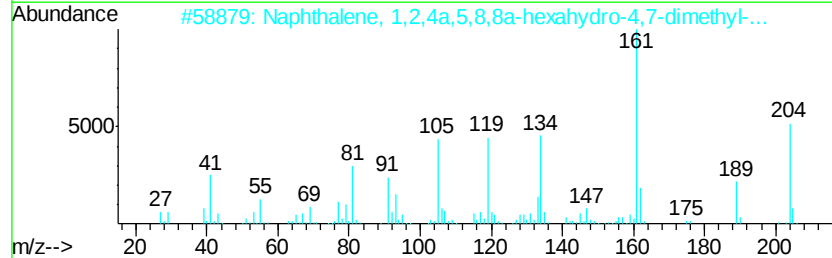
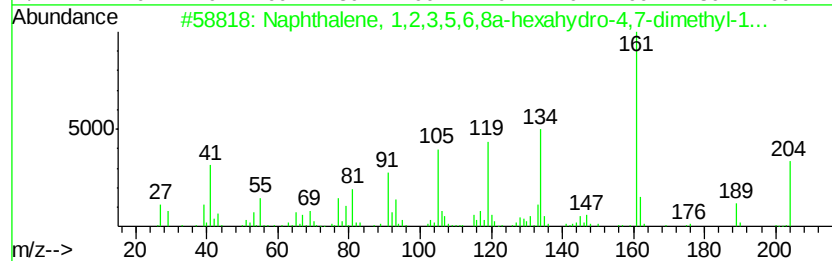
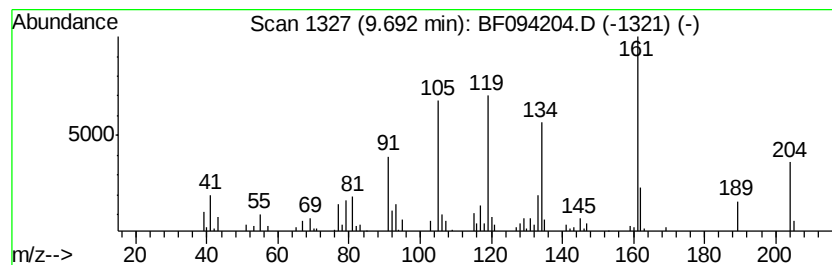
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Naphthalene, 1,2,3,5,6,8a-h... Concentration Rank 15

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.69 | 2.20 ng | 181092 | Acenaphthene-d10 | 9.50 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Naphthalene, 1,2,3,5,6,8a-hexahy... | 204 | C15H24  | 000483-76-1  | 98   |
| 2    |    | Naphthalene, 1,2,4a,5,8,8a-hexah... | 204 | C15H24  | 000523-47-7  | 93   |
| 3    |    | 1,4-Cyclohexadiene, 3-ethenyl-1,... | 134 | C10H14  | 062338-57-2  | 80   |
| 4    |    | Epizonarene                         | 204 | C15H24  | 1000156-10-7 | 74   |
| 5    |    | Copaene                             | 204 | C15H24  | 003856-25-5  | 70   |



Data Path : Z:\HPCHEM1\BNA\_F\Data\BF040517\  
Data File : BF094204.D  
Acq On : 5 Apr 2017 22:06  
Operator : SJ/MA  
Sample : I2522-17  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

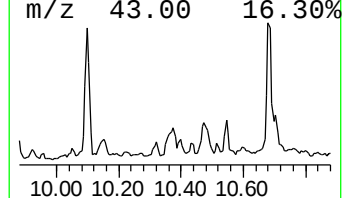
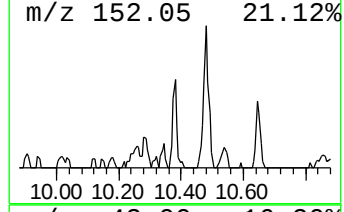
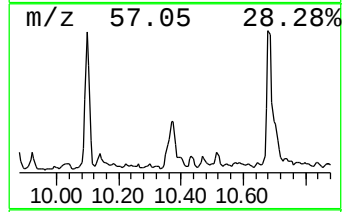
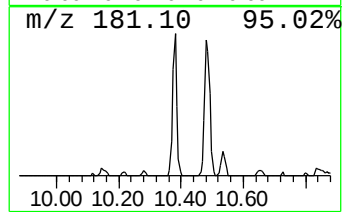
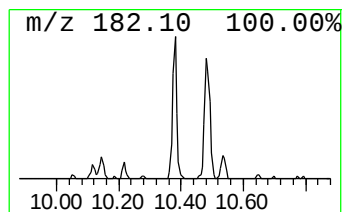
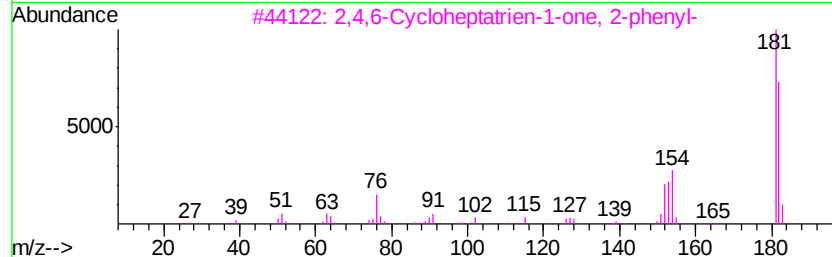
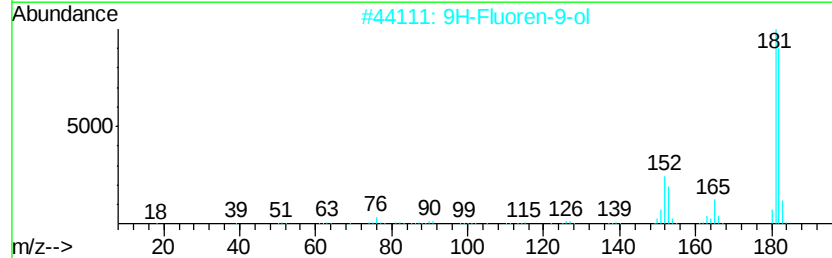
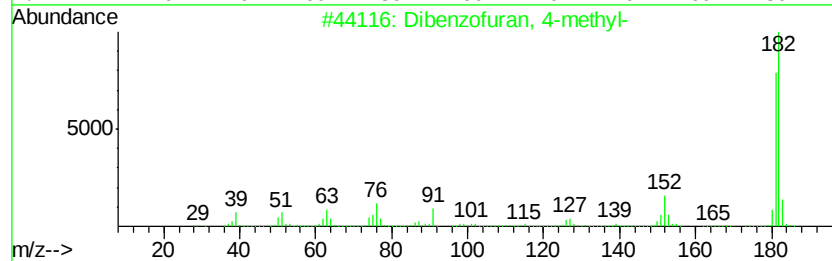
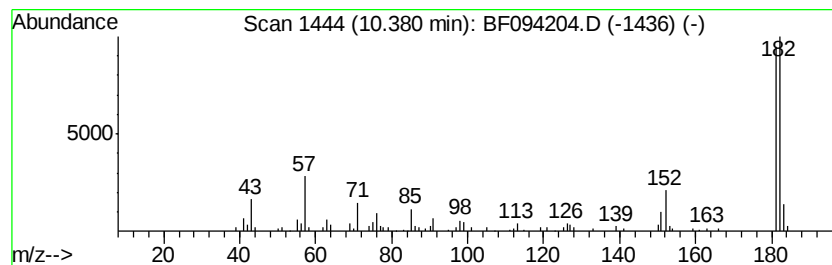
Instrument :  
BNA\_F  
ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 Dibenzofuran, 4-methyl- Concentration Rank 13

