LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF073016\
Data File : BF089446.D
Acq On : 30 Jul 2016 9:39
Operator : UM/SJ
Sample : H4222-02
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-02-COMP

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-BF072716.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 | 1.678 | 6 | 8 | 48 | rVB | 1018959 | 2265898 | 100.00% | 20.755% |
| 2 | 4.604 | 262 | 264 | 277 | rVB | 188787 | 349925 | 15.44% | 3.205% |
| 3 | 5.084 | 304 | 306 | 330 | rBV | 635196 | 935748 | 41.30% | 8.571% |
| 4 | 6.182 | 399 | 402 | 409 | rBV | 775117 | 1002507 | 44.24% | 9.183% |
| 5 | 6.284 | 409 | 411 | 422 | rVB | 1047918 | 1025645 | 45.26% | 9.395% |
| 6 | 6.513 | 430 | 431 | 442 | rBV5 | 200885 | 243461 | 10.74% | 2.230% |
| 7 | 6.673 | 442 | 445 | 447 | rBV | 755242 | 745895 | 32.92% | 6.832% |
| 8 | 7.084 | 479 | 481 | 492 | rBV | 359565 | 559383 | 24.69% | 5.124% |
| 9 | 7.805 | 542 | 544 | 546 | rBV | 293129 | 275703 | 12.17% | 2.525% |
| 10 | 8.879 | 636 | 638 | 641 | rBV | 1072211 | 925880 | 40.86% | 8.481% |
| 11 | 9.279 | 671 | 673 | 676 | rBV | 155531 | 131154 | 5.79% | 1.201% |
| 12 | 9.542 | 694 | 696 | 699 | rBV | 306912 | 278496 | 12.29% | 2.551% |
| 13 | 10.331 | 763 | 765 | 773 | rVB | 481556 | 466745 | 20.60% | 4.275% |
| 14 | 10.468 | 775 | 777 | 780 | rBV | 24560 | 27027 | 1.19% | 0.248% |
| 15 | 11.016 | 823 | 825 | 830 | rBV | 276927 | 251415 | 11.10% | 2.303% |
| 16 | 11.268 | 845 | 847 | 851 | rBV | 17010 | 25566 | 1.13% | 0.234% |
| 17 | 12.091 | 917 | 919 | 924 | rVB | 16513 | 30339 | 1.34% | 0.278% |
| 18 | 12.594 | 961 | 963 | 966 | rBV | 620292 | 784618 | 34.63% | 7.187% |
| 19 | 13.508 | 1041 | 1043 | 1052 | rBV | 34595 | 61823 | 2.73% | 0.566% |
| 20 | 13.634 | 1052 | 1054 | 1057 | rBV | 304362 | 299404 | 13.21% | 2.743% |
| 21 | 14.148 | 1095 | 1099 | 1102 | rBV2 | 12285 | 26921 | 1.19% | 0.247% |
| 22 | 14.971 | 1168 | 1171 | 1174 | rBV | 172116 | 203561 | 8.98% | 1.865% |

