LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

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-BF090115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 1.224 | 10 | 12 | 15 | rVB | 850179 | 1159919 | 26.48% | 6.276% |
| 2 | 1.304 | 17 | 19 | 29 | rVB3 | 23004 | 55557 | 1.27% | 0.301% |
| 3 | 1.430 | 29 | 30 | 36 | rBV | 218667 | 572854 | 13.08% | 3.100% |
| 4 | 1.521 | 36 | 38 | 78 | rVB | 1219172 | 4380783 | 100.00% | 23.704% |
| 5 | 4.424 | 288 | 292 | 314 | rVV | 242372 | 728864 | 16.64% | 3.944% |
| 6 | 5.304 | 366 | 369 | 391 | rBV | 595641 | 1254104 | 28.63% | 6.786% |
| 7 | 7.579 | 565 | 568 | 572 | rBV | 633459 | 1275296 | 29.11% | 6.901% |
| 8 | 7.659 | 572 | 575 | 589 | rVB | 778311 | 1399339 | 31.94% | 7.572% |
| 9 | 8.139 | 614 | 617 | 623 | rVB | 134366 | 239862 | 5.48% | 1.298% |
| 10 | 8.459 | 641 | 645 | 655 | rBB | 564546 | 947937 | 21.64% | 5.129% |
| 11 | 9.442 | 727 | 731 | 744 | rBV | 461483 | 854207 | 19.50% | 4.622% |
| 12 | 11.031 | 867 | 870 | 873 | rBV | 184567 | 298060 | 6.80% | 1.613% |
| 13 | 12.311 | 977 | 982 | 988 | rBV2 | 19801 | 47451 | 1.08% | 0.257% |
| 14 | 13.682 | 1097 | 1102 | 1105 | rBV | 789753 | 1356032 | 30.95% | 7.337% |
| 15 | 13.740 | 1105 | 1107 | 1118 | rVB | 36645 | 87071 | 1.99% | 0.471% |
| 16 | 14.734 | 1191 | 1194 | 1199 | rVV | 151054 | 240999 | 5.50% | 1.304% |
| 17 | 14.848 | 1200 | 1204 | 1207 | rVV2 | 39087 | 72769 | 1.66% | 0.394% |
| 18 | 15.157 | 1228 | 1231 | 1237 | rBV | 168728 | 302296 | 6.90% | 1.636% |
| 19 | 17.066 | 1394 | 1398 | 1404 | rBV2 | 419740 | 849074 | 19.38% | 4.594% |
| 20 | 17.728 | 1450 | 1456 | 1462 | rBV | 46298 | 123856 | 2.83% | 0.670% |
| 21 | 18.666 | 1535 | 1538 | 1542 | rBV | 171946 | 283609 | 6.47% | 1.535% |
| 22 | 18.849 | 1551 | 1554 | 1563 | rVB | 19731 | 50295 | 1.15% | 0.272% |
| 23 | 19.626 | 1616 | 1622 | 1626 | rBV | 43452 | 77367 | 1.77% | 0.419% |
| 24 | 20.437 | 1689 | 1693 | 1699 | rBV | 44326 | 101600 | 2.32% | 0.550% |
| 25 | 21.386 | 1773 | 1776 | 1782 | rBV2 | 119801 | 177942 | 4.06% | 0.963% |
| 26 | 22.049 | 1831 | 1834 | 1837 | rBV | 878087 | 1006708 | 22.98% | 5.447% |
| 27 | 23.295 | 1940 | 1943 | 1944 | rBV | 99839 | 114282 | 2.61% | 0.618% |
| 28 | 23.329 | 1944 | 1946 | 1952 | rVB | 210997 | 245103 | 5.59% | 1.326% |
| 29 | 24.769 | 2070 | 2072 | 2075 | rBV | 174021 | 177694 | 4.06% | 0.961% |

Sum of corrected areas: 18480930

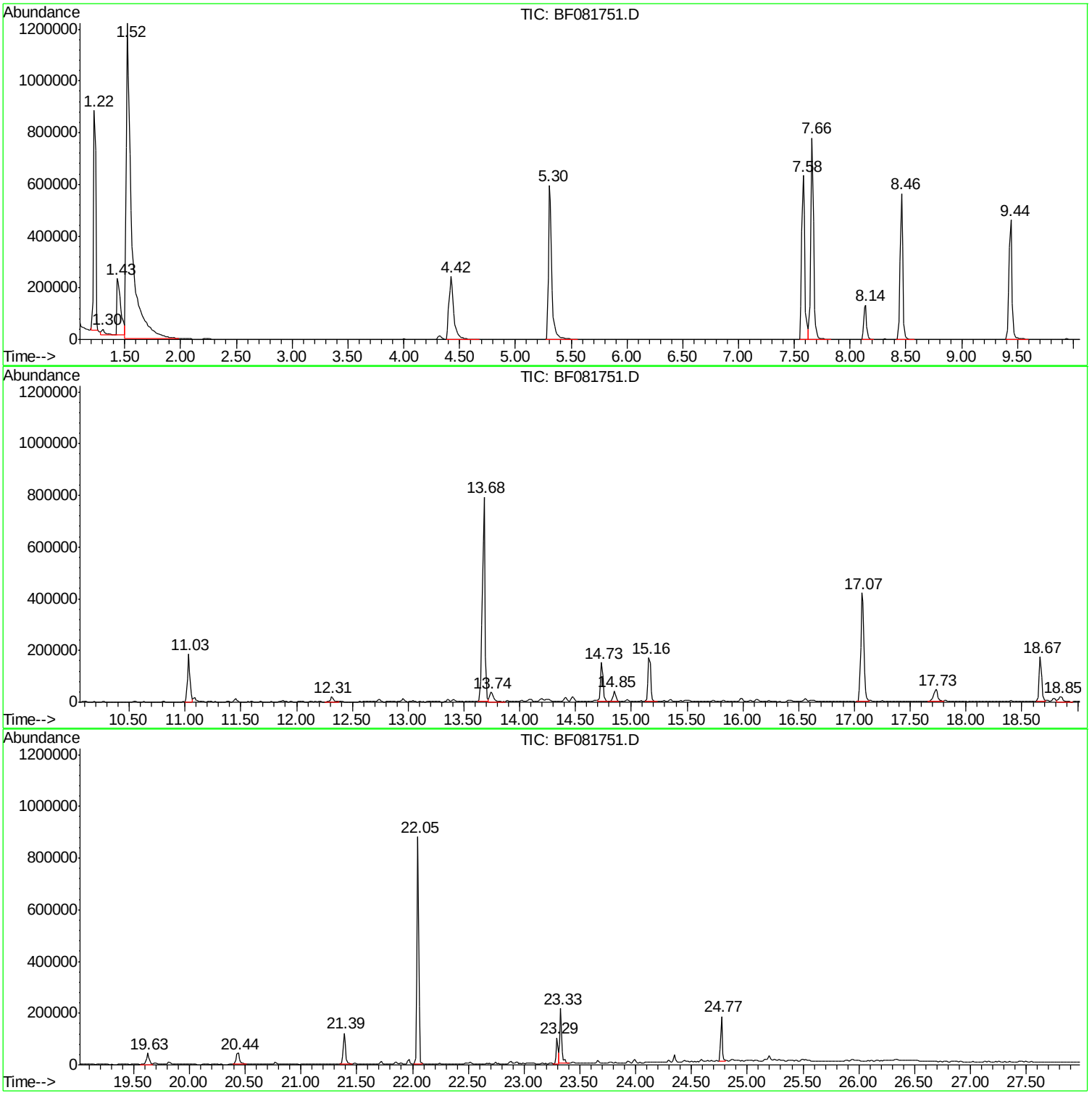
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