| R.T.    | EstConc | Area                                   | Relative to ISTD | R.T.           |
|---------|---------|----------------------------------------|------------------|----------------|
| 10.38   | 2.47 ng | 203291                                 | Acenaphthene-d10 | 9.50           |
| Hit# of | 5       | Tentative ID                           | MW MolForm       | CAS# Qual      |
| 1       |         | Dibenzofuran, 4-methyl-                | 182 C13H10O      | 007320-53-8 90 |
| 2       |         | 9H-Fluoren-9-ol                        | 182 C13H10O      | 001689-64-1 86 |
| 3       |         | 2,4,6-Cycloheptatrien-1-one, 2-phenyl- | 182 C13H10O      | 014562-09-5 78 |
| 4       |         | [1,1'-Biphenyl]-4-carboxaldehyde       | 182 C13H10O      | 003218-36-8 78 |
| 5       |         | 3-(2-Naphthyl)acrylaldehyde            | 182 C13H10O      | 020884-05-3 64 |



Data Path : Z:\HPCHEM1\BNA\_F\Data\BF040517\  
Data File : BF094204.D  
Acq On : 5 Apr 2017 22:06  
Operator : SJ/MA  
Sample : I2522-17  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

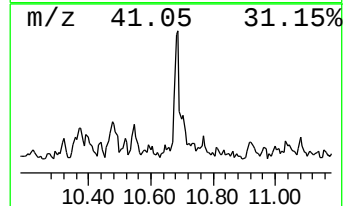
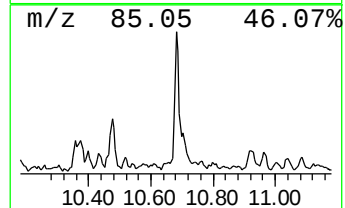
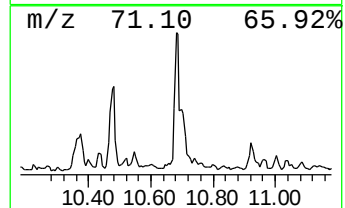
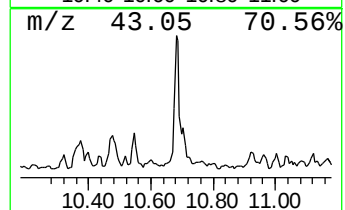
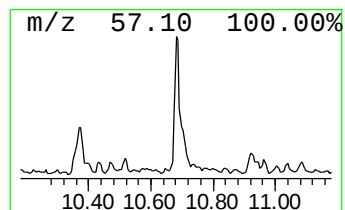
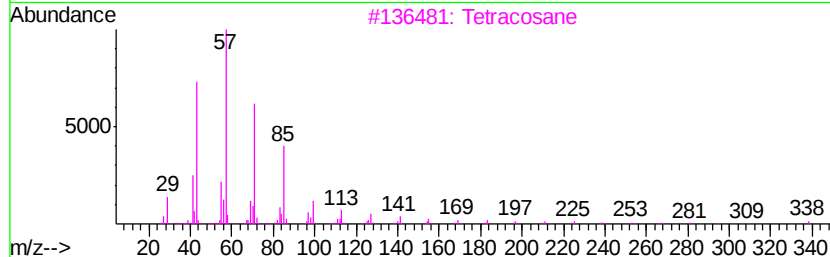
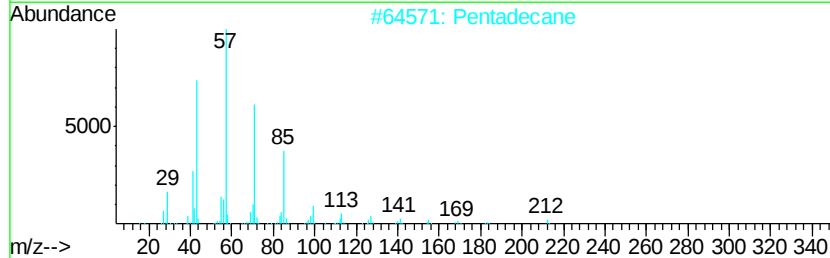
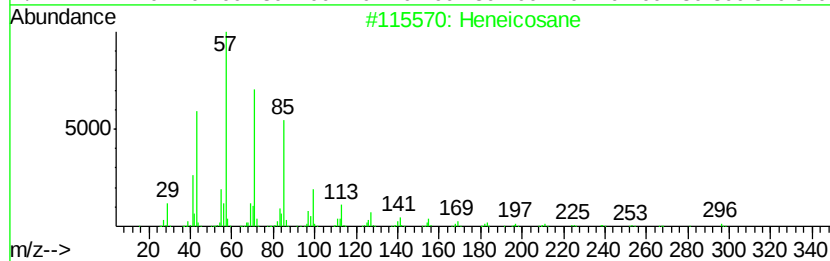
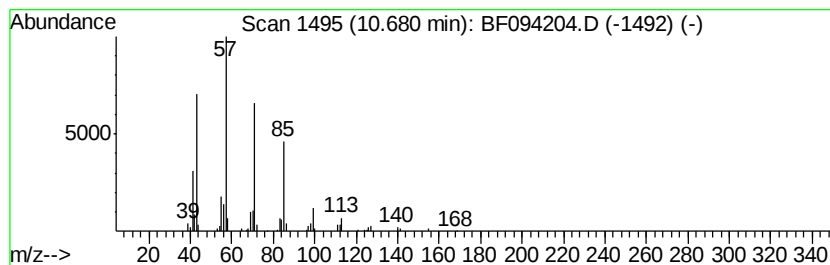
Instrument :  
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ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 Heneicosane Concentration Rank 12