Sum of corrected areas: 10917114

LSC Report - Integrated Chromatogram

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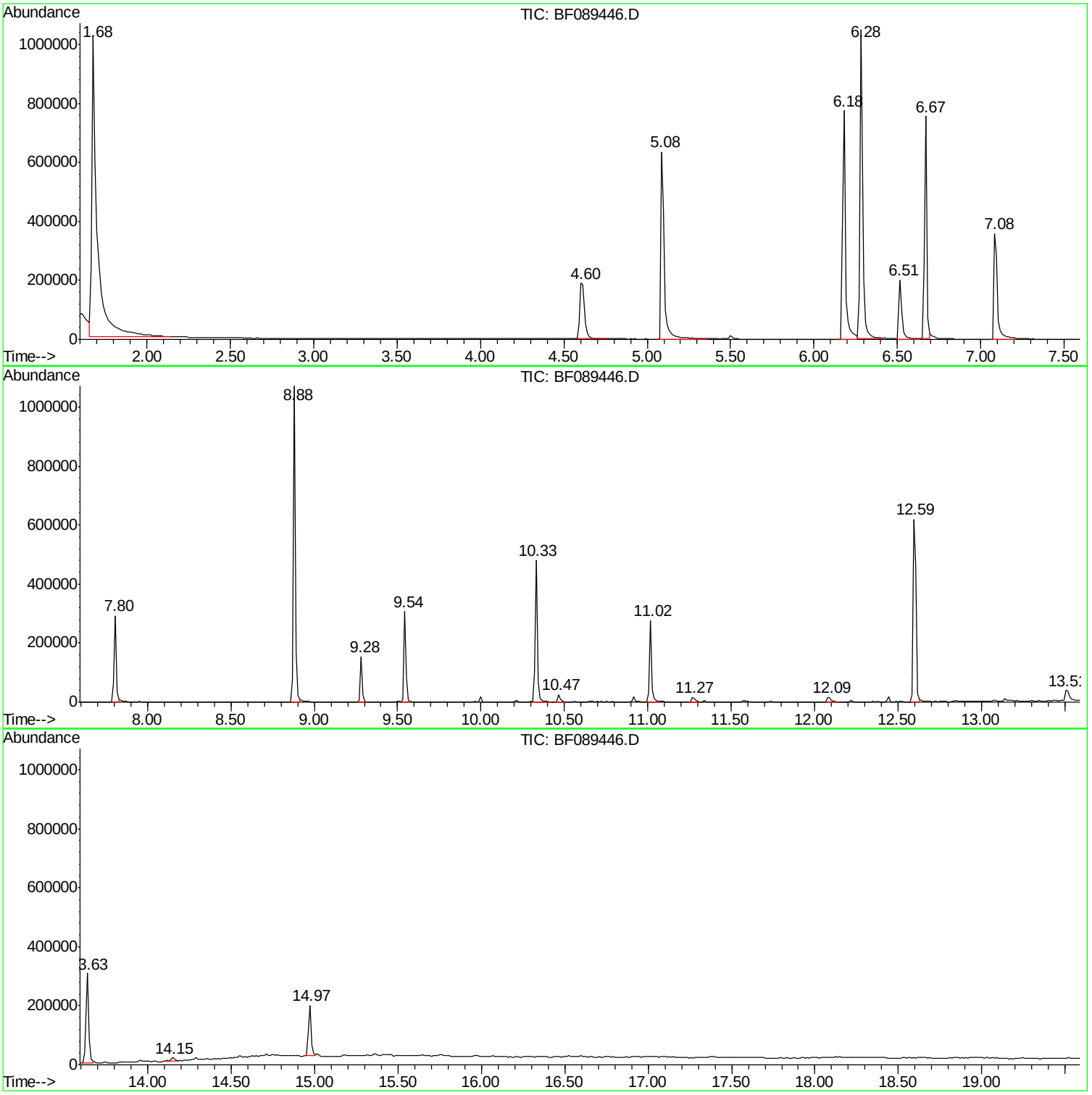
**Instrument :** BNA_F
**ClientSampleId :** SB-02-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.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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Operator : UM/SJ
Sample : H4222-02
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-02-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.