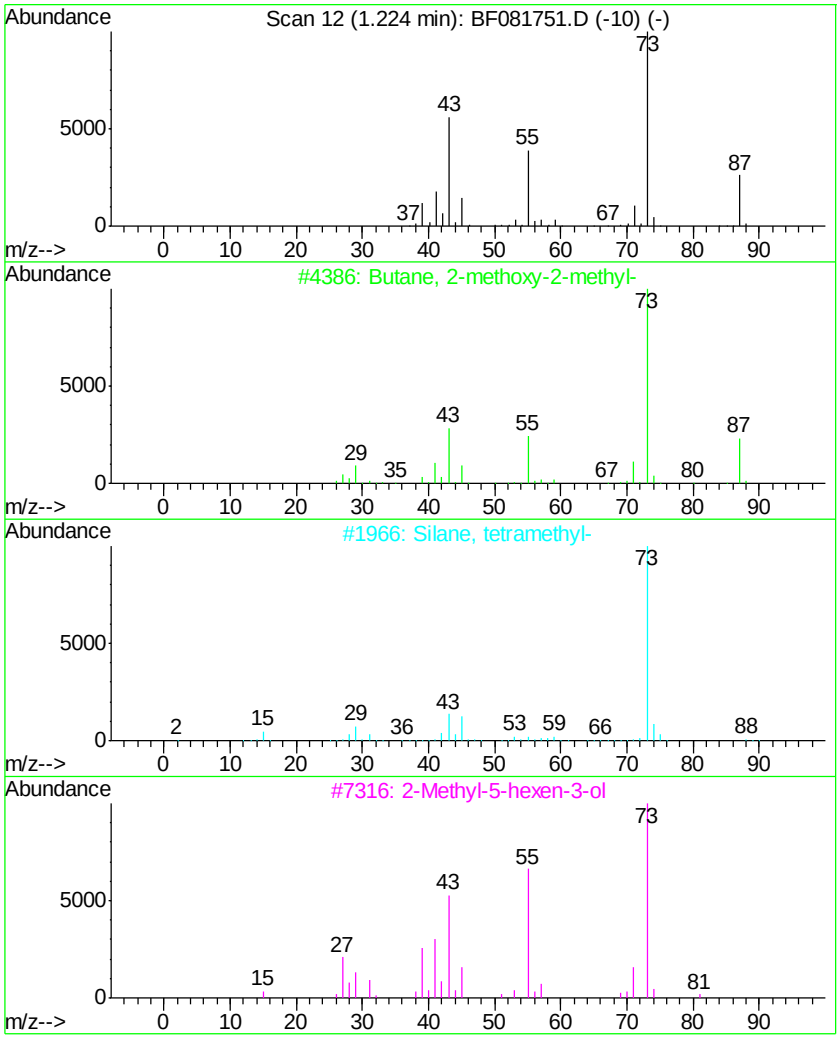
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

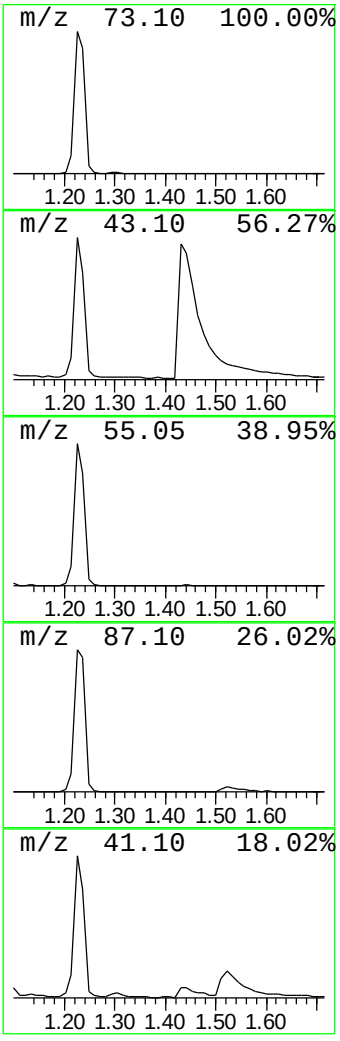
**********

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 1.22 | 96.72 ng | 1159920 | 1,4-Dichlorobenzene-d4 | 8.14 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2-methoxy-2-methyl- | 102 | C6H14O | 000994-05-8 | 83 |
| 2 | Silane, tetramethyl- | 88 | C4H12Si | 000075-76-3 | 25 |
| 3 | 2-Methyl-5-hexen-3-ol | 114 | C7H14O | 032815-70-6 | 12 |
| 4 | Methane, isothiocyanato- | 73 | C2H3NS | 000556-61-6 | 9 |
| 5 | Acetamide, N-ethyl- | 87 | C4H9NO | 000625-50-3 | 9 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

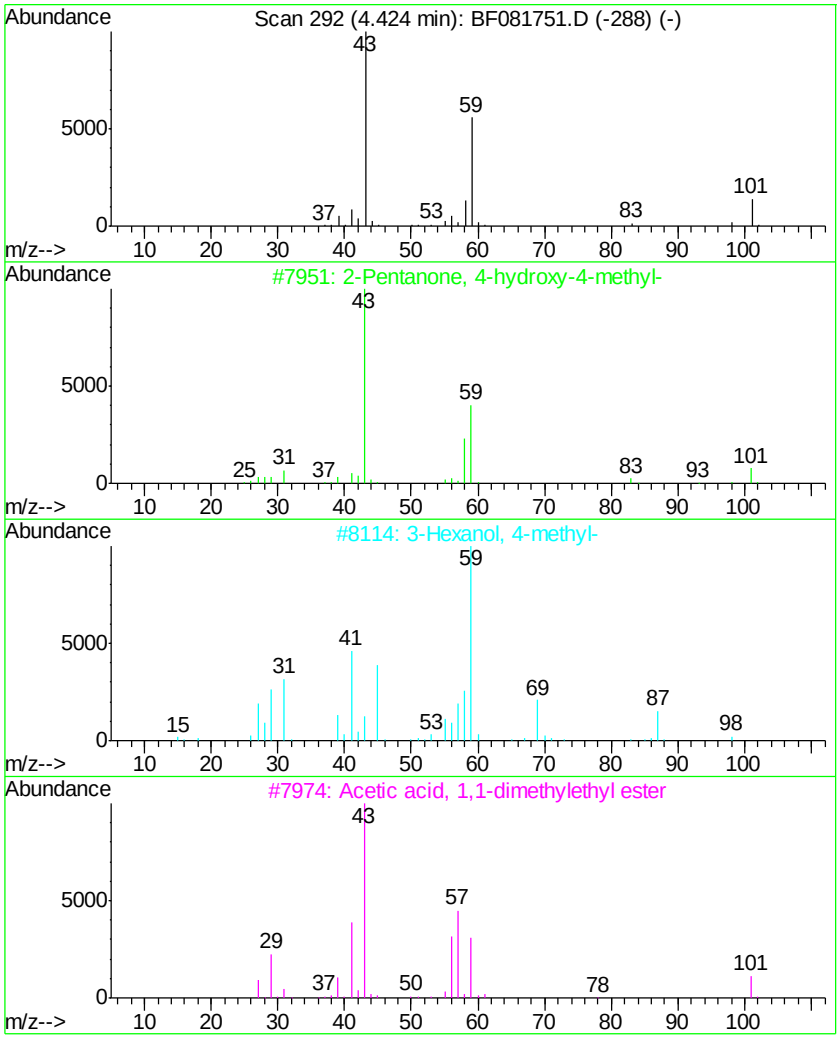
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