| R.T.      | EstConc                  | Area   | Relative to ISTD | R.T.        |      |
|-----------|--------------------------|--------|------------------|-------------|------|
| 10.68     | 3.00 ng                  | 203708 | Phenanthrene-d10 | 11.33       |      |
| Hit# of 5 | Tentative ID             | MW     | MolForm          | CAS#        | Qual |
| 1         | Heneicosane              | 296    | C21H44           | 000629-94-7 | 90   |
| 2         | Pentadecane              | 212    | C15H32           | 000629-62-9 | 86   |
| 3         | Tetracosane              | 338    | C24H50           | 000646-31-1 | 86   |
| 4         | Decane, 2,3,5-trimethyl- | 184    | C13H28           | 062238-11-3 | 83   |
| 5         | Heptadecane, 8-methyl-   | 254    | C18H38           | 013287-23-5 | 83   |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\  
Data File : BF094204.D  
Acq On : 5 Apr 2017 22:06  
Operator : SJ/MA  
Sample : I2522-17  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

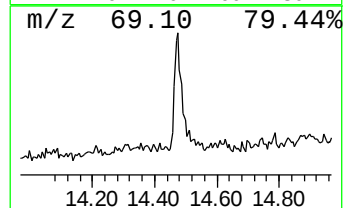
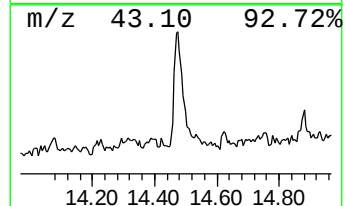
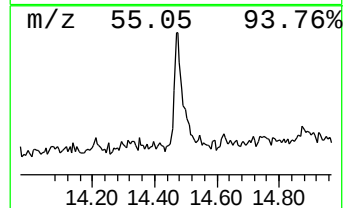
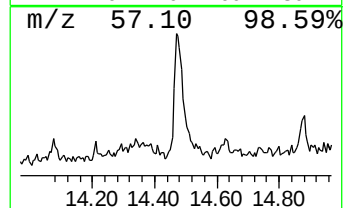
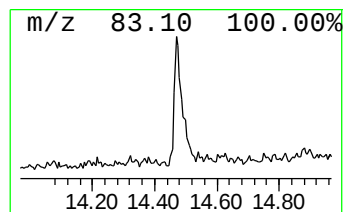
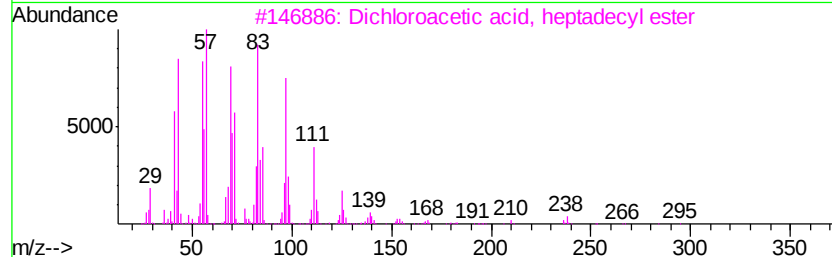
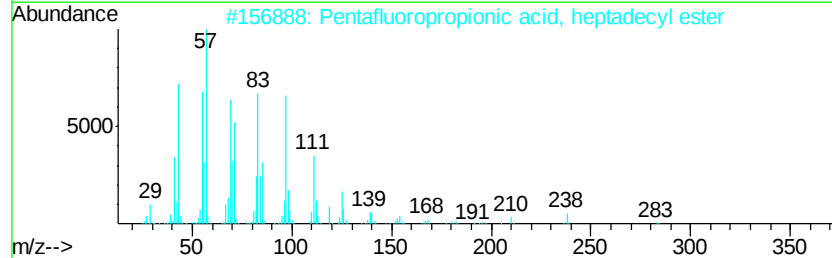
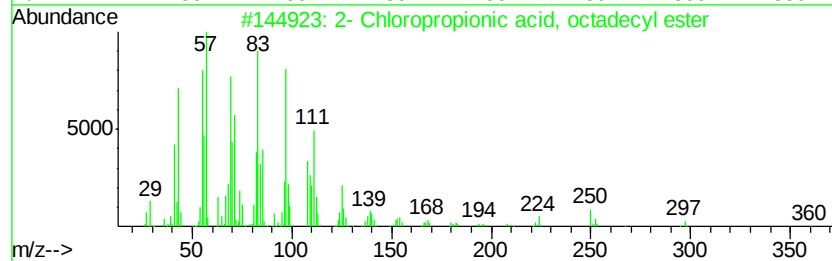
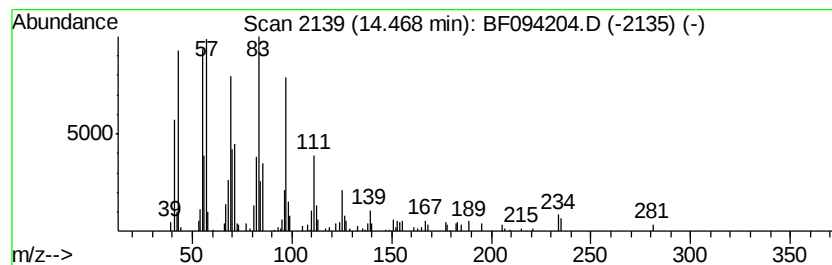
Instrument :  
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ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 2- Chloropropionic acid, oc... Concentration Rank 9

| R.T.    | EstConc                             | Area                             | Relative to ISTD | R.T.         |             |      |
|---------|-------------------------------------|----------------------------------|------------------|--------------|-------------|------|
| 14.47   | 4.20 ng                             | 224618                           | Chrysene-d12     | 14.57        |             |      |
| Hit# of | 5                                   | Tentative ID                     | MW               | MolForm      | CAS#        | Qual |
| 1       | 2-                                  | Chloropropionic acid, octadec... | 360              | C21H41ClO2   | 088104-31-8 | 94   |
| 2       | Pentafluoropropionic acid, hepta... | 402                              | C20H35F5O2       | 1000283-04-2 | 93          |      |
| 3       | Dichloroacetic acid, heptadecyl ... | 366                              | C19H36Cl2O2      | 1000282-98-2 | 93          |      |
| 4       | Cyclopentadecane                    | 210                              | C15H30           | 000295-48-7  | 92          |      |
| 5       | Heptafluorobutyric acid, n-octad... | 466                              | C22H37F7O2       | 000400-57-7  | 91          |      |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\  
Data File : BF094204.D  
Acq On : 5 Apr 2017 22:06  
Operator : SJ/MA  
Sample : I2522-17  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

