LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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-BF032217.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 | 1.916 | 3 | 5 | 10 | rVB | 81571 | 86268 | 1.91% | 0.211% |
| 2 | 4.004 | 356 | 360 | 372 | rBV | 485673 | 629816 | 13.91% | 1.540% |
| 3 | 4.404 | 422 | 428 | 431 | rBV | 3733110 | 3868161 | 85.43% | 9.459% |
| 4 | 5.475 | 603 | 610 | 613 | rBV | 3436425 | 4040968 | 89.25% | 9.881% |
| 5 | 5.610 | 628 | 633 | 636 | rBV | 3886347 | 4144557 | 91.54% | 10.134% |
| 6 | 5.863 | 672 | 676 | 680 | rBV | 1240177 | 1097640 | 24.24% | 2.684% |
| 7 | 6.045 | 702 | 707 | 710 | rBV | 3185759 | 3256312 | 71.92% | 7.962% |
| 8 | 6.516 | 780 | 787 | 790 | rBV | 2406104 | 2665103 | 58.86% | 6.517% |
| 9 | 7.063 | 876 | 880 | 887 | rBV2 | 228782 | 223355 | 4.93% | 0.546% |
| 10 | 7.363 | 926 | 931 | 935 | rBV | 1431318 | 1471362 | 32.50% | 3.598% |
| 11 | 8.133 | 1056 | 1062 | 1066 | rBV | 2341212 | 2419621 | 53.44% | 5.917% |
| 12 | 8.692 | 1150 | 1157 | 1160 | rBV | 4235056 | 4527700 | 100.00% | 11.071% |
| 13 | 9.186 | 1236 | 1241 | 1244 | rBV | 342661 | 342046 | 7.55% | 0.836% |
| 14 | 9.504 | 1287 | 1295 | 1301 | rBV2 | 1573233 | 1588019 | 35.07% | 3.883% |
| 15 | 9.769 | 1334 | 1340 | 1344 | rBV5 | 22016 | 52364 | 1.16% | 0.128% |
| 16 | 10.098 | 1391 | 1396 | 1401 | rBV | 127960 | 117020 | 2.58% | 0.286% |
| 17 | 10.375 | 1434 | 1443 | 1445 | rBV | 35382 | 62233 | 1.37% | 0.152% |
| 18 | 10.480 | 1455 | 1461 | 1464 | rBV | 2047596 | 2259523 | 49.90% | 5.525% |
| 19 | 10.680 | 1492 | 1495 | 1498 | rBV | 123345 | 133486 | 2.95% | 0.326% |
| 20 | 11.239 | 1585 | 1590 | 1593 | rBV | 70253 | 73126 | 1.62% | 0.179% |
| 21 | 11.327 | 1600 | 1605 | 1609 | rBV | 1399697 | 1338036 | 29.55% | 3.272% |
| 22 | 13.310 | 1936 | 1942 | 1945 | rBV | 2702605 | 3109493 | 68.68% | 7.603% |
| 23 | 14.468 | 2135 | 2139 | 2149 | rBV3 | 74385 | 147127 | 3.25% | 0.360% |
| 24 | 14.574 | 2152 | 2157 | 2161 | rBV | 969717 | 1011864 | 22.35% | 2.474% |
| 25 | 15.968 | 2391 | 2394 | 2398 | rBV | 85602 | 92916 | 2.05% | 0.227% |
| 26 | 16.204 | 2429 | 2434 | 2441 | rBV | 807830 | 919784 | 20.31% | 2.249% |
| 27 | 16.327 | 2451 | 2455 | 2462 | rVB | 151163 | 192229 | 4.25% | 0.470% |
| 28 | 16.721 | 2518 | 2522 | 2529 | rVB | 174998 | 242168 | 5.35% | 0.592% |
| 29 | 17.180 | 2595 | 2600 | 2608 | rVB2 | 161428 | 290768 | 6.42% | 0.711% |
| 30 | 17.703 | 2684 | 2689 | 2698 | rVB | 105510 | 215421 | 4.76% | 0.527% |
| 31 | 18.327 | 2790 | 2795 | 2804 | rVB3 | 77850 | 182326 | 4.03% | 0.446% |
| 32 | 19.050 | 2912 | 2918 | 2926 | rVB4 | 40288 | 95180 | 2.10% | 0.233% |

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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-BF032217.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Sum of corrected areas: 40895992

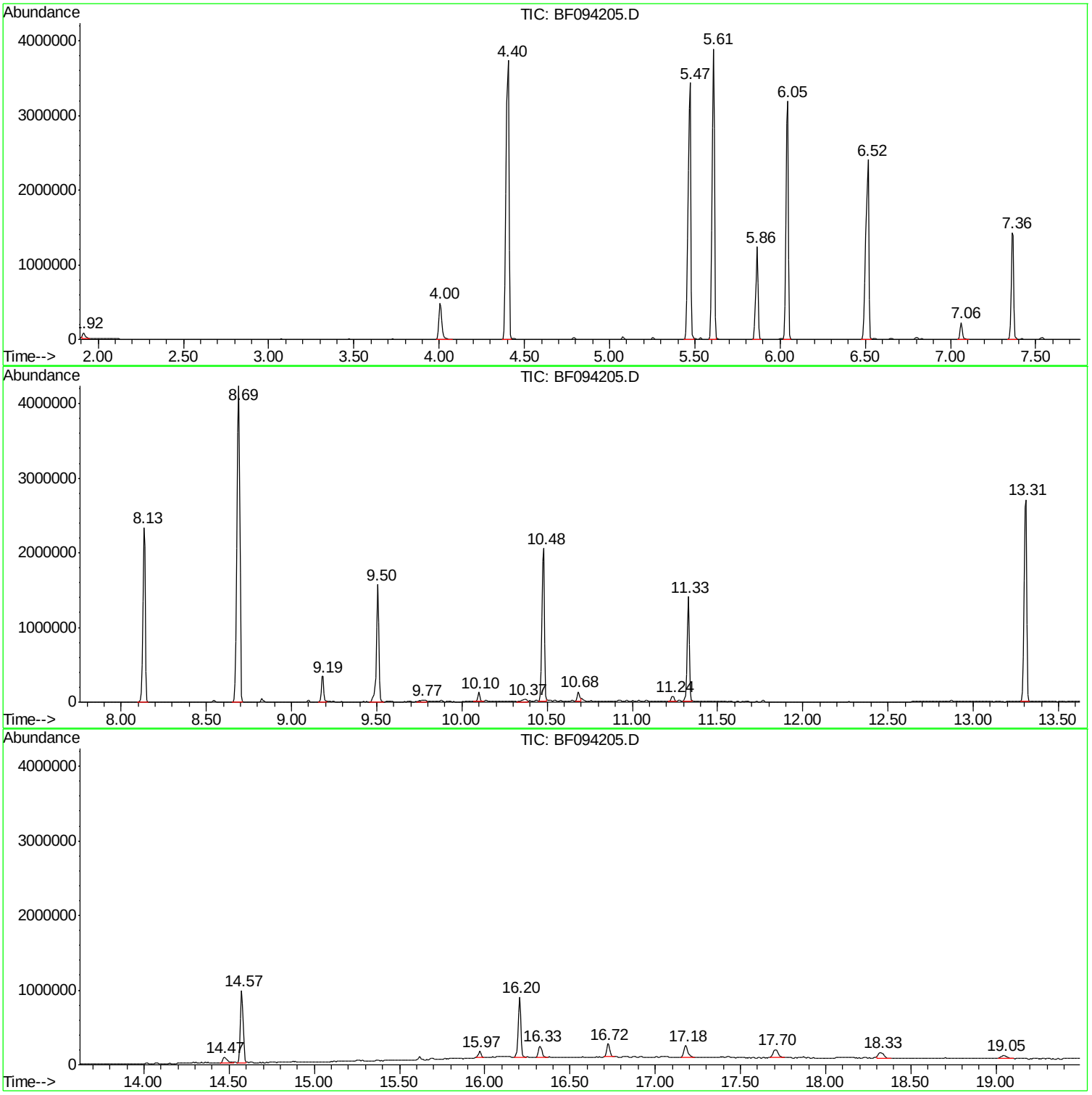
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(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