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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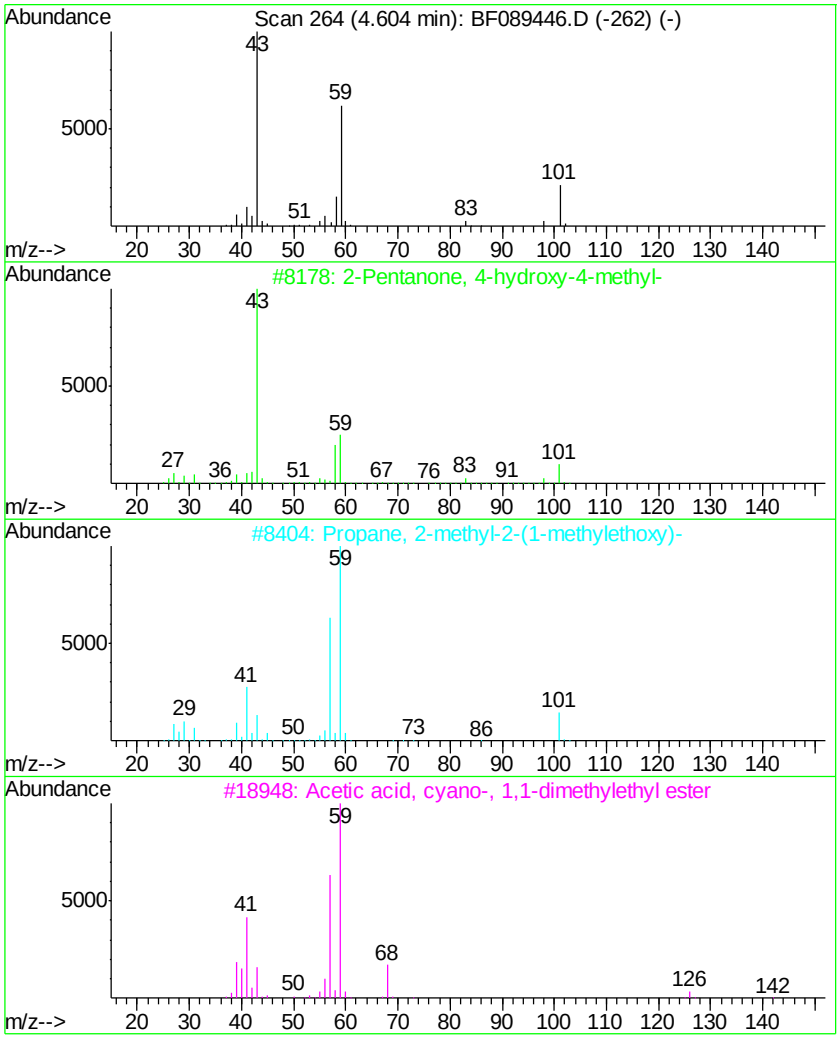
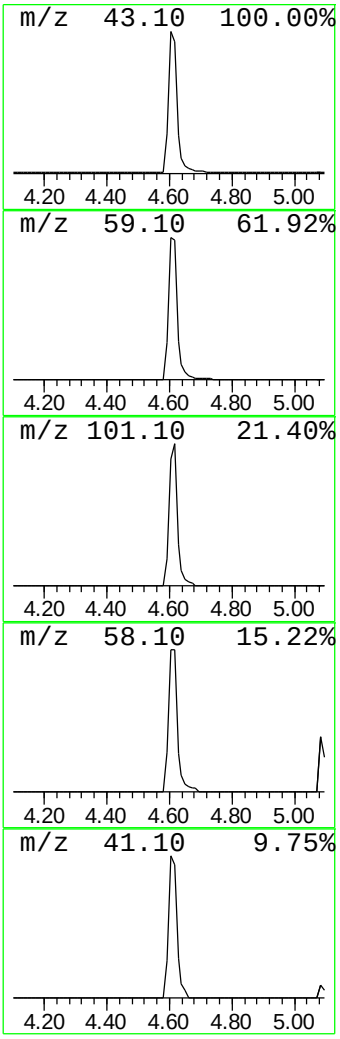
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.60 | 28.75 ng | 349925 | 1,4-Dichlorobenzene-d4 | 6.51 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 56 |
| 2 | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O | 017348-59-3 | 42 |
| 3 | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 23 |
| 4 | Morpholine, 4-methyl- | 101 | C5H11NO | 000109-02-4 | 9 |
| 5 | Butane, 1-ethoxy- | 102 | C6H14O | 000628-81-9 | 9 |