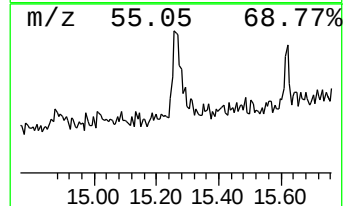
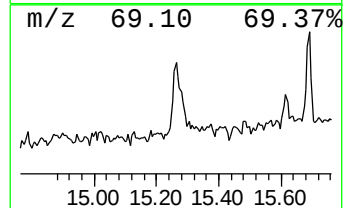
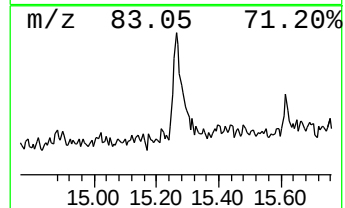
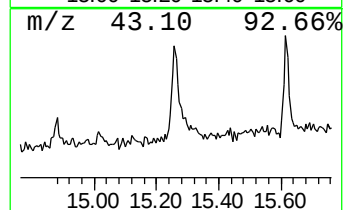
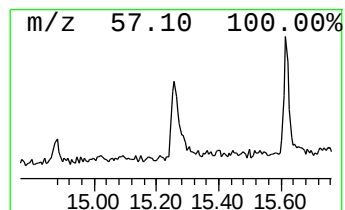
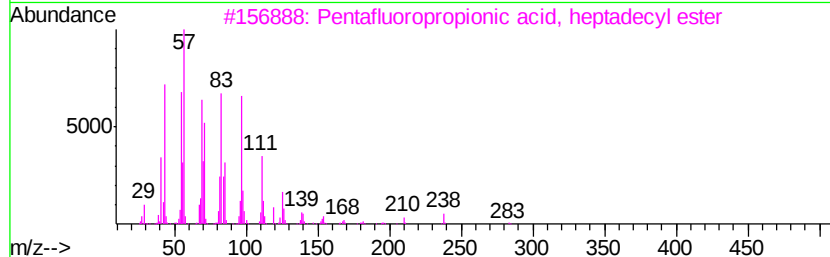
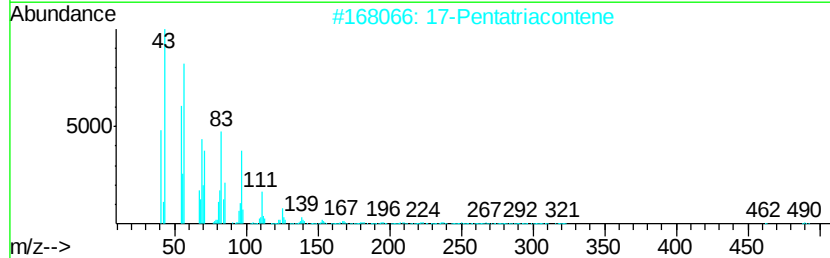
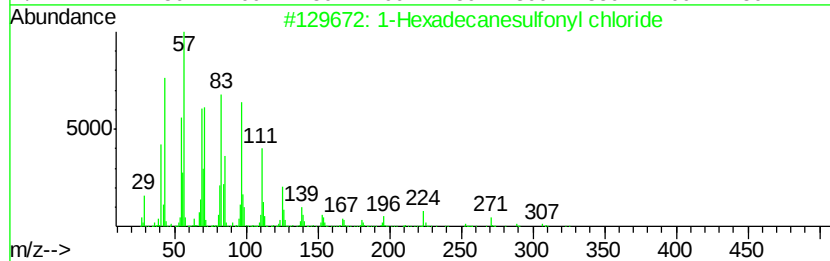
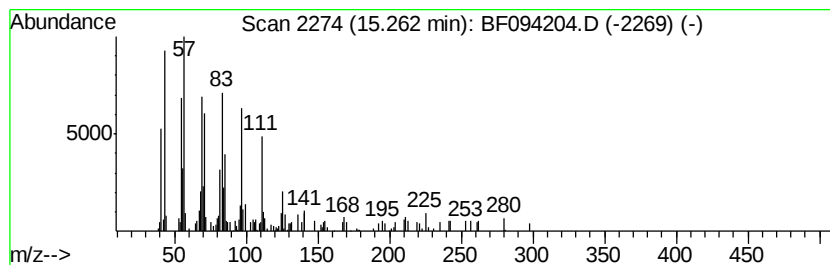
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ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 1-Hexadecanesulfonyl chloride Concentration Rank 11

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 15.26     | 3.24 ng                             | 173171 | Chrysene-d12     | 14.57        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1-Hexadecanesulfonyl chloride       | 324    | C16H33ClO2S      | 038775-38-1  | 58   |
| 2         | 17-Pentatriacontene                 | 491    | C35H70           | 006971-40-0  | 49   |
| 3         | Pentafluoropropionic acid, hepta... | 402    | C20H35F5O2       | 1000283-04-2 | 49   |
| 4         | 1-Decanol, 2-hexyl-                 | 242    | C16H34O          | 002425-77-6  | 46   |
| 5         | Heptafluorobutanoic acid, heptad... | 452    | C21H35F7O2       | 1000282-97-3 | 46   |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\  
Data File : BF094204.D  
Acq On : 5 Apr 2017 22:06  
Operator : SJ/MA  
Sample : I2522-17  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