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
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Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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

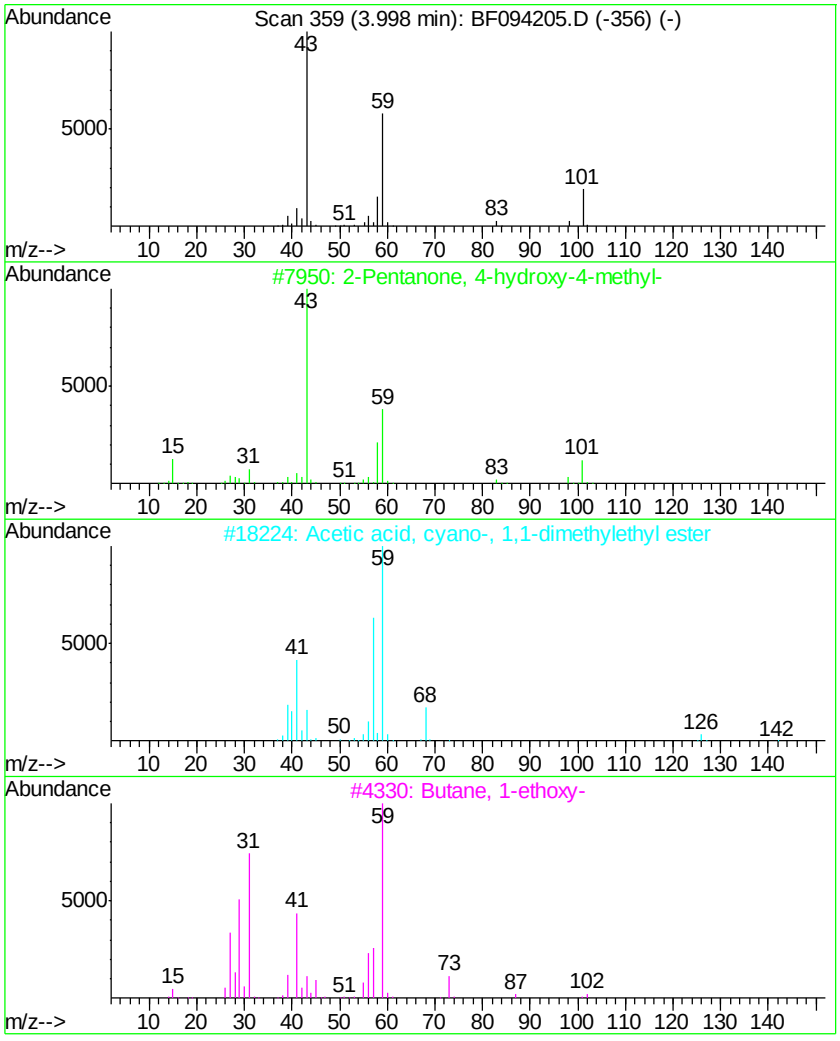
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

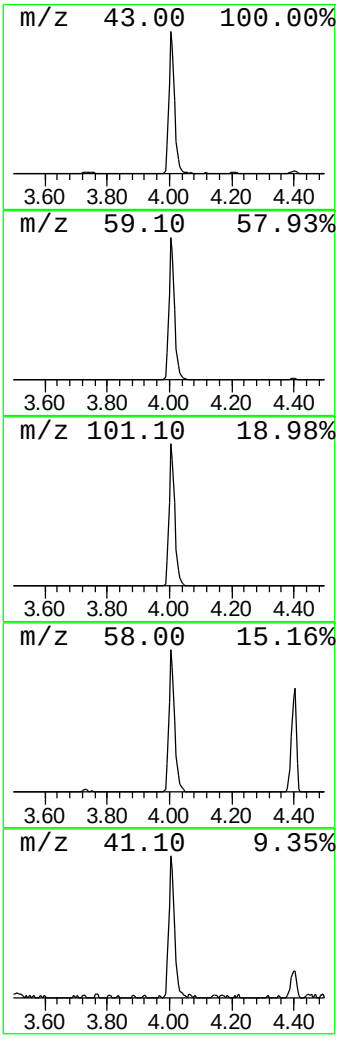
**********

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

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.00 | 11.48 ng | 629816 | 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 | 56 |
| 2 | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 17 |
| 3 | Butane, 1-ethoxy- | 102 | C6H14O | 000628-81-9 | 9 |
| 4 | Morpholine, 4-methyl- | 101 | C5H11NO | 000109-02-4 | 9 |
| 5 | 2,3-Butanedione, monooxime | 101 | C4H7NO2 | 000057-71-6 | 9 |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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