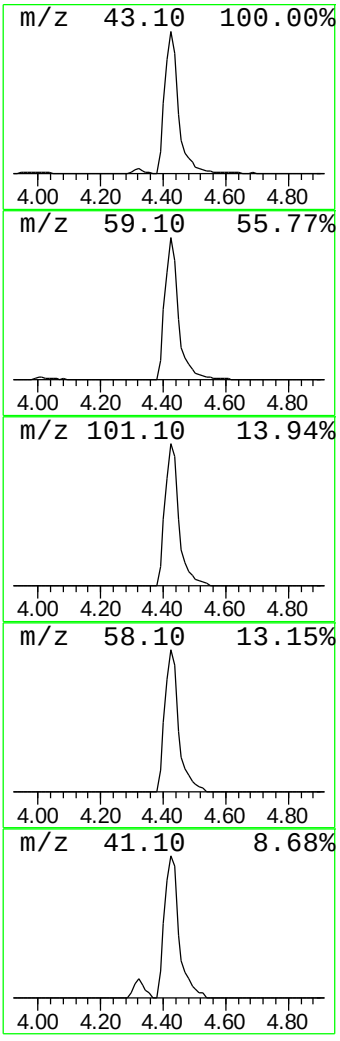
**********

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.42 | 60.77 ng | 728864 | 1,4-Dichlorobenzene-d4 | 8.14 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 50 |
| 2 | 3-Hexanol, 4-methyl- | 116 | C7H16O | 000615-29-2 | 33 |
| 3 | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5 | 28 |
| 4 | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 25 |
| 5 | 1-Propen-2-ol, acetate | 100 | C5H8O2 | 000108-22-5 | 10 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

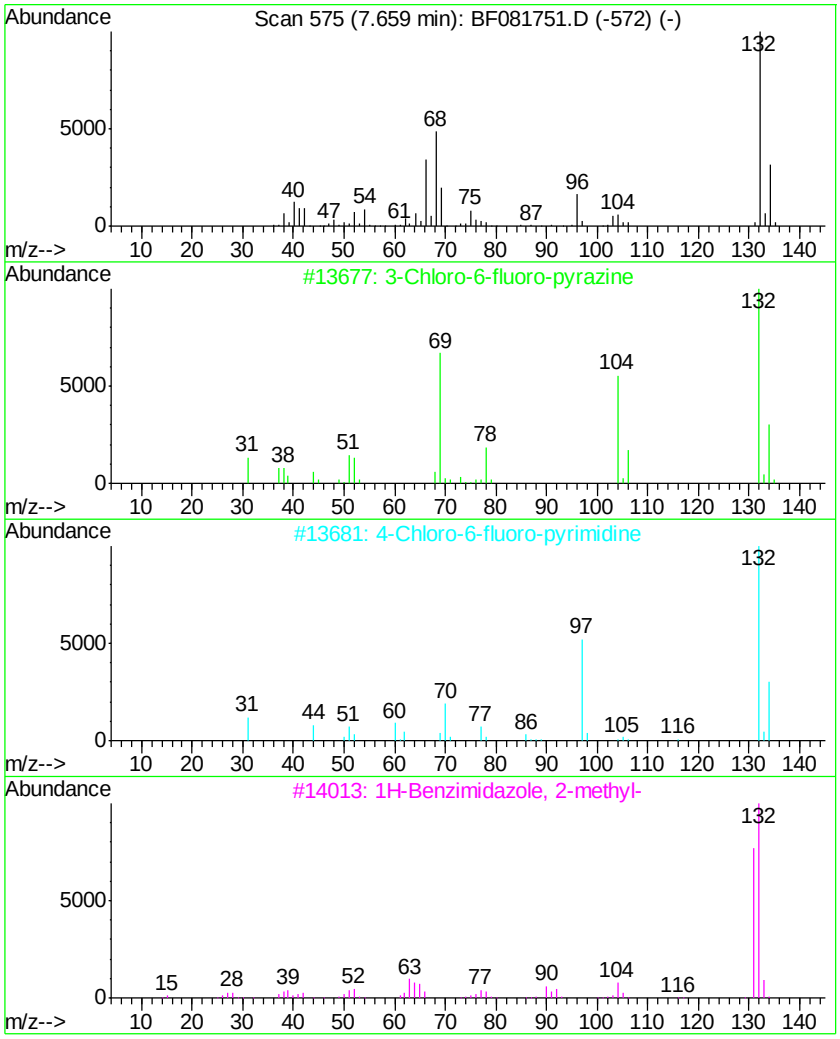
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

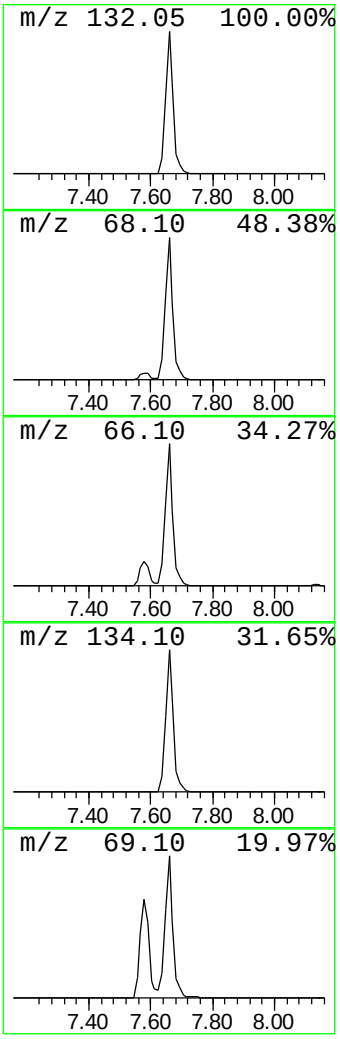
**********

Peak Number 6 unknown7.66 Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.66 | 116.68 ng | 1399340 | 1,4-Dichlorobenzene-d4 | 8.14 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Chloro-6-fluoro-pyrazine | 132 | C4H2ClFN2 | 1000146-10-7 | 27 |
| 2 | 4-Chloro-6-fluoro-pyrimidine | 132 | C4H2ClFN2 | 051422-01-6 | 27 |
| 3 | 1H-Benzimidazole, 2-methyl- | 132 | C8H8N2 | 000615-15-6 | 22 |
| 4 | 5-Fluoro-2-chloropyrimidine | 132 | C4H2ClFN2 | 062802-42-0 | 12 |
| 5 | (5-METHYL-2-PYRIDYL)ACETONITRILE | 132 | C8H8N2 | 1000241-93-9 | 10 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