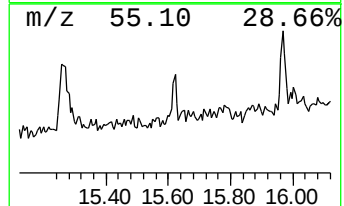
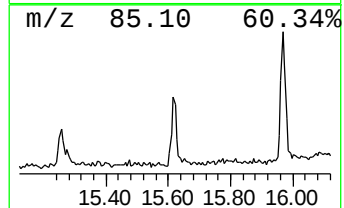
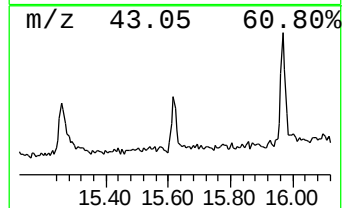
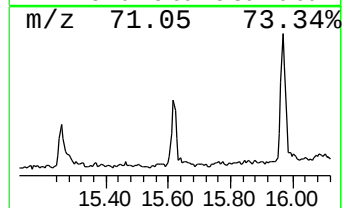
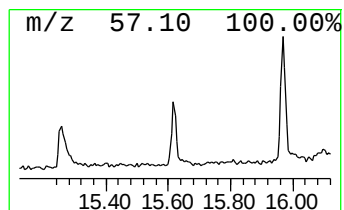
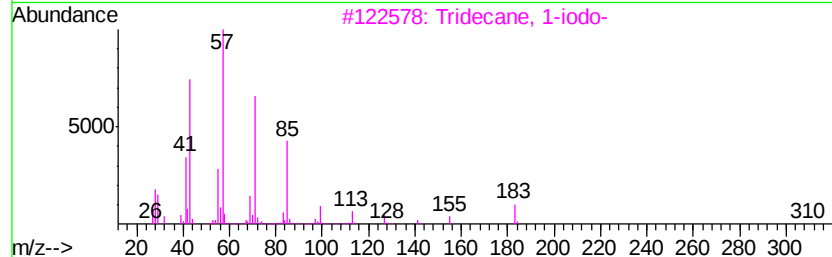
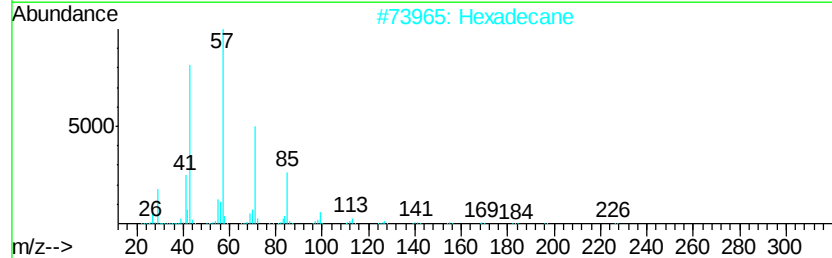
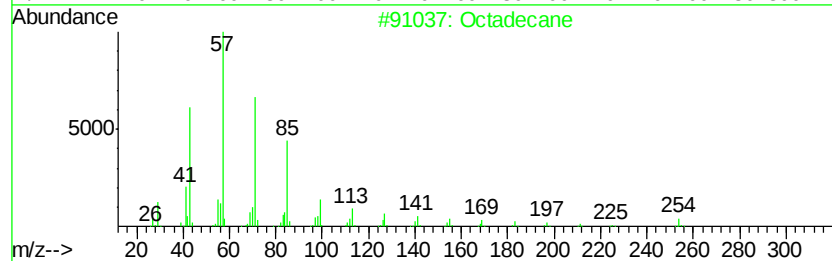
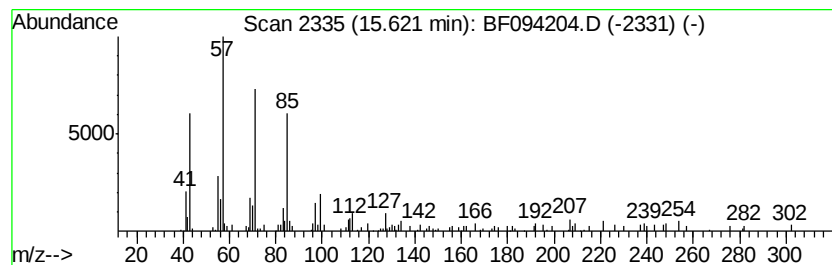
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BNA\_F  
ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 Octadecane Concentration Rank 17

| R.T.    | EstConc            | Area         | Relative to ISTD | R.T.      |
|---------|--------------------|--------------|------------------|-----------|
| 15.62   | 2.03 ng            | 90920        | Perylene-d12     | 16.20     |
| Hit# of | 5                  | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Octadecane         | 254 C18H38   | 000593-45-3      | 87        |
| 2       | Hexadecane         | 226 C16H34   | 000544-76-3      | 83        |
| 3       | Tridecane, 1-iodo- | 310 C13H27I  | 035599-77-0      | 80        |
| 4       | Dodecane, 1-iodo-  | 296 C12H25I  | 004292-19-7      | 80        |
| 5       | Heptadecane        | 240 C17H36   | 000629-78-7      | 80        |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\  
Data File : BF094204.D  
Acq On : 5 Apr 2017 22:06  
Operator : SJ/MA  
Sample : I2522-17  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

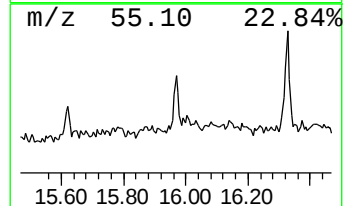
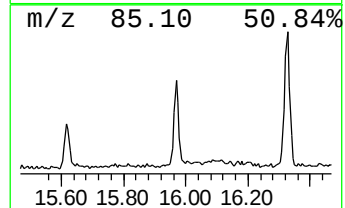
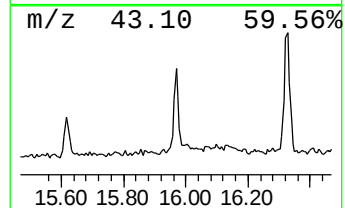
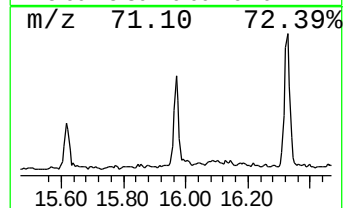
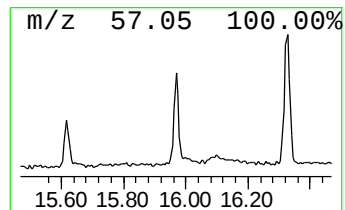
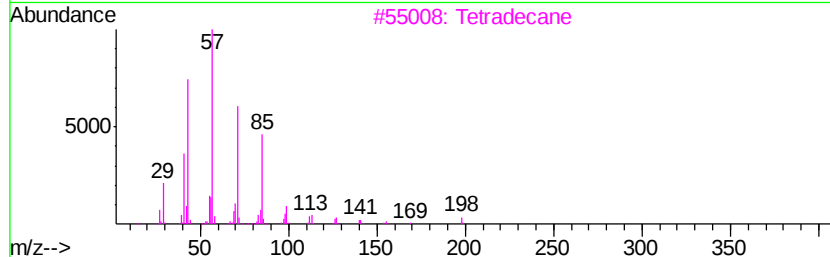
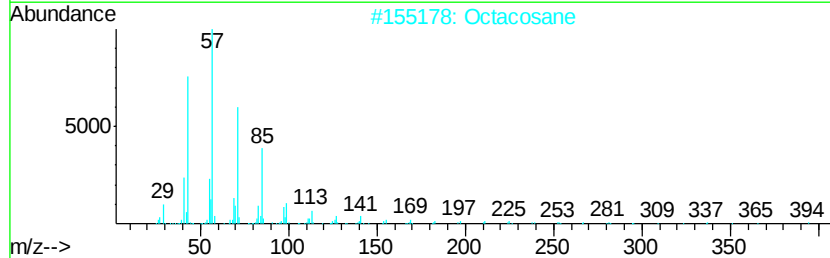
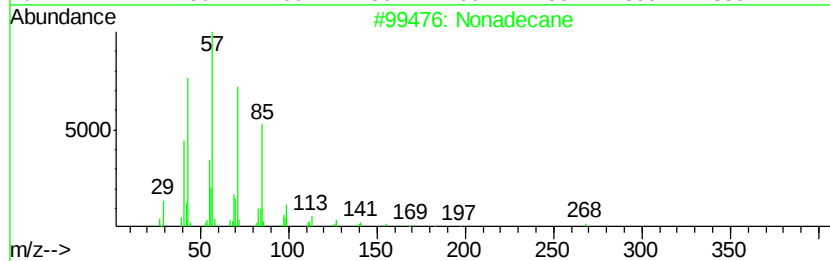
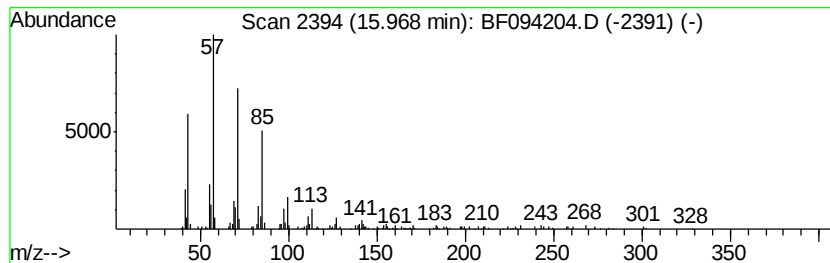
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Nonadecane Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.97 | 3.77 ng | 169046 | Perylene-d12     | 16.20 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 94   |
| 2    |    |   | Octacosane   | 394 | C28H58  | 000630-02-4 | 90   |
| 3    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 87   |
| 4    |    |   | Tetracosane  | 338 | C24H50  | 000646-31-1 | 87   |
| 5    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 87   |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\  
Data File : BF094204.D  
Acq On : 5 Apr 2017 22:06  
Operator : SJ/MA  
Sample : I2522-17  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