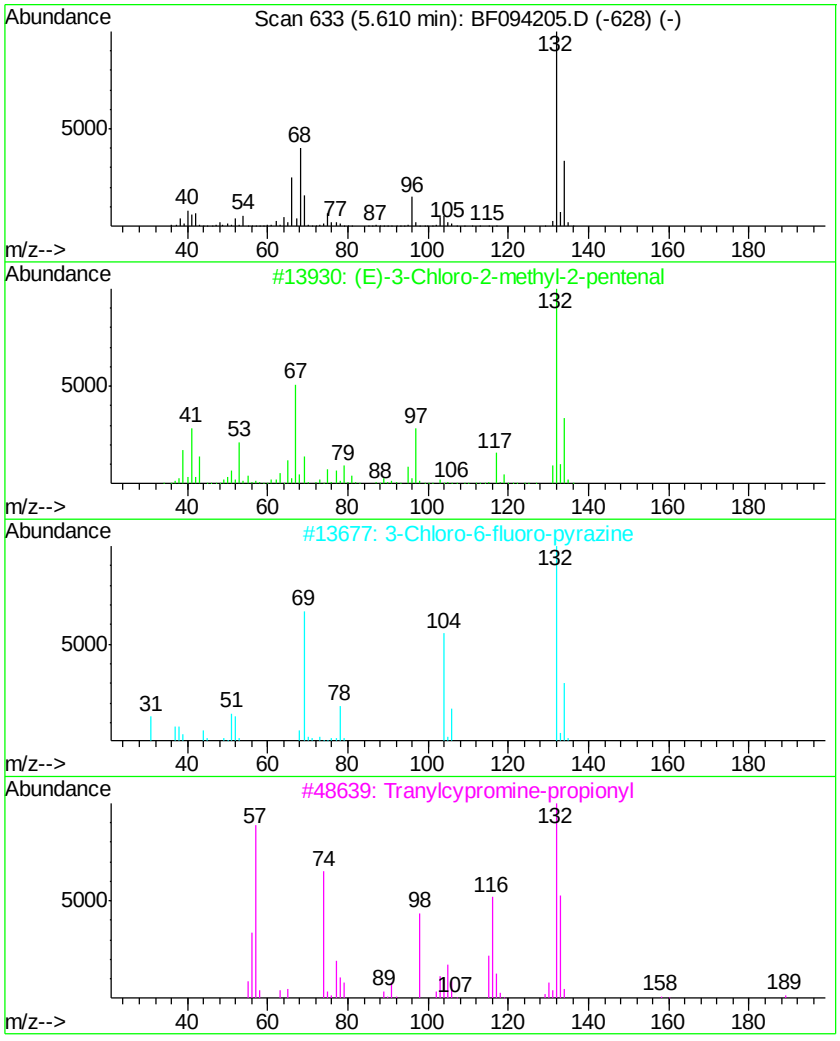
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

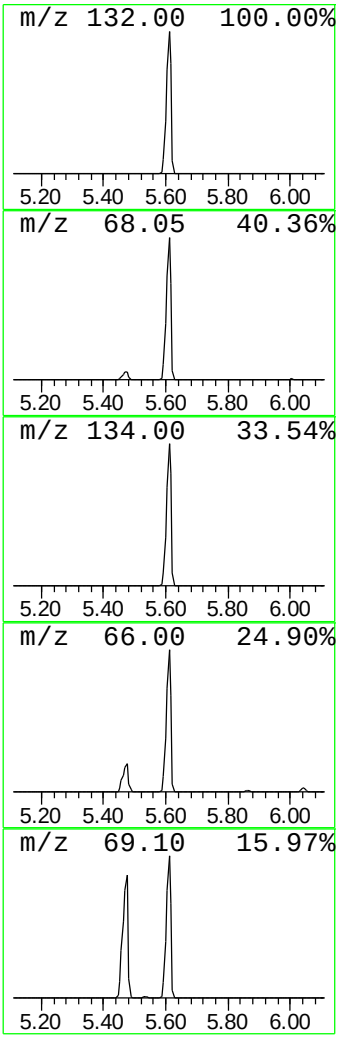
**********

Peak Number 2 unknown5.61 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.61 | 75.52 ng | 4144560 | 1,4-Dichlorobenzene-d4 | 5.86 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO | 031357-76-3 | 16 |
| 2 | 3-Chloro-6-fluoro-pyrazine | 132 | C4H2ClFN2 | 1000146-10-7 | 16 |
| 3 | Tranylcypromine-propionyl | 189 | C12H15NO | 1000123-86-3 | 10 |
| 4 | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12 | 028749-81-7 | 9 |
| 5 | 4-Chloro-2-fluoro-pyrimidine | 132 | C4H2ClFN2 | 051422-00-5 | 9 |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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

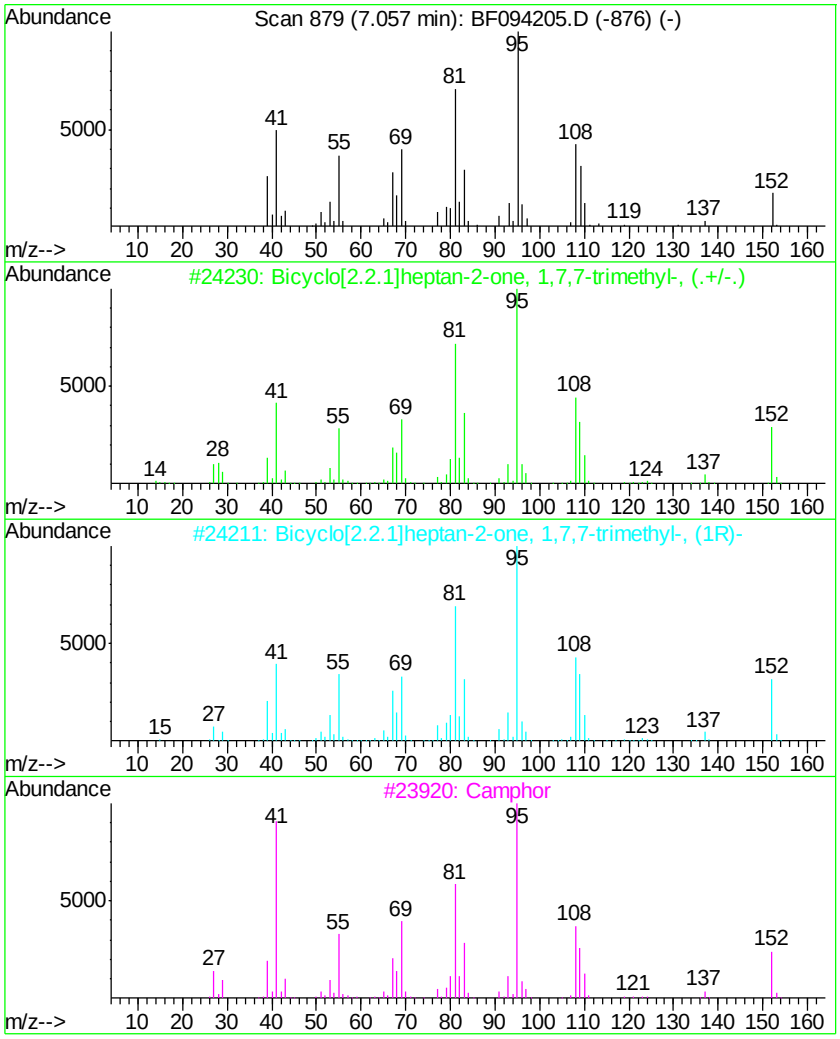
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

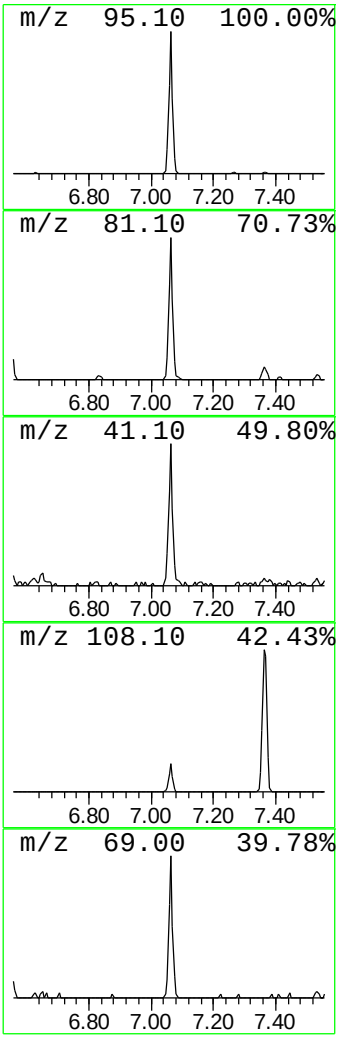
**********

Peak Number 4 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.06 | 3.04 ng | 223355 | Naphthalene-d8 | 7.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 021368-68-3 | 97 |
| 2 | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 000464-49-3 | 97 |
| 3 | Camphor | 152 | C10H16O | 000076-22-2 | 96 |
| 4 | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 000464-48-2 | 93 |
| 5 | 5,7-Octadien-4-one, 2,6-dimethyl... | 152 | C10H16O | 006752-80-3 | 49 |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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