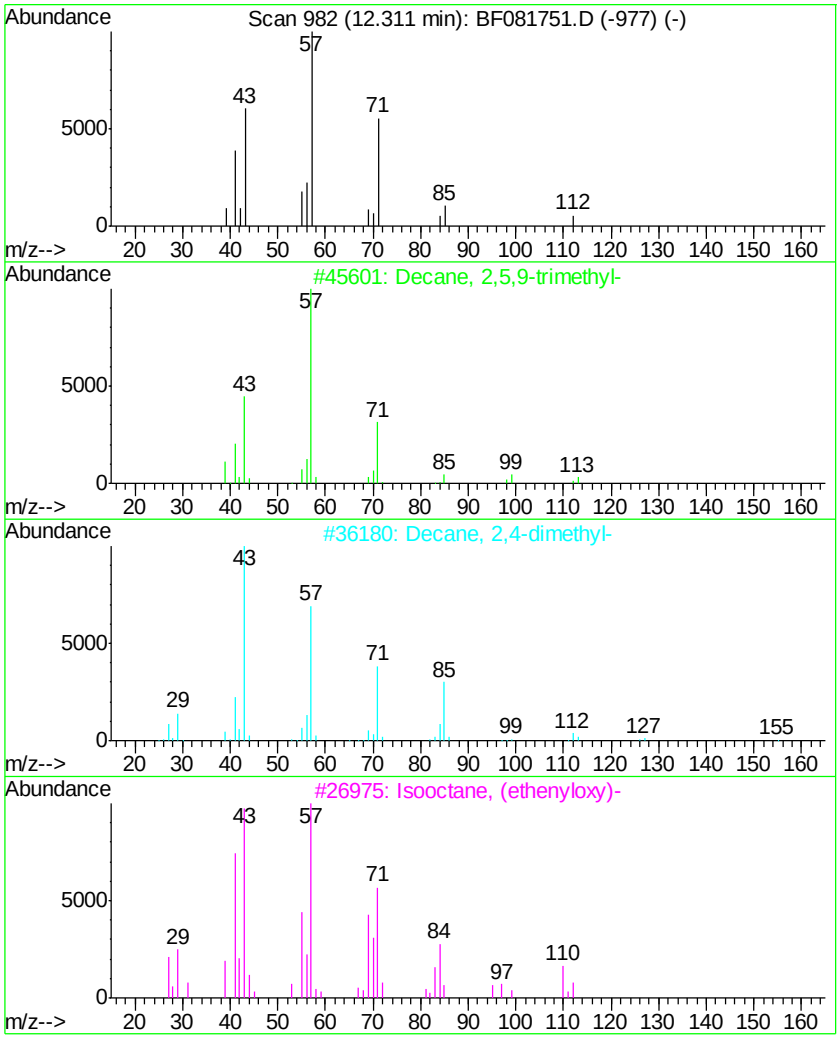
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

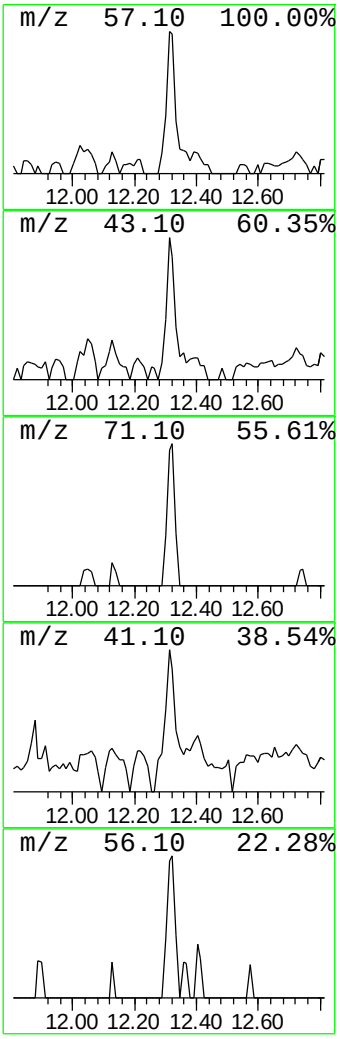
**********

Peak Number 8 Decane, 2,5,9-trimethyl- Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.31 | 3.18 ng | 47451 | Naphthalene-d8 | 11.03 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Decane, 2,5,9-trimethyl- | 184 | C13H28 | 062108-22-9 | 59 |
| 2 | Decane, 2,4-dimethyl- | 170 | C12H26 | 002801-84-5 | 39 |
| 3 | Isooctane, (ethenyloxy)- | 156 | C10H20O | 037769-62-3 | 39 |
| 4 | Hexadecane | 226 | C16H34 | 000544-76-3 | 36 |
| 5 | Trifluoroacetic acid, 4-methylpe... | 198 | C8H13F3O2 | 000327-71-9 | 36 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

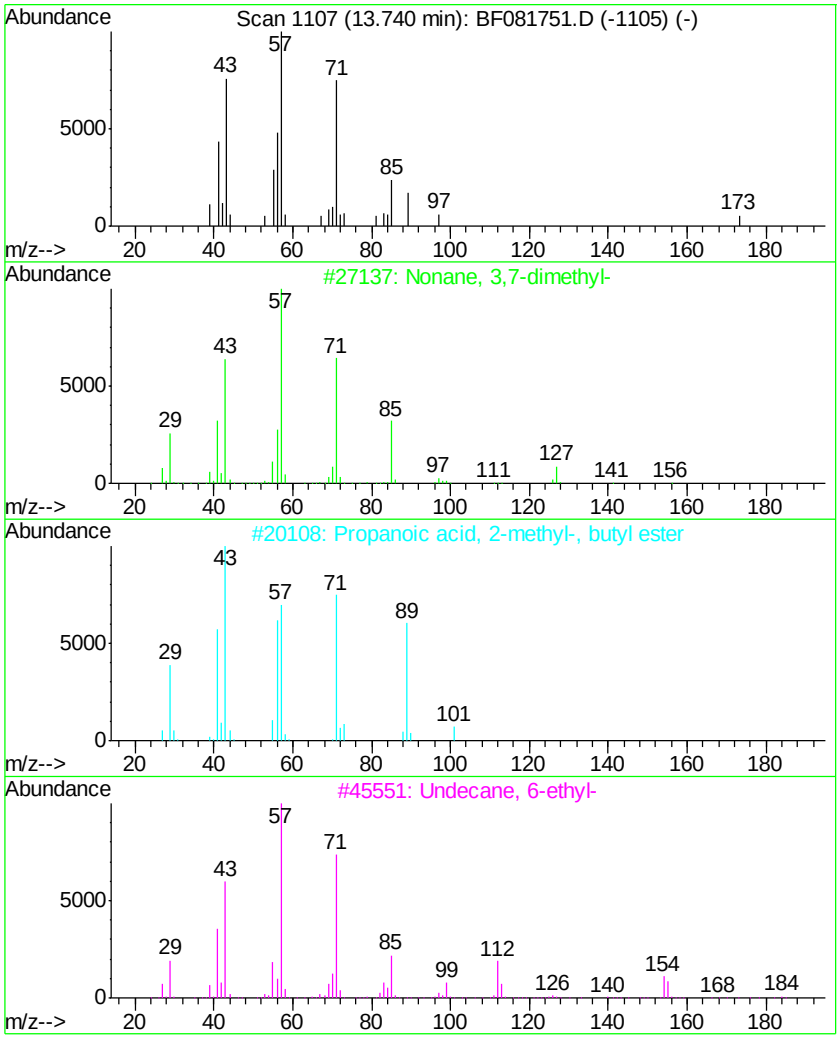
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

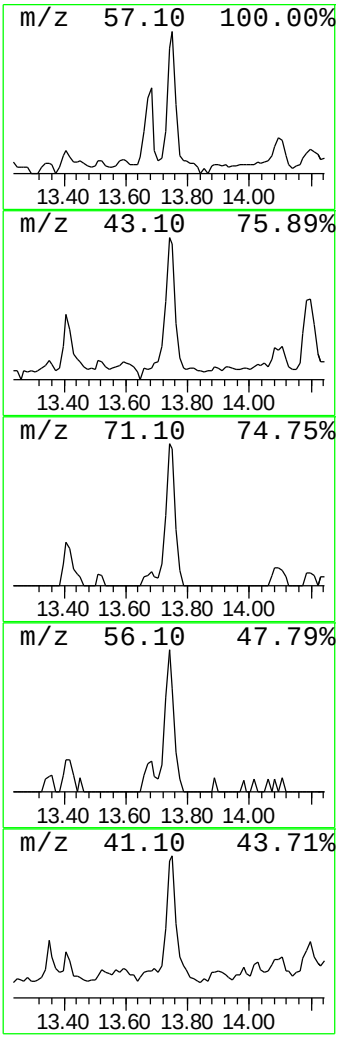
**********

Peak Number 9 Nonane, 3,7-dimethyl- Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.74 | 5.76 ng | 87071 | Acenaphthene-d10 | 15.16 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Nonane, 3,7-dimethyl- | 156 | C11H24 | 017302-32-8 | 59 |
| 2 | Propanoic acid, 2-methyl-, butyl... | 144 | C8H16O2 | 000097-87-0 | 59 |
| 3 | Undecane, 6-ethyl- | 184 | C13H28 | 017312-60-6 | 53 |
| 4 | Dodecane, 2,6,11-trimethyl- | 212 | C15H32 | 031295-56-4 | 53 |
| 5 | Decane, 2-methyl- | 156 | C11H24 | 006975-98-0 | 53 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