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Data File : BF089446.D
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Operator : UM/SJ
Sample : H4222-02
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-02-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.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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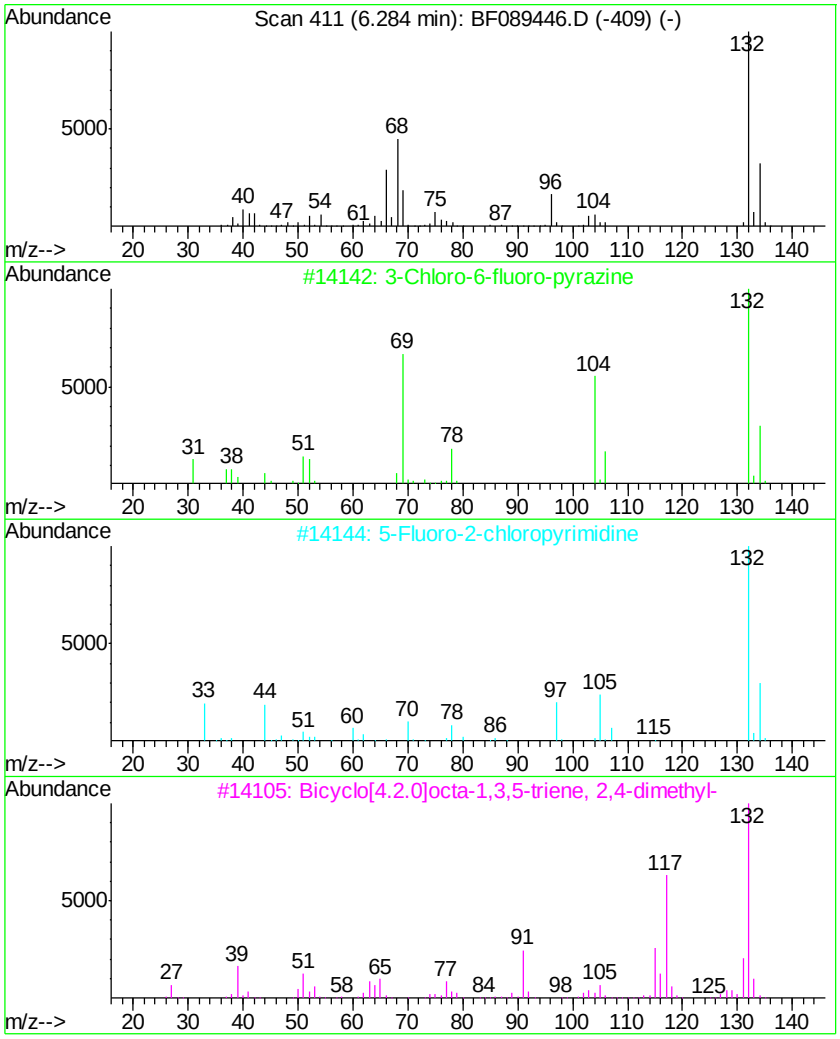
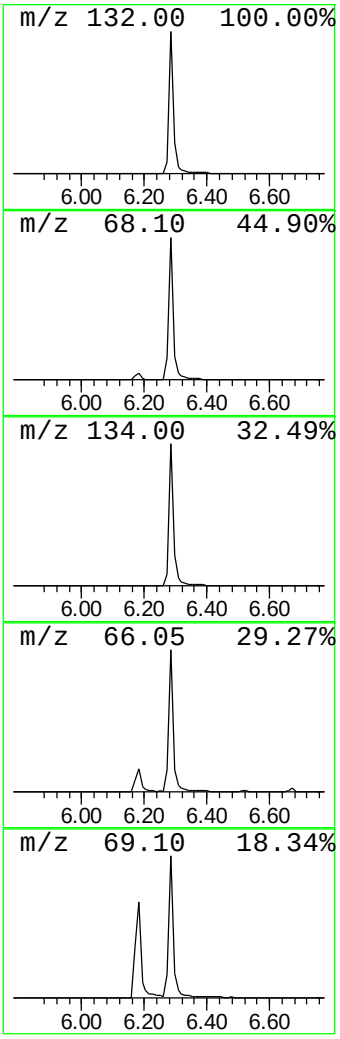
Peak Number 3 unknown6.28 Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.28 | 84.26 ng | 1025650 | 1,4-Dichlorobenzene-d4 | 6.51 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Chloro-6-fluoro-pyrazine | 132 | C4H2ClFN2 | 1000146-10-7 | 27 |
| 2 | 5-Fluoro-2-chloropyrimidine | 132 | C4H2ClFN2 | 062802-42-0 | 12 |
| 3 | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12 | 028749-81-7 | 9 |
| 4 | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO | 031357-76-3 | 9 |
| 5 | 1-Methylpyrrolo[1,2-a]pyrazine | 132 | C8H8N2 | 064608-59-9 | 9 |