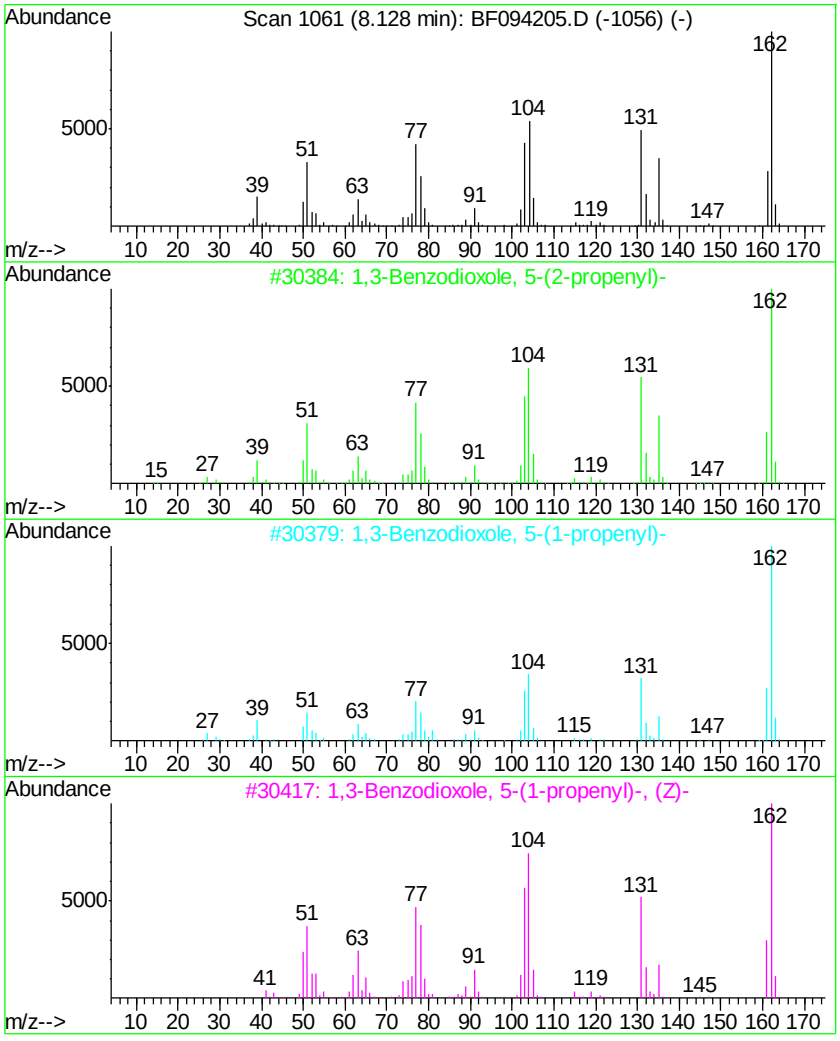
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

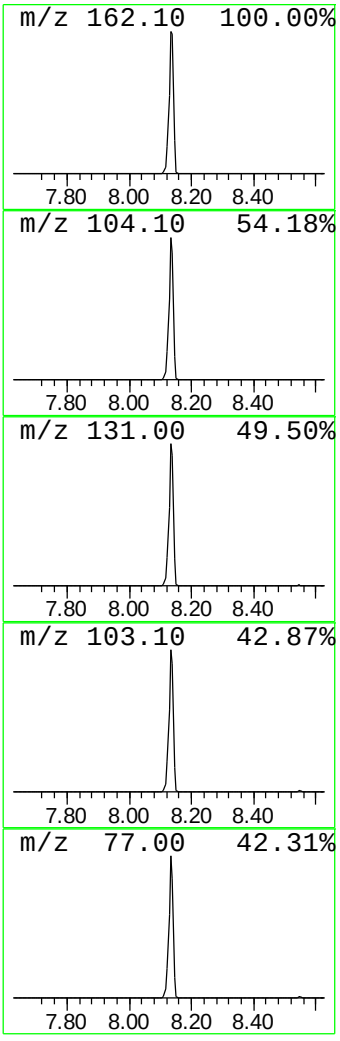
**********

Peak Number 5 1,3-Benzodioxole, 5-(2-prop... Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.13 | 32.89 ng | 2419620 | Naphthalene-d8 | 7.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,3-Benzodioxole, 5-(2-propenyl)- | 162 | C10H10O2 | 000094-59-7 | 98 |
| 2 | 1,3-Benzodioxole, 5-(1-propenyl)- | 162 | C10H10O2 | 000120-58-1 | 96 |
| 3 | 1,3-Benzodioxole, 5-(1-propenyl)... | 162 | C10H10O2 | 017627-76-8 | 94 |
| 4 | Benzene, 1,2-(methylenedioxy)-4-... | 162 | C10H10O2 | 004043-71-4 | 94 |
| 5 | 1,4-Dioxo-1,2,3,4-tetrahydrophth... | 162 | C8H6N2O2 | 001445-69-8 | 41 |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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