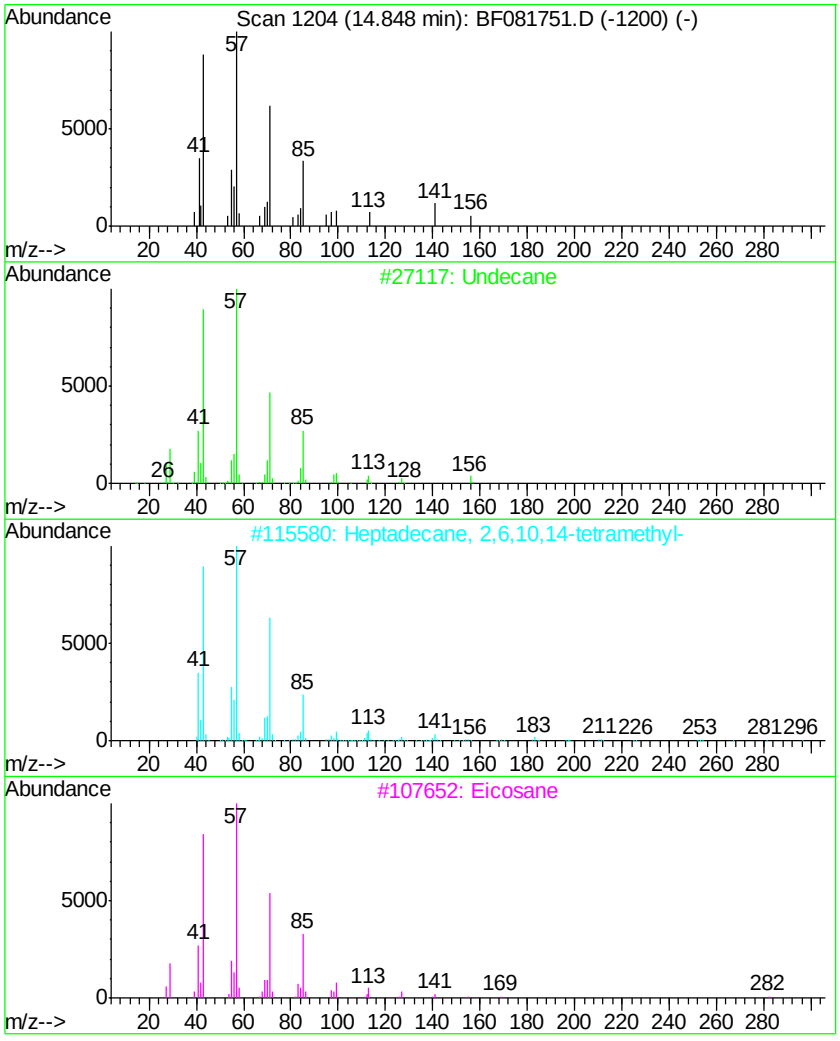
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

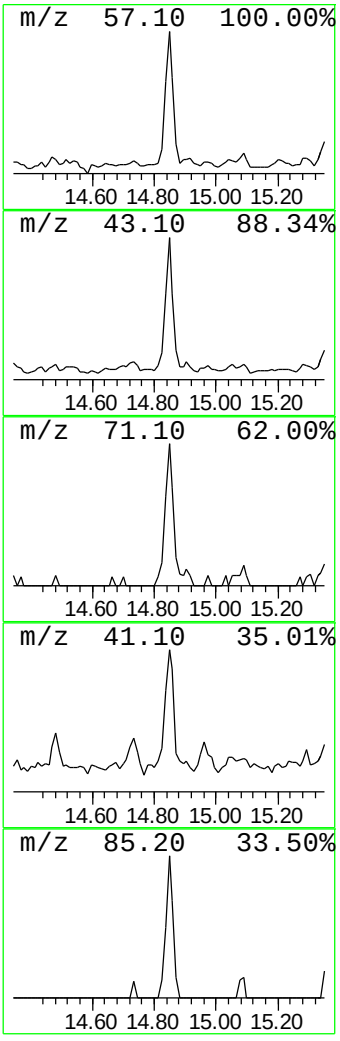
**********

Peak Number 10 Undecane Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.85 | 4.81 ng | 72769 | Acenaphthene-d10 | 15.16 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Undecane | 156 | C11H24 | 001120-21-4 | 90 |
| 2 | Heptadecane, 2,6,10,14-tetramethyl- | 296 | C21H44 | 018344-37-1 | 72 |
| 3 | Eicosane | 282 | C20H42 | 000112-95-8 | 72 |
| 4 | Hexadecane | 226 | C16H34 | 000544-76-3 | 72 |
| 5 | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 72 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

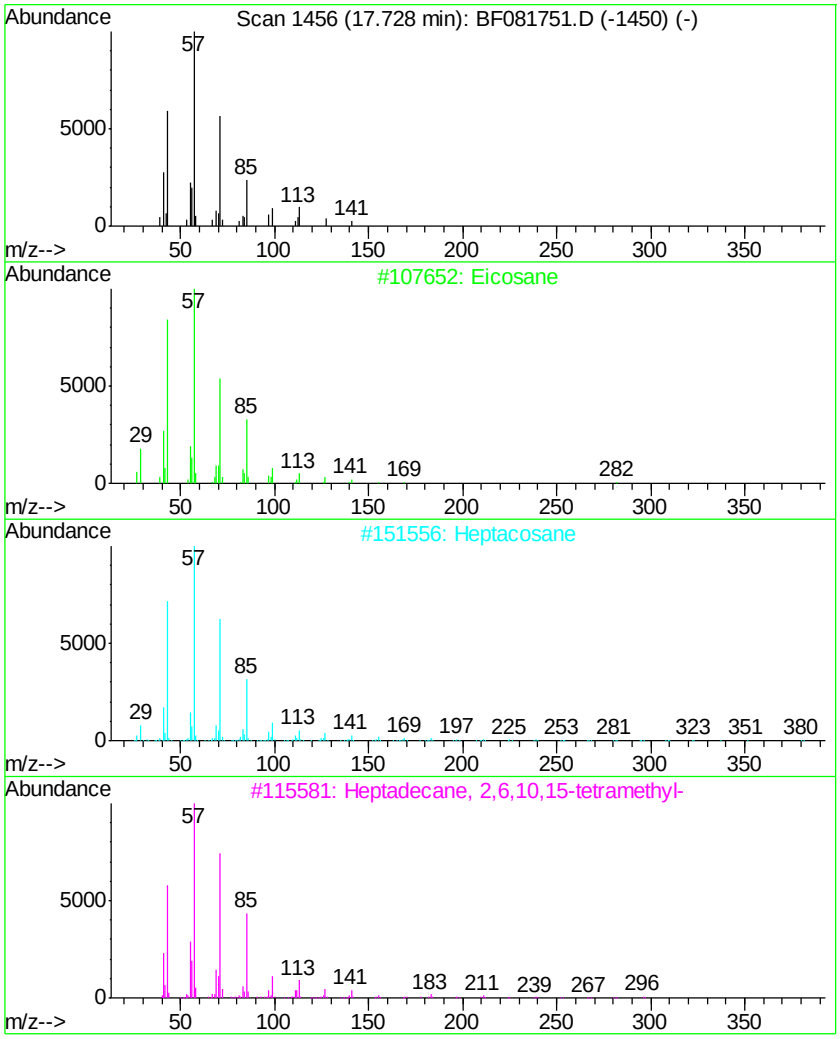
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