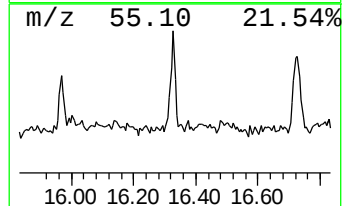
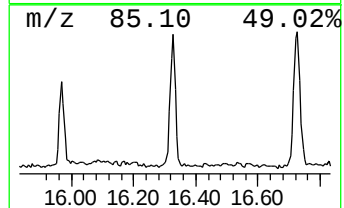
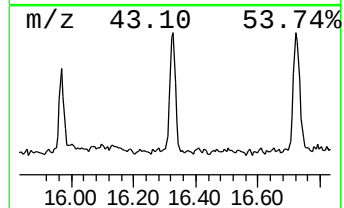
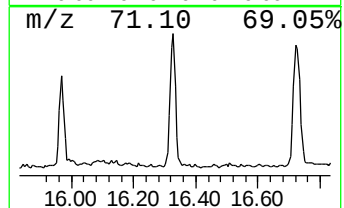
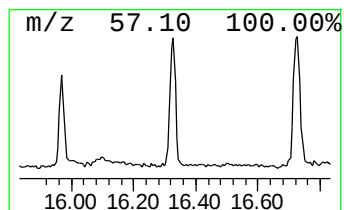
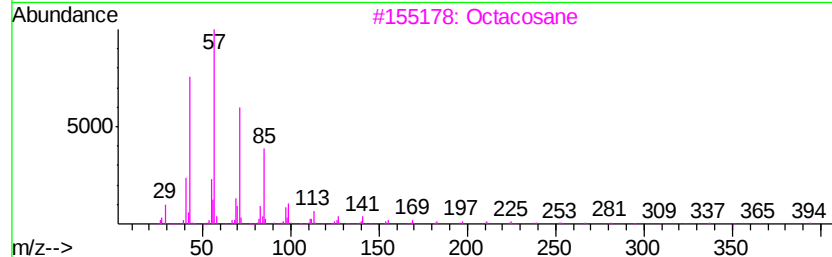
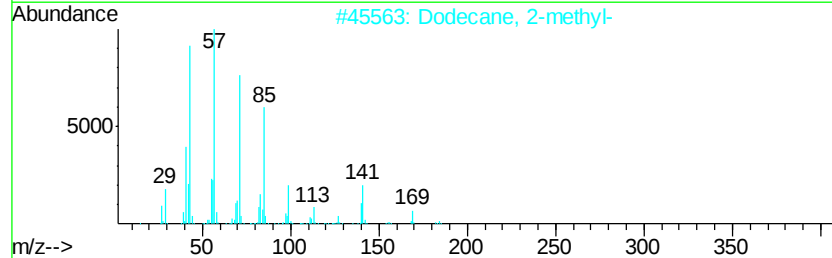
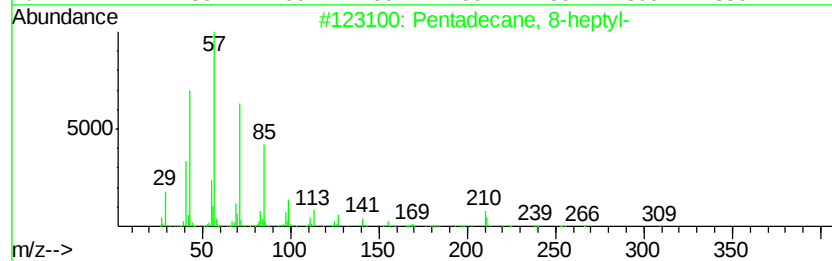
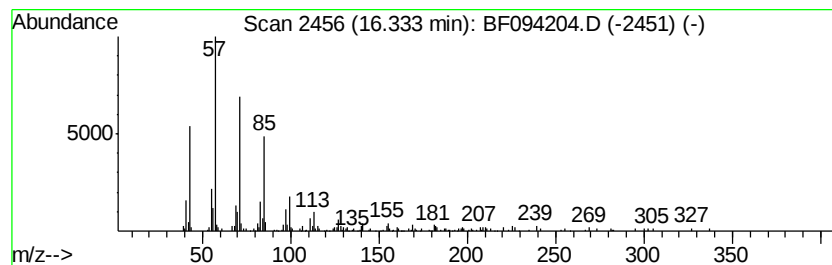
Instrument :  
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ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 Pentadecane, 8-heptyl- Concentration Rank 6