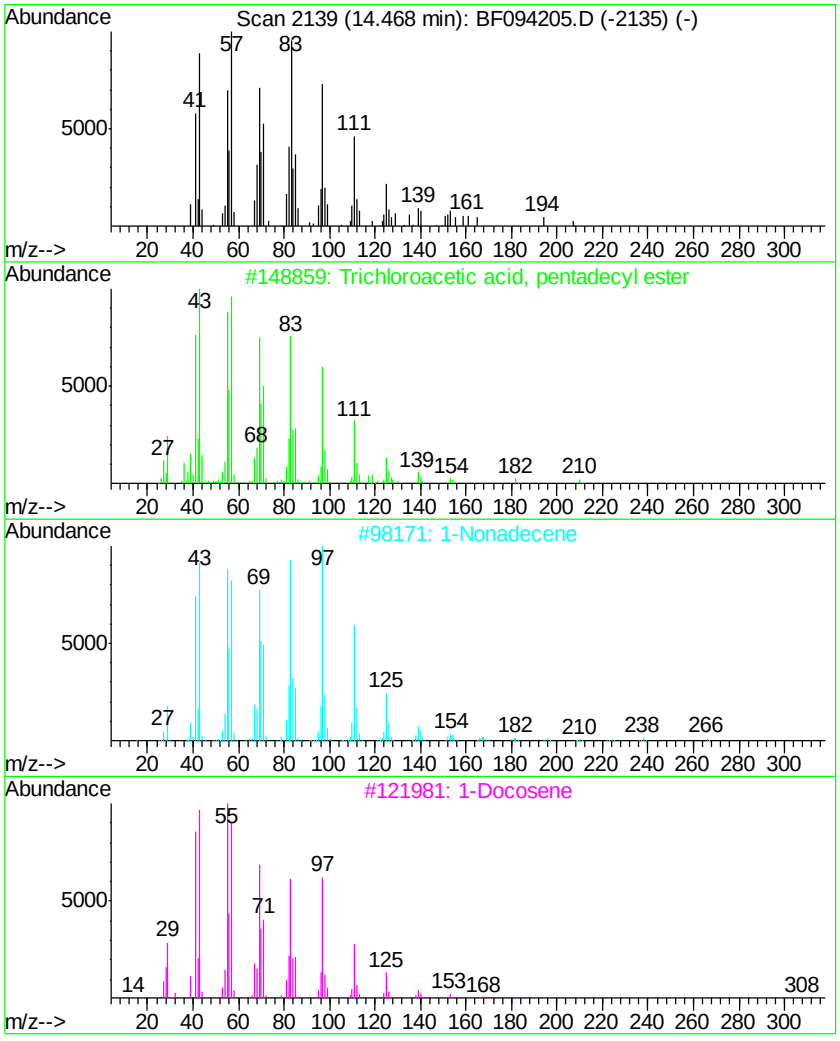
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

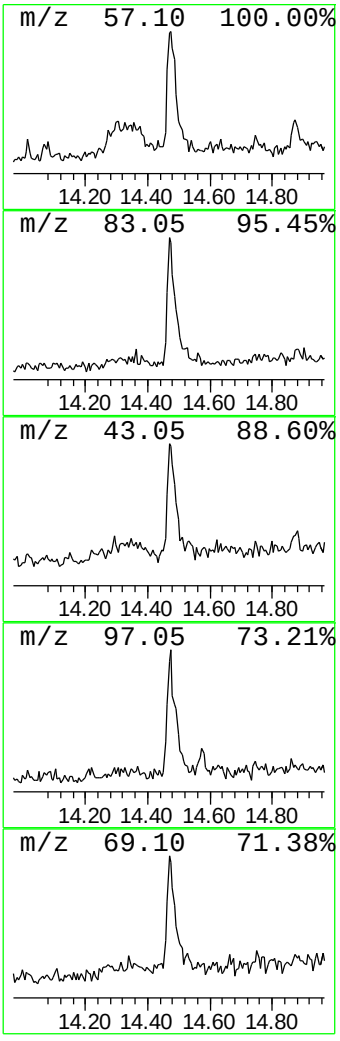
**********

Peak Number 6 Trichloroacetic acid, penta... Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.47 | 2.91 ng | 147127 | Chrysene-d12 | 14.57 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Trichloroacetic acid, pentadecyl... | 372 | C17H31Cl3O2 | 074339-53-0 | 91 |
| 2 | 1-Nonadecene | 266 | C19H38 | 018435-45-5 | 91 |
| 3 | 1-Docosene | 308 | C22H44 | 001599-67-3 | 91 |
| 4 | Dichloroacetic acid, tetradecyl ... | 324 | C16H30Cl2O2 | 083005-02-1 | 91 |
| 5 | 1-Eicosanol | 298 | C20H42O | 000629-96-9 | 91 |



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Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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

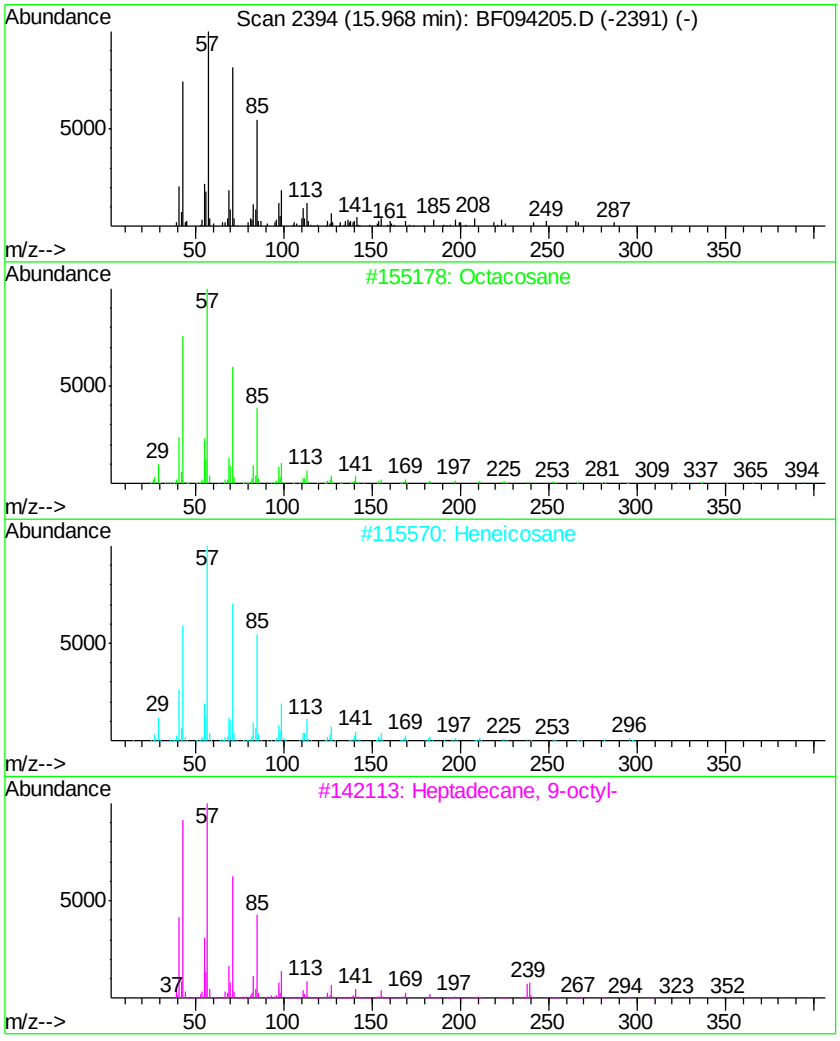
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