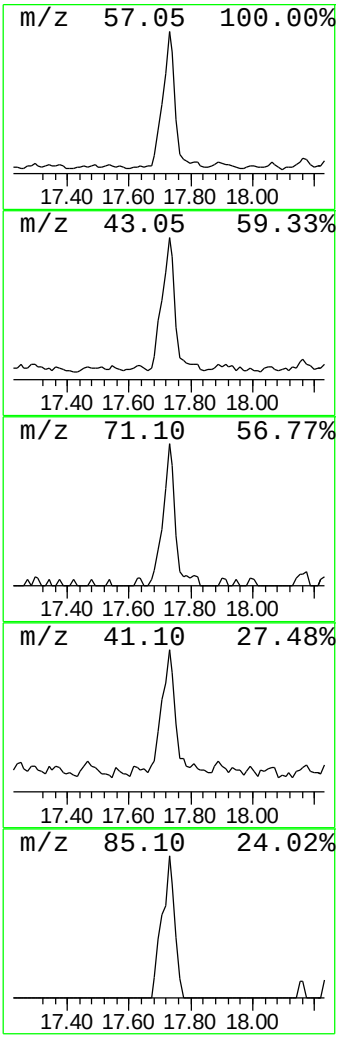
**********

Peak Number 11 Eicosane Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.73 | 8.73 ng | 123856 | Phenanthrene-d10 | 18.67 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Eicosane | 282 | C20H42 | 000112-95-8 | 80 |
| 2 | Heptacosane | 380 | C27H56 | 000593-49-7 | 72 |
| 3 | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44 | 054833-48-6 | 72 |
| 4 | Hexadecane | 226 | C16H34 | 000544-76-3 | 64 |
| 5 | Tetracosane | 338 | C24H50 | 000646-31-1 | 59 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

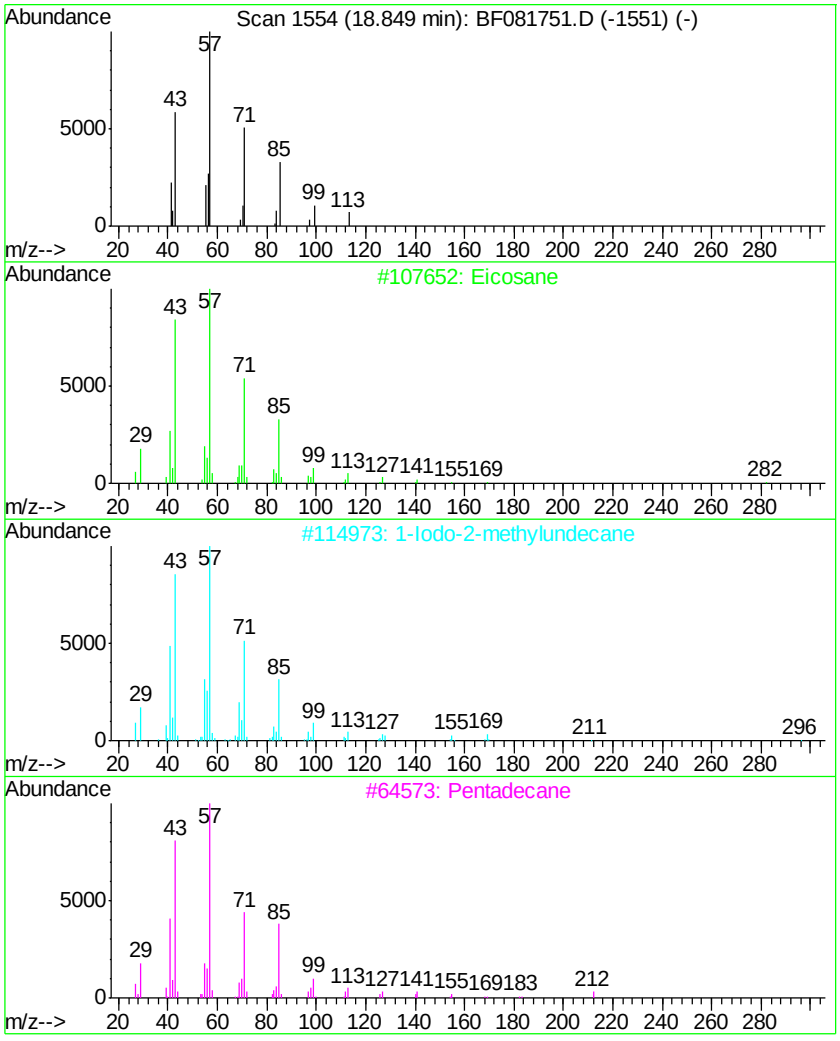
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

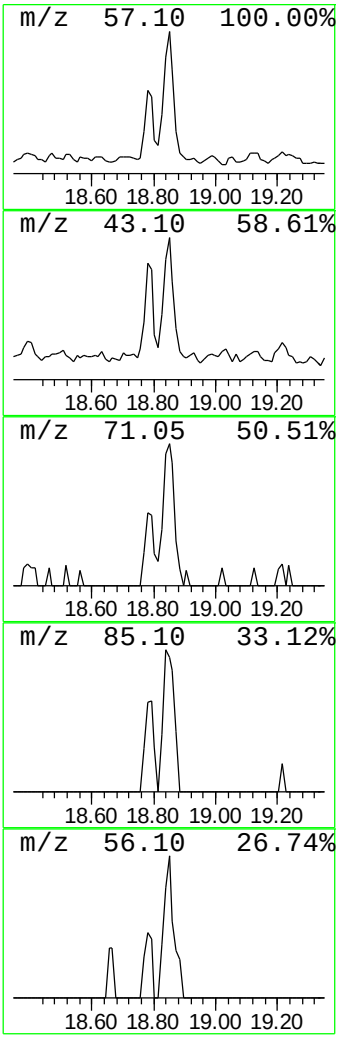
**********

Peak Number 12 1-Iodo-2-methylundecane Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.85 | 3.55 ng | 50295 | Phenanthrene-d10 | 18.67 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Eicosane | 282 | C20H42 | 000112-95-8 | 78 |
| 2 | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 72 |
| 3 | Pentadecane | 212 | C15H32 | 000629-62-9 | 72 |
| 4 | Hexadecane | 226 | C16H34 | 000544-76-3 | 72 |
| 5 | Tetradecane | 198 | C14H30 | 000629-59-4 | 64 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

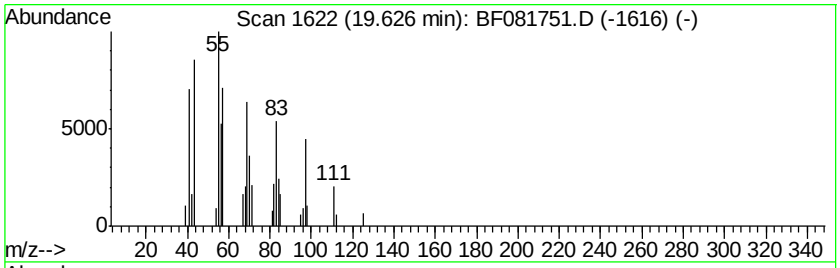
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