| R.T.      | EstConc                | Area   | Relative to ISTD |             | R.T.  |
|-----------|------------------------|--------|------------------|-------------|-------|
| 16.33     | 5.75 ng                | 257898 | Perylene-d12     |             | 16.20 |
| Hit# of 5 | Tentative ID           | MW     | MolForm          | CAS#        | Qual  |
| 1         | Pentadecane, 8-heptyl- | 310    | C22H46           | 071005-15-7 | 91    |
| 2         | Dodecane, 2-methyl-    | 184    | C13H28           | 001560-97-0 | 90    |
| 3         | Octacosane             | 394    | C28H58           | 000630-02-4 | 87    |
| 4         | Octadecane, 1-iodo-    | 380    | C18H37I          | 000629-93-6 | 87    |
| 5         | Heptadecane, 9-octyl-  | 352    | C25H52           | 007225-64-1 | 87    |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\  
Data File : BF094204.D  
Acq On : 5 Apr 2017 22:06  
Operator : SJ/MA  
Sample : I2522-17  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

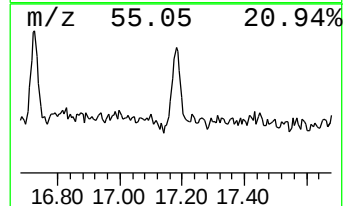
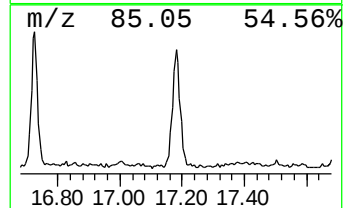
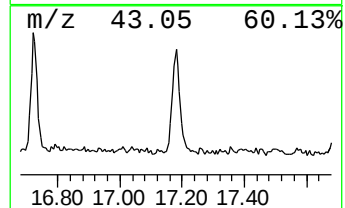
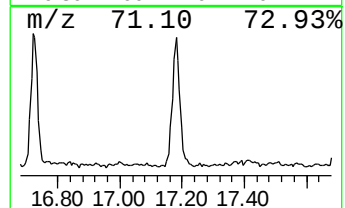
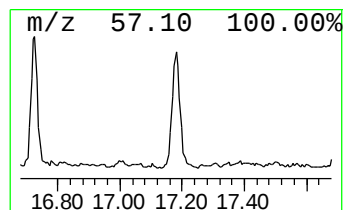
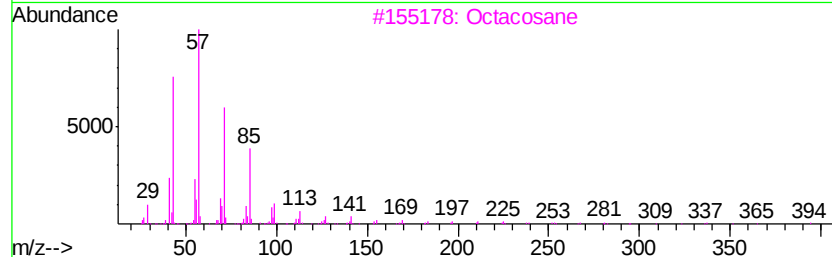
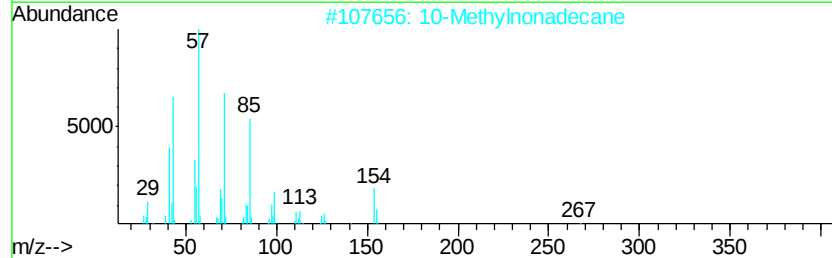
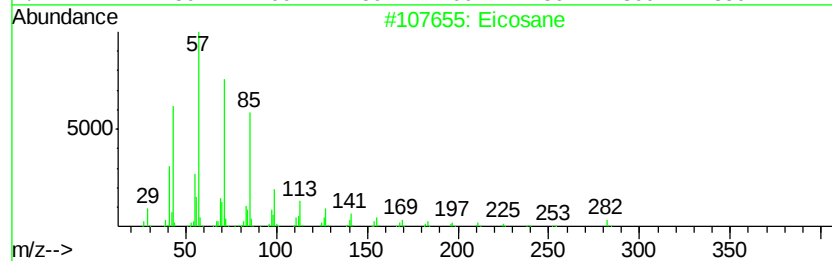
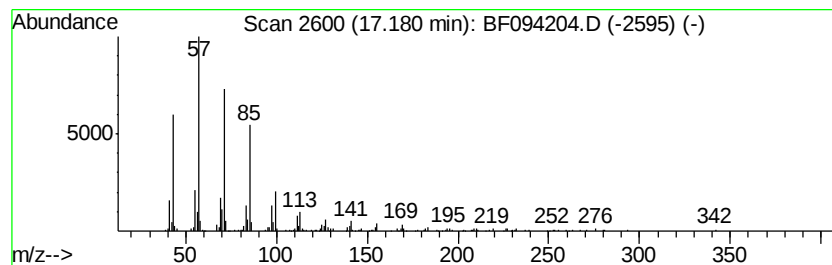
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Eicosane Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.18 | 7.32 ng | 328421 | Perylene-d12     | 16.20 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane              | 282 | C20H42  | 000112-95-8 | 86   |
| 2    |    | 10-Methylnonadecane   | 282 | C20H42  | 056862-62-5 | 86   |
| 3    |    | Octacosane            | 394 | C28H58  | 000630-02-4 | 83   |
| 4    |    | Decane, 3,8-dimethyl- | 170 | C12H26  | 017312-55-9 | 81   |
| 5    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 80   |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\  
Data File : BF094204.D  
Acq On : 5 Apr 2017 22:06  
Operator : SJ/MA  
Sample : I2522-17  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