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF073016\
Data File : BF089446.D
Acq On : 30 Jul 2016 9:39
Operator : UM/SJ
Sample : H4222-02
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-02-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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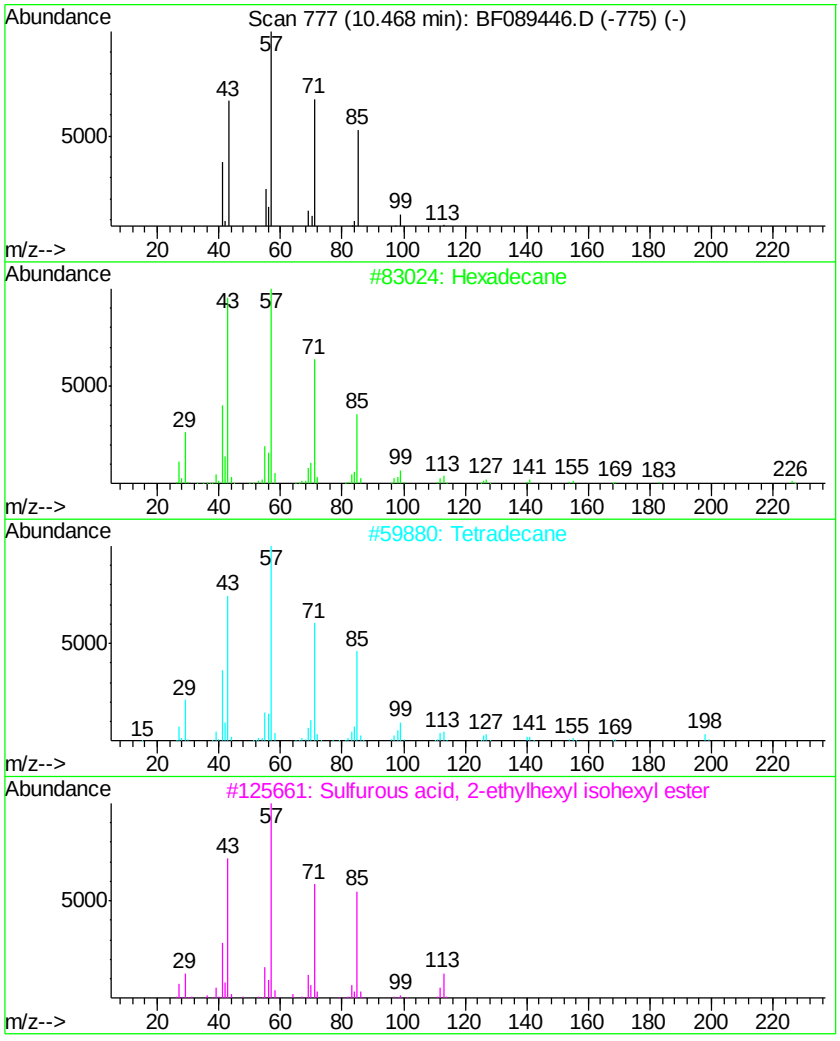
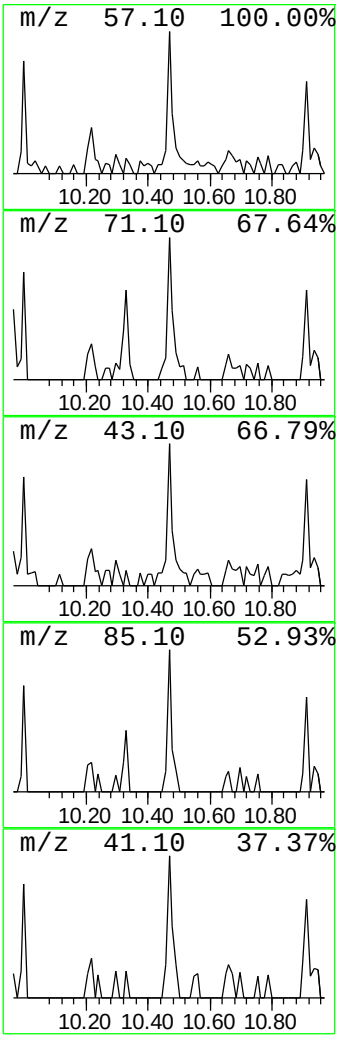
Peak Number 5 Hexadecane Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.47 | 2.15 ng | 27027 | Phenanthrene-d10 | 11.02 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexadecane | 226 | C16H34 | 000544-76-3 | 78 |
| 2 | Tetradecane | 198 | C14H30 | 000629-59-4 | 78 |
| 3 | Sulfurous acid, 2-ethylhexyl iso... | 278 | C14H30O3S | 1000309-19-0 | 72 |
| 4 | Sulfurous acid, 2-ethylhexyl hex... | 278 | C14H30O3S | 1000309-20-2 | 64 |
| 5 | Sulfurous acid, hexyl pentyl ester | 236 | C11H24O3S | 1000309-14-1 | 64 |