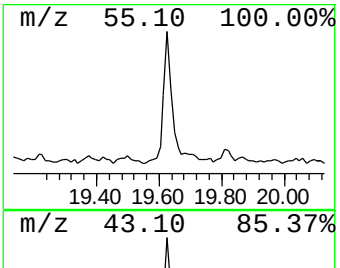
**********

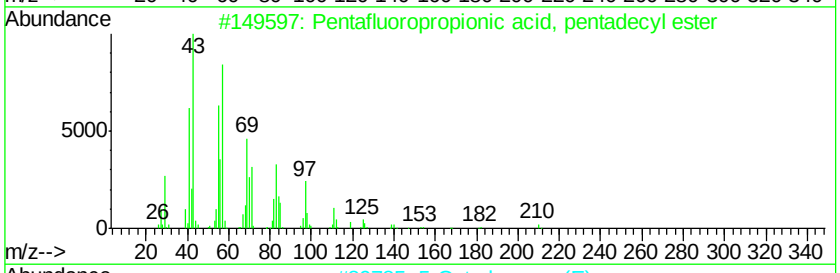
Peak Number 13 Pentafluoropropionic acid, ... Concentration Rank 11

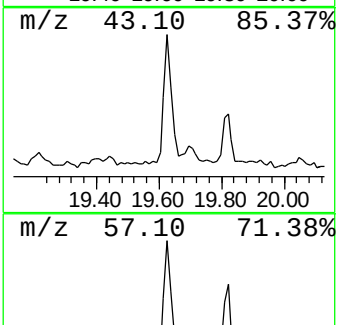
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.63 | 5.46 ng | 77367 | Phenanthrene-d10 | 18.67 |

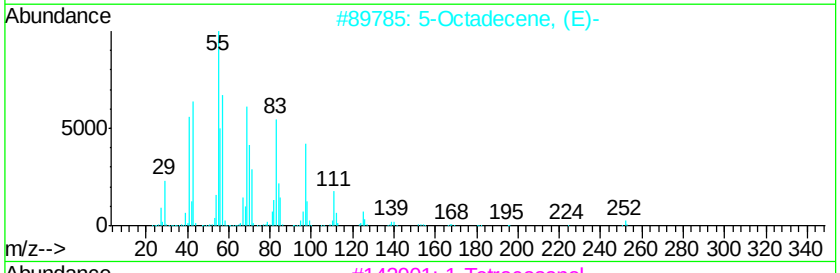
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentafluoropropionic acid, penta... | 374 | C18H31F5O2 | 1000280-07-5 | 80 |
| 2 | 5-Octadecene, (E)- | 252 | C18H36 | 007206-21-5 | 74 |
| 3 | 1-Tetracosanol | 354 | C24H50O | 000506-51-4 | 68 |
| 4 | Pentafluoropropionic acid, tetra... | 360 | C17H29F5O2 | 006222-06-6 | 64 |
| 5 | Pentafluoropropionic acid, octad... | 416 | C21H37F5O2 | 1000280-07-7 | 64 |

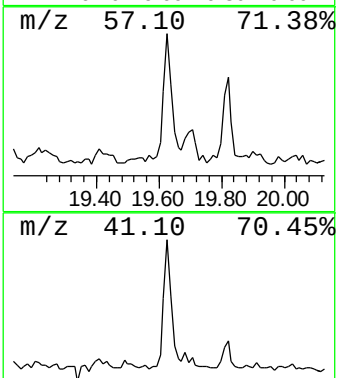


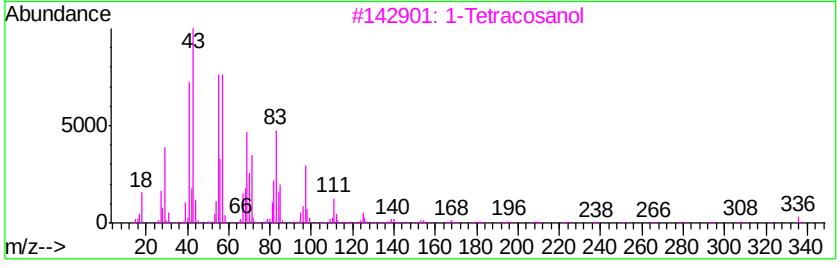



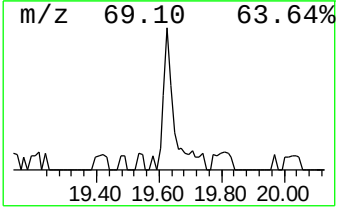



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

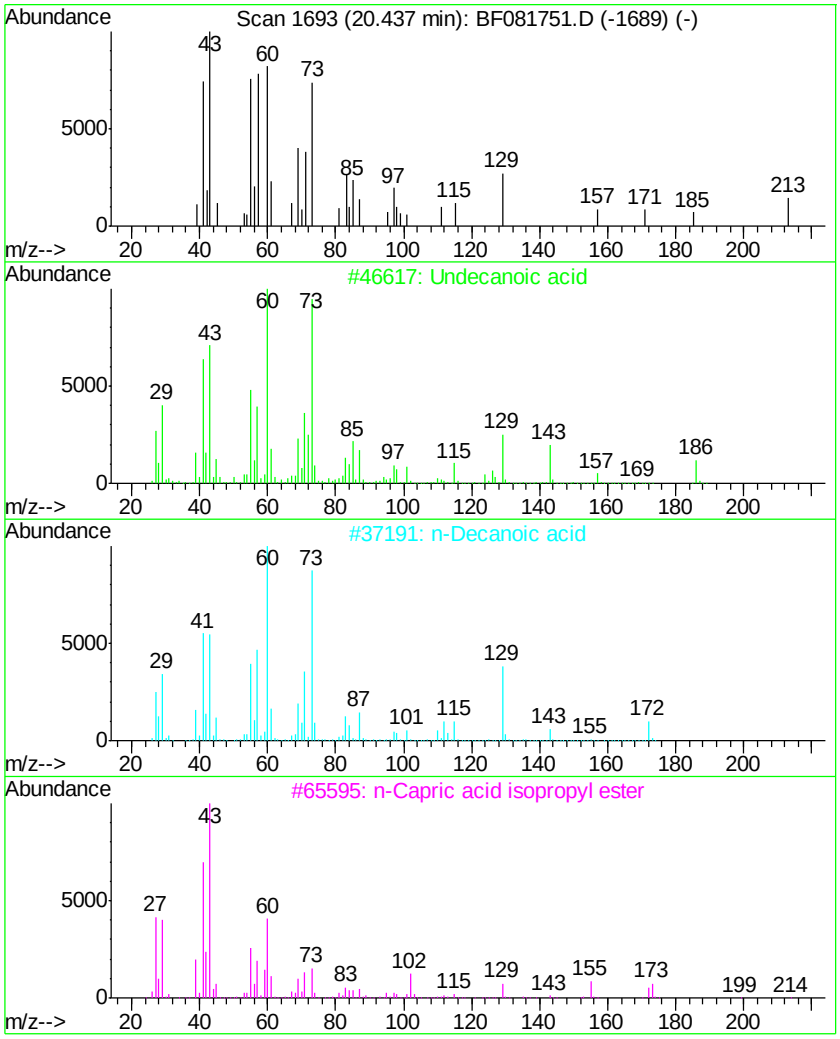
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