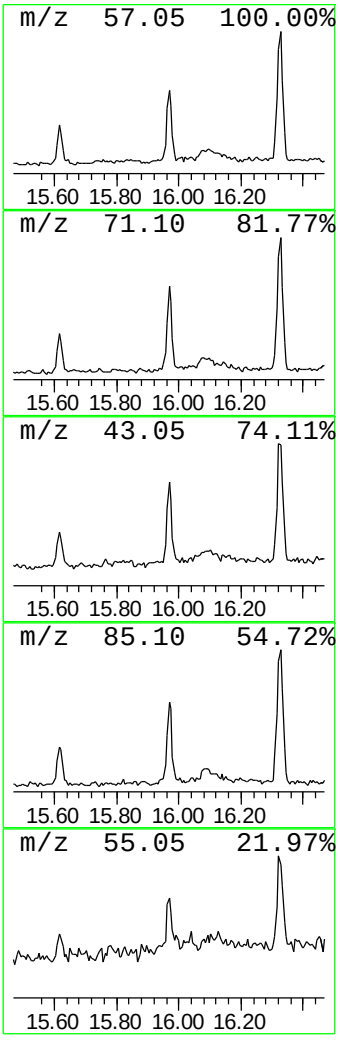
**********

Peak Number 7 Octacosane Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.97 | 2.02 ng | 92916 | Perylene-d12 | 16.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octacosane | 394 | C28H58 | 000630-02-4 | 90 |
| 2 | Heneicosane | 296 | C21H44 | 000629-94-7 | 87 |
| 3 | Heptadecane, 9-octyl- | 352 | C25H52 | 007225-64-1 | 87 |
| 4 | Nonadecane | 268 | C19H40 | 000629-92-5 | 86 |
| 5 | Eicosane, 7-hexyl- | 366 | C26H54 | 055333-99-8 | 86 |



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Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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

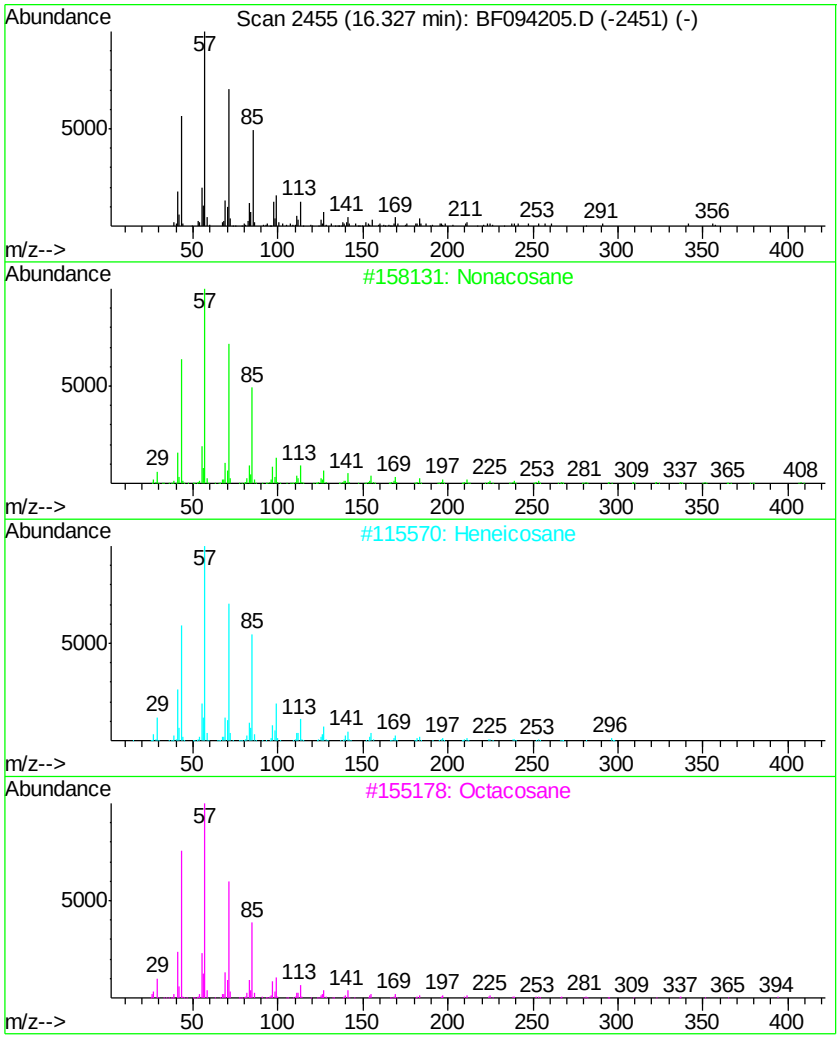
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

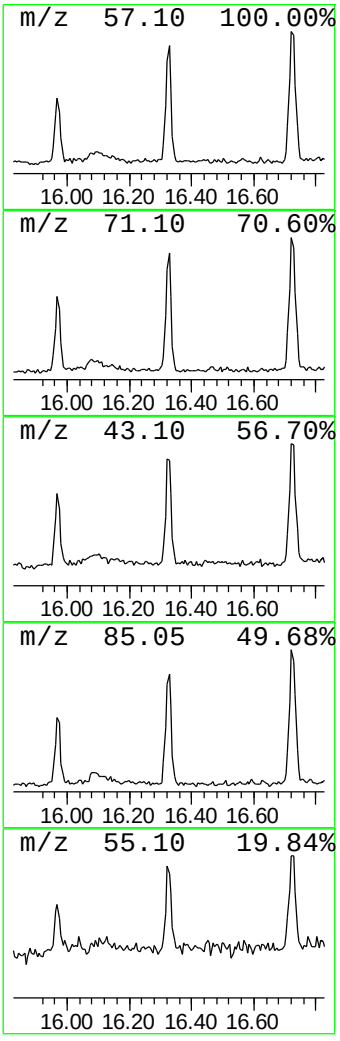
**********

Peak Number 8 Nonacosane Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.33 | 4.18 ng | 192229 | Perylene-d12 | 16.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Nonacosane | 408 | C29H60 | 000630-03-5 | 90 |
| 2 | Heneicosane | 296 | C21H44 | 000629-94-7 | 90 |
| 3 | Octacosane | 394 | C28H58 | 000630-02-4 | 87 |
| 4 | Tetracosane | 338 | C24H50 | 000646-31-1 | 87 |
| 5 | Docosane | 310 | C22H46 | 000629-97-0 | 87 |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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