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Data File : BF089446.D
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Operator : UM/SJ
Sample : H4222-02
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-02-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
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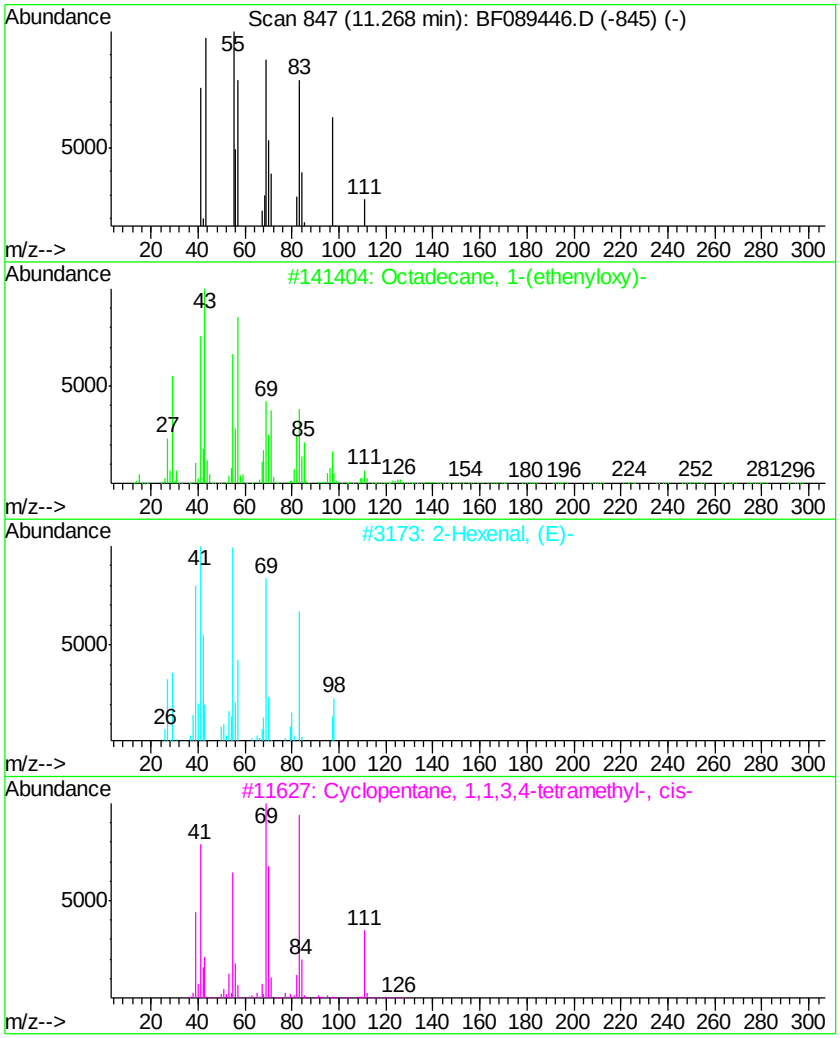
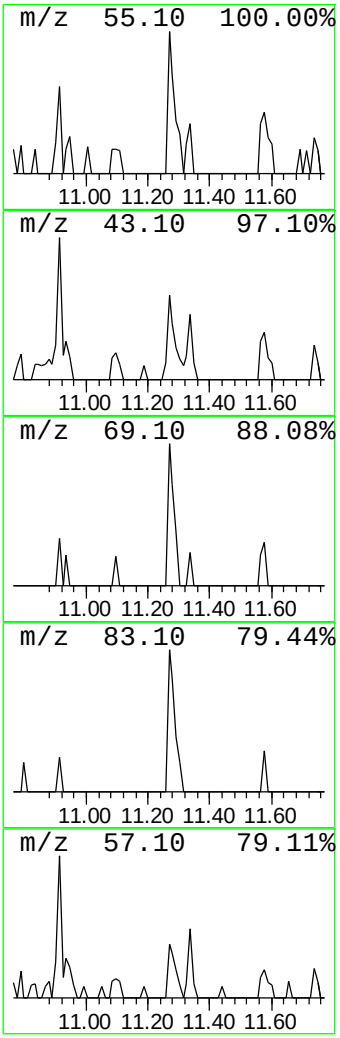
Peak Number 6 unknown11.27 Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.27 | 2.03 ng | 25566 | Phenanthrene-d10 | 11.02 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecane, 1-(ethenyloxy)- | 296 | C20H40O | 000930-02-9 | 45 |
| 2 | 2-Hexenal, (E)- | 98 | C6H10O | 006728-26-3 | 43 |
| 3 | Cyclopentane, 1,1,3,4-tetramethy... | 126 | C9H18 | 053907-60-1 | 38 |
| 4 | 1-Fluorononane | 146 | C9H19F | 000463-18-3 | 38 |
| 5 | 2,4,6,8-Tetramethyl-1-undecene | 210 | C15H30 | 059920-26-2 | 38 |