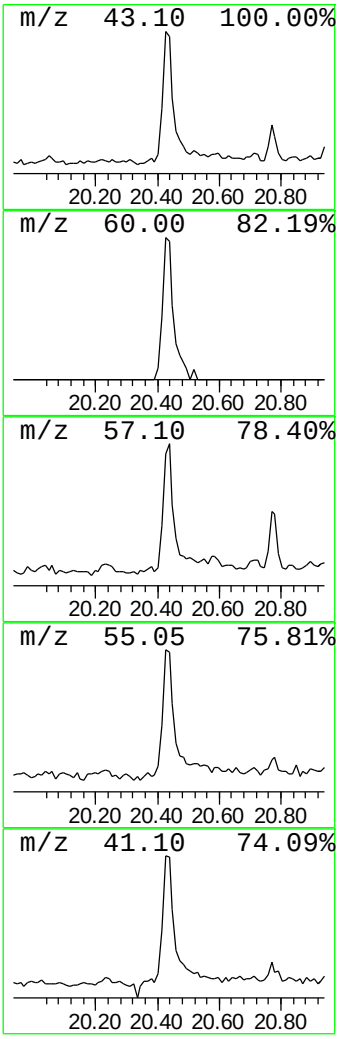
**********

Peak Number 14 Undecanoic acid Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.44 | 7.16 ng | 101600 | Phenanthrene-d10 | 18.67 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Undecanoic acid | 186 | C11H22O2 | 000112-37-8 | 59 |
| 2 | n-Decanoic acid | 172 | C10H20O2 | 000334-48-5 | 53 |
| 3 | n-Capric acid isopropyl ester | 214 | C13H26O2 | 002311-59-3 | 47 |
| 4 | d-Glucohexodialdose | 178 | C6H10O6 | 1000149-19-1 | 38 |
| 5 | 4-Propionyloxytetradecane | 270 | C17H34O2 | 1000245-62-9 | 14 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

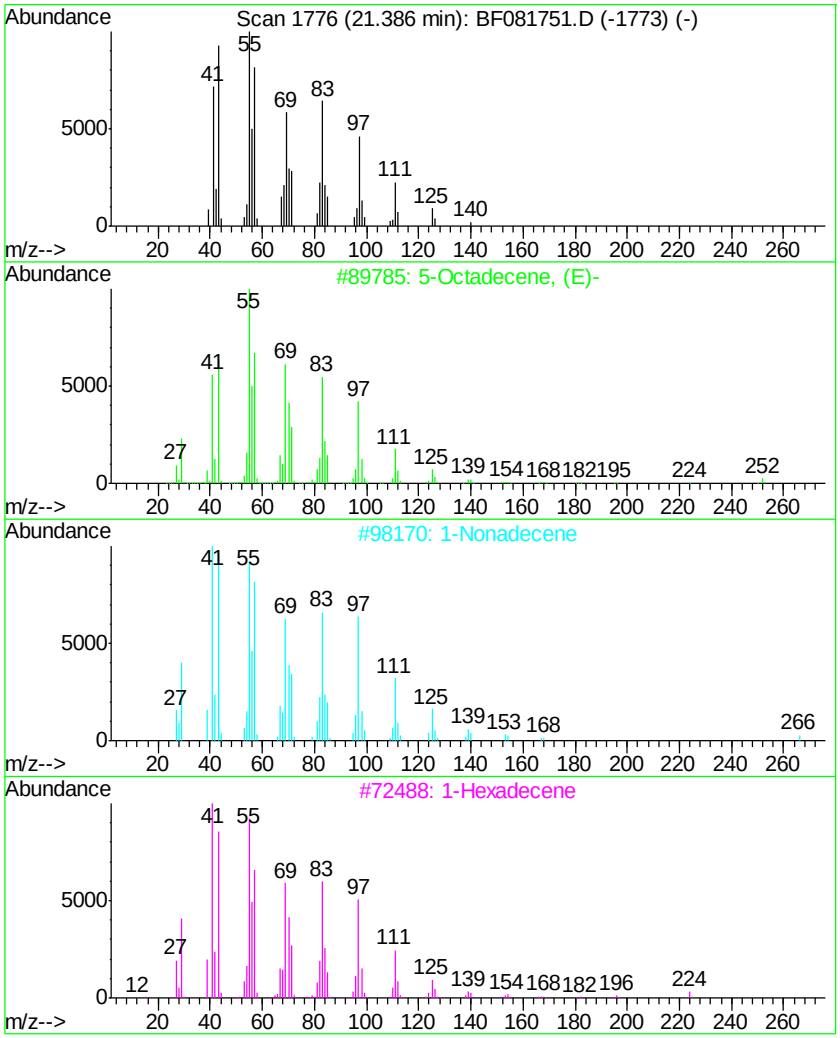
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

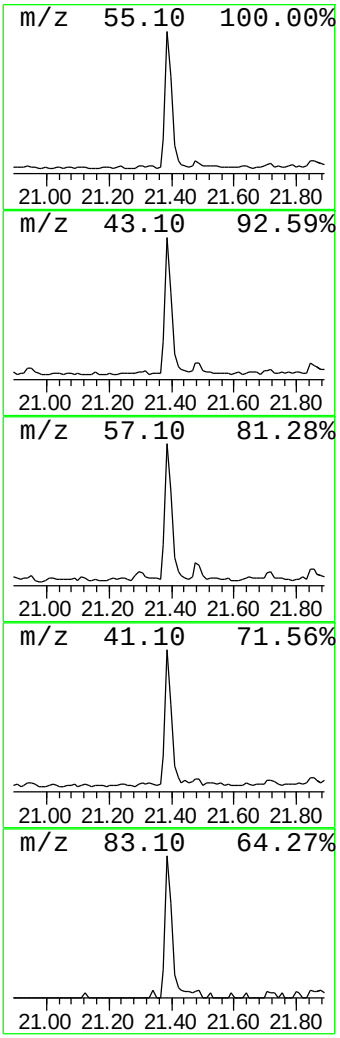
**********

Peak Number 15 5-Octadecene, (E)- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.39 | 14.52 ng | 177942 | Chrysene-d12 | 23.33 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 5-Octadecene, (E)- | 252 | C18H36 | 007206-21-5 | 91 |
| 2 | 1-Nonadecene | 266 | C19H38 | 018435-45-5 | 91 |
| 3 | 1-Hexadecene | 224 | C16H32 | 000629-73-2 | 91 |
| 4 | 5-Eicosene, (E)- | 280 | C20H40 | 074685-30-6 | 90 |
| 5 | 1-Octadecene | 252 | C18H36 | 000112-88-9 | 90 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092215\
Data File : BF081751.D
Acq On : 22 Sep 2015 21:52
Operator : UM/IZ
Sample : G3747-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TRENCH-1-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| | | | | | Internal Standard | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| Butane, 2-methoxy... | 1.22 | 96.7 | ng | 1159920 | 1 | 8.14 | 239862 | 20.0 |
| 2-Pentanone, 4-hy... | 4.42 | 60.8 | ng | 728864 | 1 | 8.14 | 239862 | 20.0 |
| unknown7.66 | 7.66 | 116.7 | ng | 1399340 | 1 | 8.14 | 239862 | 20.0 |
| Decane, 2,5,9-tri... | 12.31 | 3.2 | ng | 47451 | 2 | 11.03 | 298060 | 20.0 |
| Nonane, 3,7-dimet... | 13.74 | 5.8 | ng | 87071 | 3 | 15.16 | 302296 | 20.0 |
| Undecane | 14.85 | 4.8 | ng | 72769 | 3 | 15.16 | 302296 | 20.0 |
| Eicosane | 17.73 | 8.7 | ng | 123856 | 4 | 18.67 | 283609 | 20.0 |
| 1-Iodo-2-methylun... | 18.85 | 3.5 | ng | 50295 | 4 | 18.67 | 283609 | 20.0 |
| Pentafluoropropio... | 19.63 | 5.5 | ng | 77367 | 4 | 18.67 | 283609 | 20.0 |
| Undecanoic acid | 20.44 | 7.2 | ng | 101600 | 4 | 18.67 | 283609 | 20.0 |
| 5-Octadecene, (E)- | 21.39 | 14.5 | ng | 177942 | 5 | 23.33 | 245103 | 20.0 |