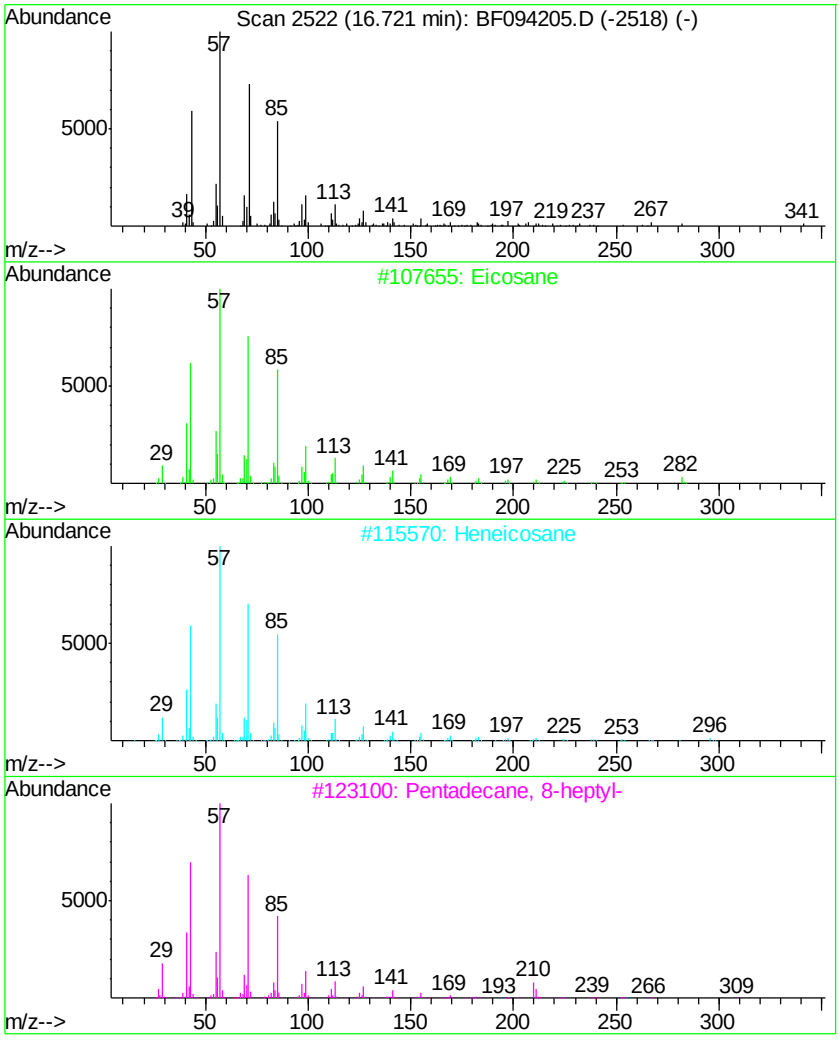
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

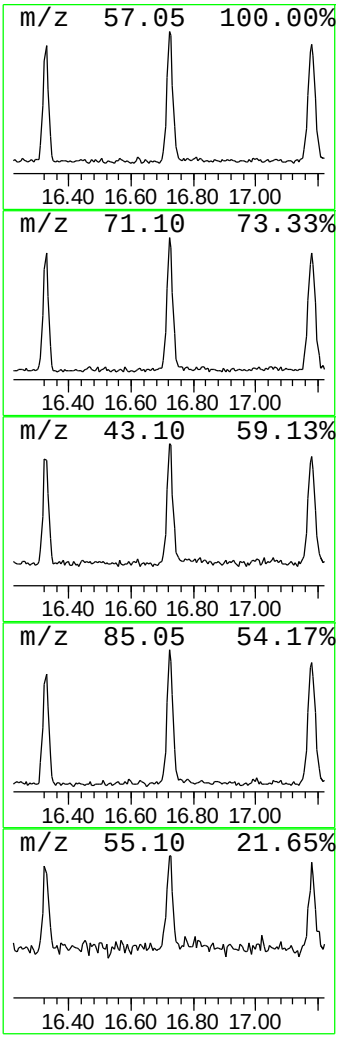
**********

Peak Number 9 Eicosane Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.72 | 5.27 ng | 242168 | Perylene-d12 | 16.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Eicosane | 282 | C20H42 | 000112-95-8 | 94 |
| 2 | Heneicosane | 296 | C21H44 | 000629-94-7 | 91 |
| 3 | Pentadecane, 8-heptyl- | 310 | C22H46 | 071005-15-7 | 91 |
| 4 | Octacosane | 394 | C28H58 | 000630-02-4 | 91 |
| 5 | Tetracosane | 338 | C24H50 | 000646-31-1 | 91 |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(0-2)

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

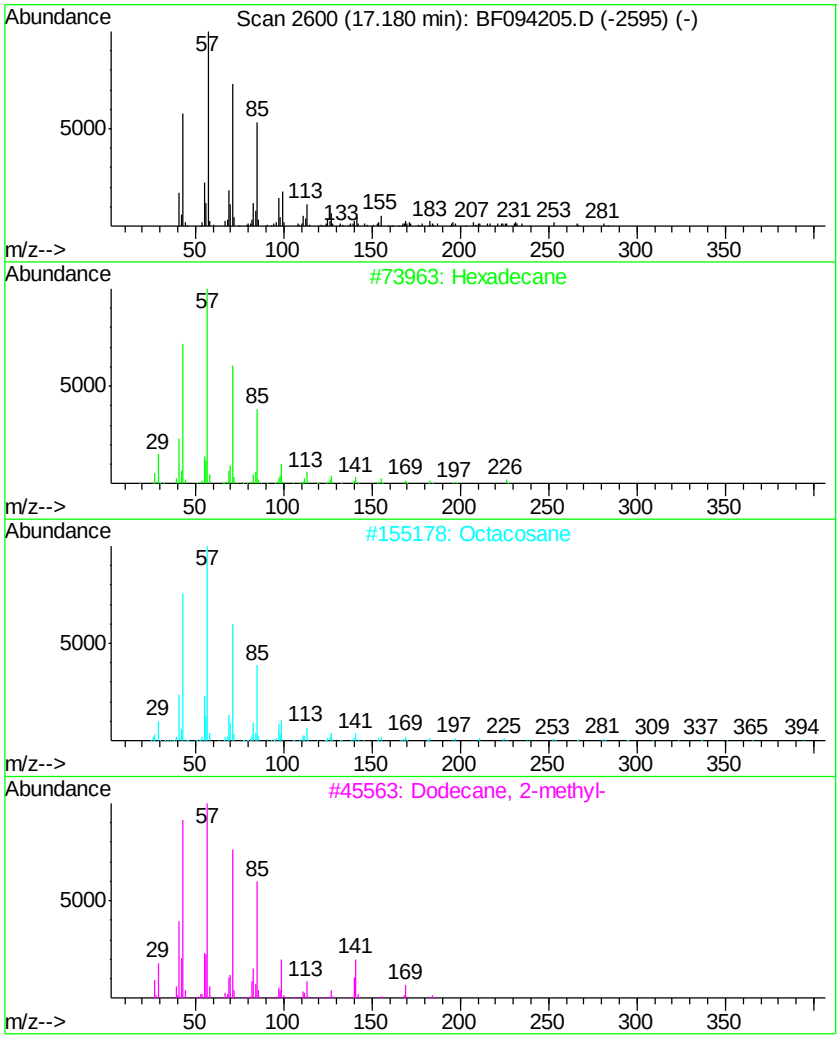
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

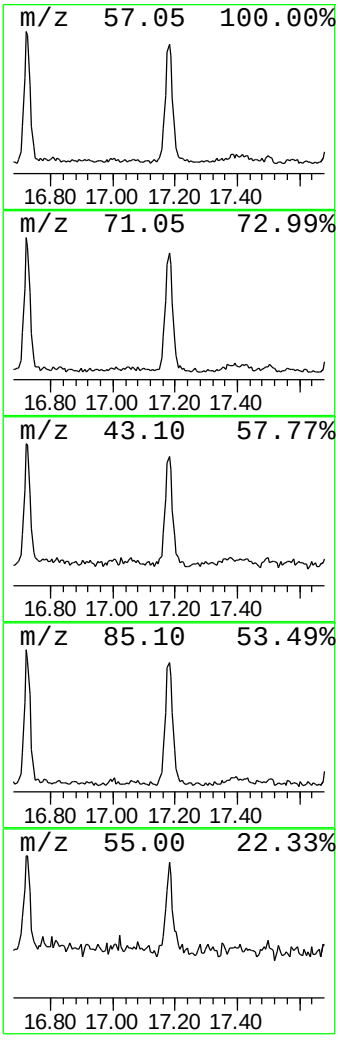
**********

Peak Number 10 Hexadecane Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.18 | 6.32 ng | 290768 | Perylene-d12 | 16.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexadecane | 226 | C16H34 | 000544-76-3 | 94 |
| 2 | Octacosane | 394 | C28H58 | 000630-02-4 | 90 |
| 3 | Dodecane, 2-methyl- | 184 | C13H28 | 001560-97-0 | 90 |
| 4 | Heneicosane | 296 | C21H44 | 000629-94-7 | 90 |
| 5 | Heptadecane, 9-octyl- | 352 | C25H52 | 007225-64-1 | 87 |