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Operator : UM/SJ
Sample : H4222-02
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-02-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

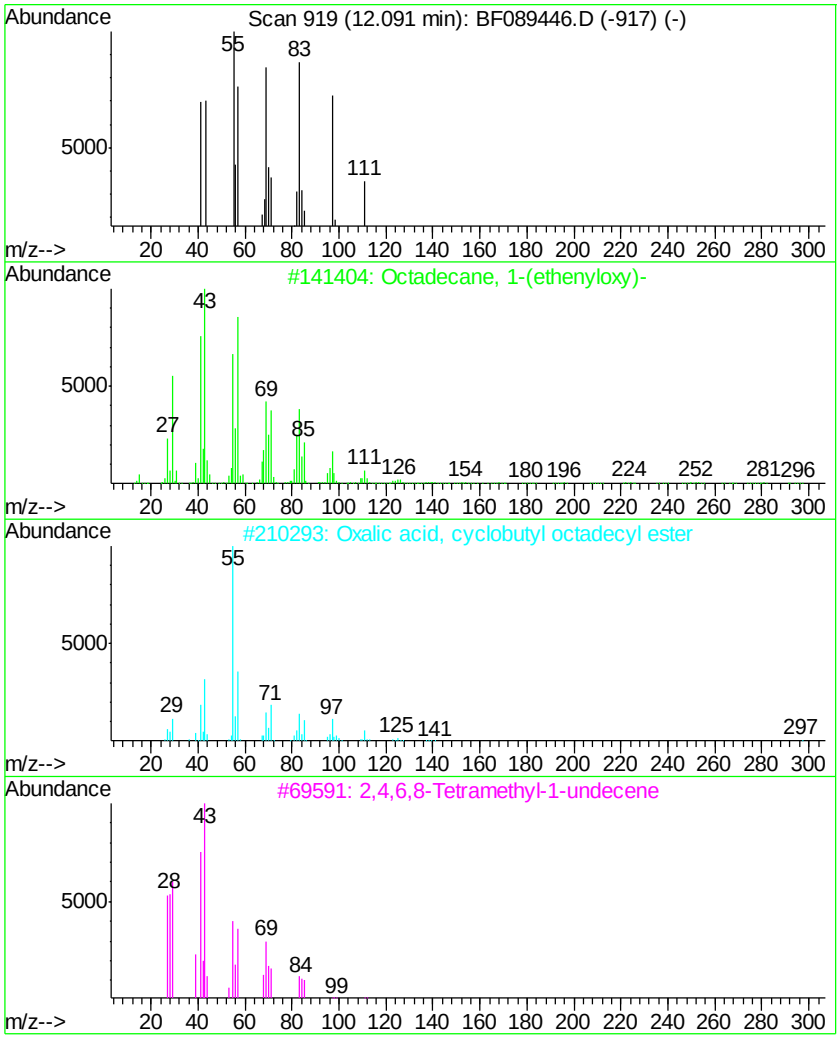
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

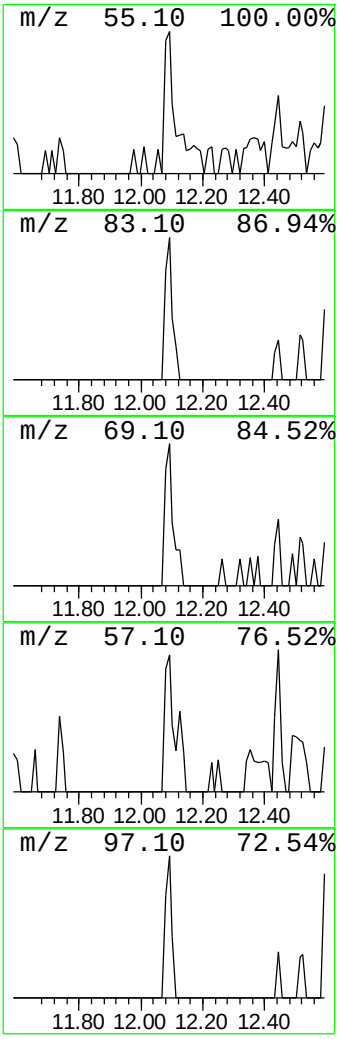
**********

Peak Number 7 Octadecane, 1-(ethenyloxy)- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.09 | 2.41 ng | 30339 | Phenanthrene-d10 | 11.02 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecane, 1-(ethenyloxy)- | 296 | C20H40O | 000930-02-9 | 78 |
| 2 | Oxalic acid, cyclobutyl octadecy... | 396 | C24H44O4 | 1000309-70-8 | 47 |
| 3 | 2,4,6,8-Tetramethyl-1-undecene | 210 | C15H30 | 059920-26-2 | 47 |
| 4 | Oxalic acid, allyl octadecyl ester | 382 | C23H42O4 | 1000309-24-5 | 45 |
| 5 | (S)-(+)-5-Methyl-1-heptanol | 130 | C8H18O | 057803-73-3 | 43 |