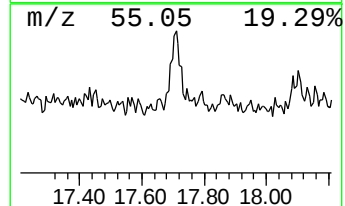
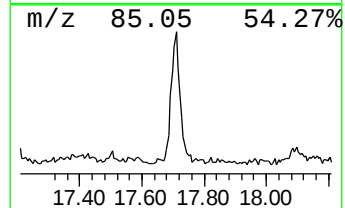
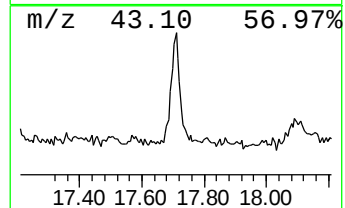
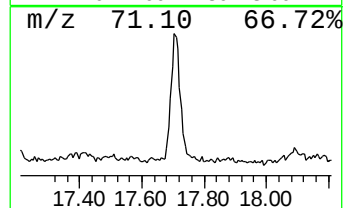
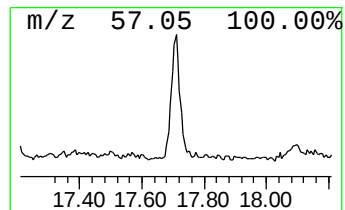
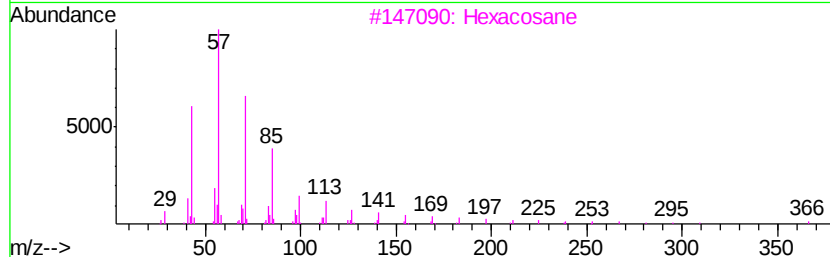
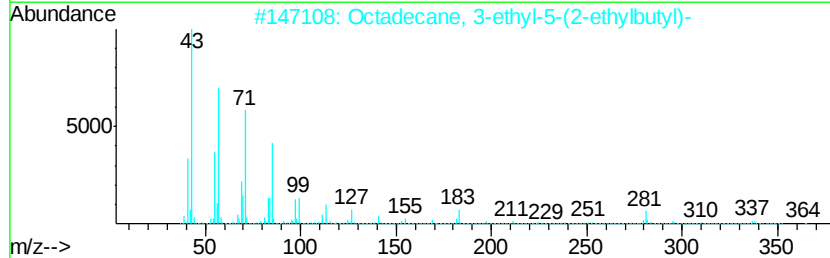
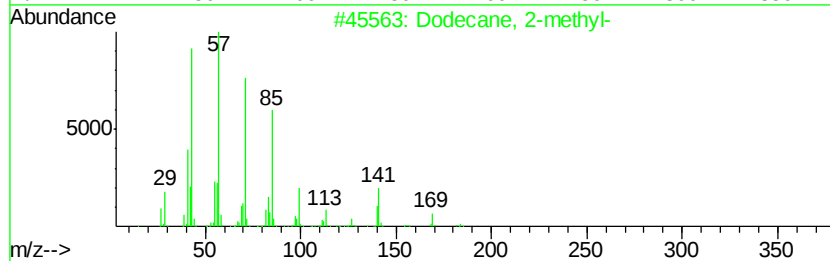
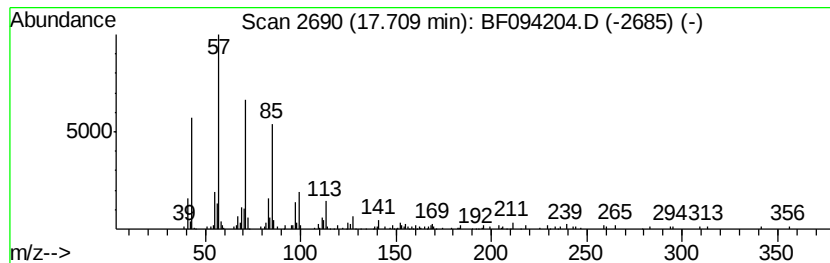
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Dodecane, 2-methyl- Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.71 | 5.48 ng | 245715 | Perylene-d12     | 16.20 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2-methyl-                 | 184 | C13H28  | 001560-97-0 | 90   |
| 2    |    |   | Octadecane, 3-ethyl-5-(2-ethylbu... | 366 | C26H54  | 055282-12-7 | 90   |
| 3    |    |   | Hexacosane                          | 366 | C26H54  | 000630-01-3 | 90   |
| 4    |    |   | Decane, 3,8-dimethyl-               | 170 | C12H26  | 017312-55-9 | 87   |
| 5    |    |   | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 87   |



Data Path : Z:\HPCHEM1\BNA\_F\Data\BF040517\  
 Data File : BF094204.D  
 Acq On : 5 Apr 2017 22:06  
 Operator : SJ/MA  
 Sample : I2522-17  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 BASING-9(0-2)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032217.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2-Pentanone, 4-hy... | 4.01  | 11.0    | ng    | 601390   | 1                     | 5.86  | 1089850 | 20.0 |
| unknown5.61          | 5.61  | 76.8    | ng    | 4183170  | 1                     | 5.86  | 1089850 | 20.0 |
| Naphthalene, 1,3-... | 9.10  | 2.1     | ng    | 175829   | 3                     | 9.50  | 1647000 | 20.0 |
| Naphthalene, 1,2,... | 9.57  | 2.4     | ng    | 196449   | 3                     | 9.50  | 1647000 | 20.0 |
| Naphthalene, 1,2,... | 9.69  | 2.2     | ng    | 181092   | 3                     | 9.50  | 1647000 | 20.0 |
| Dibenzofuran, 4-m... | 10.38 | 2.5     | ng    | 203291   | 3                     | 9.50  | 1647000 | 20.0 |
| Heneicosane          | 10.68 | 3.0     | ng    | 203708   | 4                     | 11.33 | 1357860 | 20.0 |
| 2-Chloropropioni...  | 14.47 | 4.2     | ng    | 224618   | 5                     | 14.57 | 1068900 | 20.0 |
| 1-Hexadecanesulfo... | 15.26 | 3.2     | ng    | 173171   | 5                     | 14.57 | 1068900 | 20.0 |
| Octadecane           | 15.62 | 2.0     | ng    | 90920    | 6                     | 16.20 | 897342  | 20.0 |
| Nonadecane           | 15.97 | 3.8     | ng    | 169046   | 6                     | 16.20 | 897342  | 20.0 |
| Pentadecane, 8-he... | 16.33 | 5.8     | ng    | 257898   | 6                     | 16.20 | 897342  | 20.0 |
| Eicosane             | 17.18 | 7.3     | ng    | 328421   | 6                     | 16.20 | 897342  | 20.0 |
| Dodecane, 2-methyl-  | 17.71 | 5.5     | ng    | 245715   | 6                     | 16.20 | 897342  | 20.0 |