Data Path : Z:\HPCHEM1\BNA F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(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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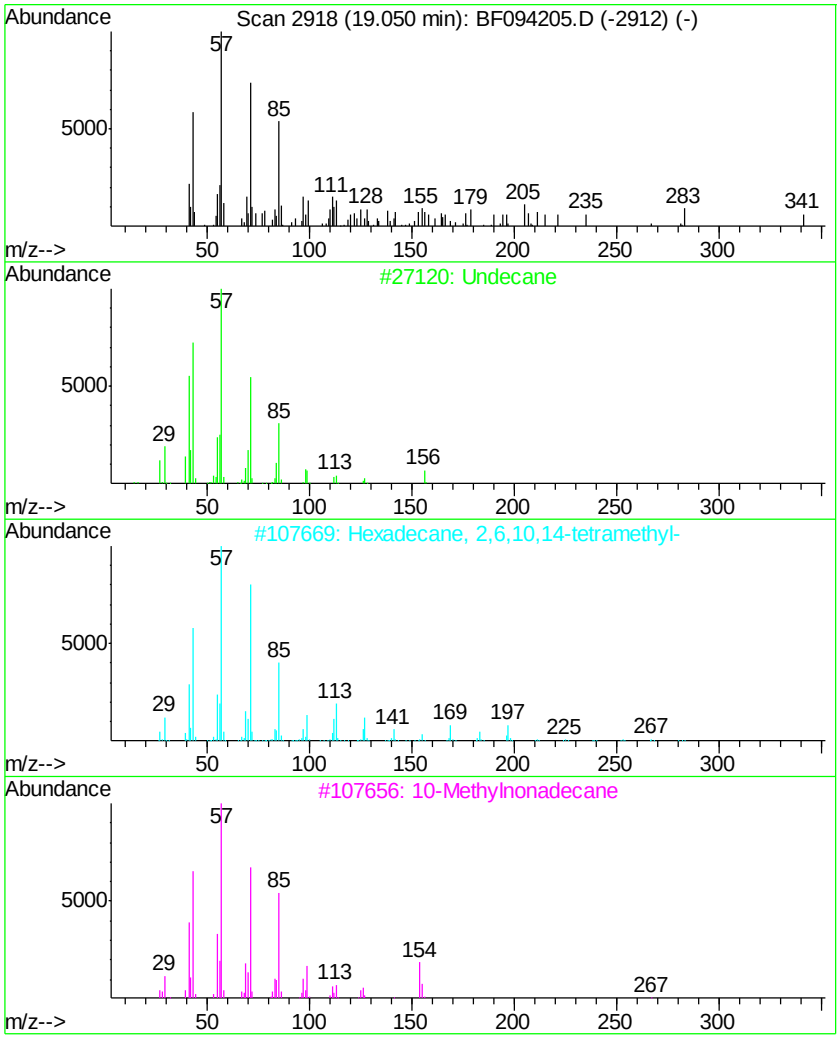
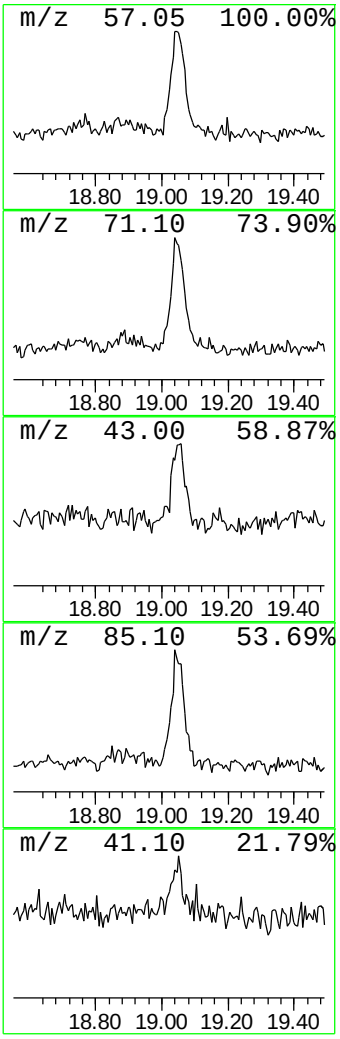
Peak Number 13 Undecane Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.05 | 2.07 ng | 95180 | Perylene-d12 | 16.20 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Undecane | 156 | C11H24 | 001120-21-4 | 68 |
| 2 | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42 | 000638-36-8 | 62 |
| 3 | 10-Methylnonadecane | 282 | C20H42 | 056862-62-5 | 59 |
| 4 | Octadecane, 2,6-dimethyl- | 282 | C20H42 | 075163-97-2 | 58 |
| 5 | Nonadecane | 268 | C19H40 | 000629-92-5 | 58 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\Data\BF040517\
Data File : BF094205.D
Acq On : 5 Apr 2017 22:33
Operator : SJ/MA
Sample : I2522-07
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BASING-4(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

| | | | | | Internal Standard | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| 2-Pentanone, 4-hy... | 4.00 | 11.5 | ng | 629816 | 1 | 5.86 | 1097640 | 20.0 |
| unknown5.61 | 5.61 | 75.5 | ng | 4144560 | 1 | 5.86 | 1097640 | 20.0 |
| Bicyclo[2.2.1]hep... | 7.06 | 3.0 | ng | 223355 | 2 | 7.36 | 1471360 | 20.0 |
| 1,3-Benzodioxole,... | 8.13 | 32.9 | ng | 2419620 | 2 | 7.36 | 1471360 | 20.0 |
| Trichloroacetic a... | 14.47 | 2.9 | ng | 147127 | 5 | 14.57 | 1011860 | 20.0 |
| Octacosane | 15.97 | 2.0 | ng | 92916 | 6 | 16.20 | 919784 | 20.0 |
| Nonacosane | 16.33 | 4.2 | ng | 192229 | 6 | 16.20 | 919784 | 20.0 |
| Eicosane | 16.72 | 5.3 | ng | 242168 | 6 | 16.20 | 919784 | 20.0 |
| Hexadecane | 17.18 | 6.3 | ng | 290768 | 6 | 16.20 | 919784 | 20.0 |
| Undecane | 19.05 | 2.1 | ng | 95180 | 6 | 16.20 | 919784 | 20.0 |