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Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
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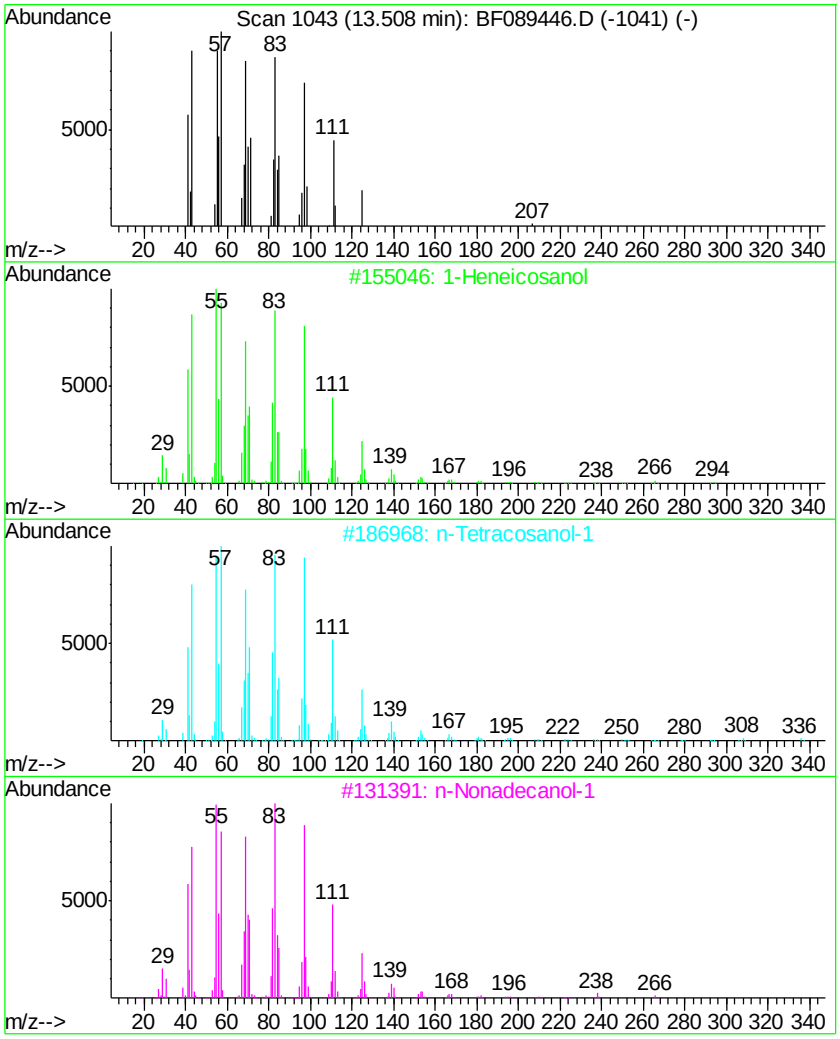
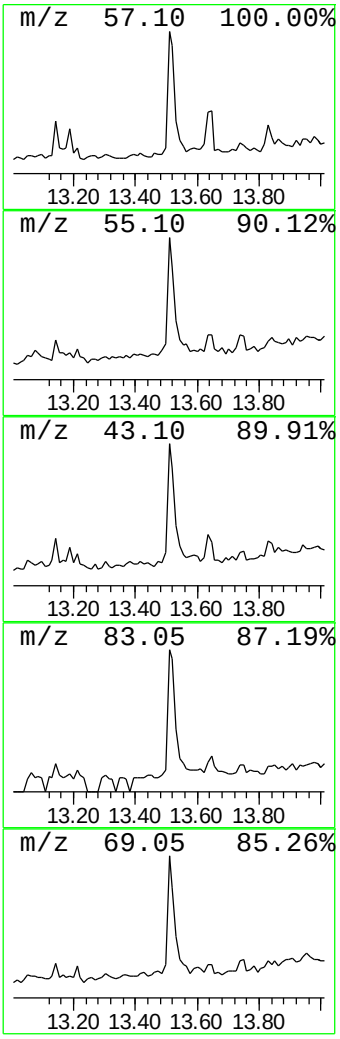
Peak Number 8 1-Heneicosanol Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.51 | 4.13 ng | 61823 | Chrysene-d12 | 13.63 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Heneicosanol | 312 | C21H44O | 015594-90-8 | 94 |
| 2 | n-Tetracosanol-1 | 354 | C24H50O | 000506-51-4 | 91 |
| 3 | n-Nonadecanol-1 | 284 | C19H40O | 001454-84-8 | 91 |
| 4 | Trichloroacetic acid, tetradecyl... | 358 | C16H29Cl3O2 | 074339-52-9 | 91 |
| 5 | Pentafluoropropionic acid, tetra... | 360 | C17H29F5O2 | 006222-06-6 | 91 |



# Tentatively Identified Compound (LSC) summary

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

| | | | | | Internal Standard | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| 2-Pentanone, 4-hy... | 4.60 | 28.8 | ng | 349925 | 1 | 6.51 | 243461 | 20.0 |
| unknown6.28 | 6.28 | 84.3 | ng | 1025650 | 1 | 6.51 | 243461 | 20.0 |
| Hexadecane | 10.47 | 2.1 | ng | 27027 | 4 | 11.02 | 251415 | 20.0 |
| unknown11.27 | 11.27 | 2.0 | ng | 25566 | 4 | 11.02 | 251415 | 20.0 |
| Octadecane, 1-(et... | 12.09 | 2.4 | ng | 30339 | 4 | 11.02 | 251415 | 20.0 |
| 1-Heneicosanol | 13.51 | 4.1 | ng | 61823 | 5 | 13.63 | 299404 | 20.0 |