LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** 15-VE-PILE-WALL(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_G\METHODS\8270-BG060616.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 | 2.933 | 11 | 16 | 36 | rVB | 302153 | 502553 | 24.88% | 3.445% |
| 2 | 3.086 | 36 | 42 | 55 | rVB2 | 22269 | 55007 | 2.72% | 0.377% |
| 3 | 5.154 | 388 | 394 | 405 | rVB2 | 28645 | 55758 | 2.76% | 0.382% |
| 4 | 5.636 | 468 | 476 | 490 | rBV | 540658 | 1016888 | 50.34% | 6.971% |
| 5 | 7.257 | 742 | 752 | 768 | rBV | 620204 | 1206065 | 59.70% | 8.268% |
| 6 | 7.621 | 806 | 814 | 829 | rBV | 673432 | 1283815 | 63.55% | 8.801% |
| 7 | 8.086 | 886 | 893 | 901 | rBV | 111470 | 211529 | 10.47% | 1.450% |
| 8 | 8.415 | 941 | 949 | 970 | rVB | 510016 | 986997 | 48.86% | 6.766% |
| 9 | 9.261 | 1083 | 1093 | 1107 | rBV | 372798 | 753840 | 37.32% | 5.168% |
| 10 | 9.390 | 1111 | 1115 | 1131 | rVB5 | 9497 | 26087 | 1.29% | 0.179% |
| 11 | 10.906 | 1365 | 1373 | 1385 | rBV | 178191 | 365198 | 18.08% | 2.503% |
| 12 | 13.344 | 1780 | 1788 | 1805 | rBV | 1159387 | 2004834 | 99.24% | 13.743% |
| 13 | 14.167 | 1921 | 1928 | 1936 | rBV | 220653 | 350477 | 17.35% | 2.403% |
| 14 | 14.590 | 1995 | 2000 | 2006 | rVB | 16919 | 23910 | 1.18% | 0.164% |
| 15 | 14.725 | 2016 | 2023 | 2032 | rVB2 | 271218 | 485777 | 24.05% | 3.330% |
| 16 | 15.536 | 2157 | 2161 | 2167 | rVB2 | 26809 | 40438 | 2.00% | 0.277% |
| 17 | 16.211 | 2269 | 2276 | 2291 | rBV | 718268 | 1196483 | 59.23% | 8.202% |
| 18 | 16.399 | 2303 | 2308 | 2317 | rBV2 | 25445 | 49643 | 2.46% | 0.340% |
| 19 | 17.193 | 2437 | 2443 | 2448 | rBV | 17127 | 27575 | 1.37% | 0.189% |
| 20 | 17.469 | 2483 | 2490 | 2504 | rBV | 292142 | 508181 | 25.16% | 3.484% |
| 21 | 19.737 | 2871 | 2876 | 2881 | rBV | 144200 | 188946 | 9.35% | 1.295% |
| 22 | 19.813 | 2885 | 2889 | 2892 | rBV3 | 15107 | 22481 | 1.11% | 0.154% |
| 23 | 19.848 | 2892 | 2895 | 2903 | rVB2 | 37381 | 58907 | 2.92% | 0.404% |
| 24 | 20.060 | 2925 | 2931 | 2939 | rBV | 1374557 | 2020108 | 100.00% | 13.848% |
| 25 | 20.771 | 3048 | 3052 | 3059 | rVB | 101595 | 139310 | 6.90% | 0.955% |
| 26 | 20.835 | 3060 | 3063 | 3066 | rBV | 23632 | 29661 | 1.47% | 0.203% |
| 27 | 20.871 | 3067 | 3069 | 3073 | rVB3 | 19133 | 20742 | 1.03% | 0.142% |
| 28 | 21.341 | 3146 | 3149 | 3156 | rVB2 | 37690 | 63907 | 3.16% | 0.438% |
| 29 | 21.740 | 3211 | 3217 | 3229 | rVB | 234308 | 452310 | 22.39% | 3.101% |
| 30 | 21.893 | 3239 | 3243 | 3249 | rVB | 31878 | 48739 | 2.41% | 0.334% |
| 31 | 23.203 | 3461 | 3466 | 3474 | rVB4 | 26671 | 61148 | 3.03% | 0.419% |
| 32 | 25.019 | 3769 | 3775 | 3792 | rVB2 | 95815 | 330245 | 16.35% | 2.264% |

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
15-VE-PILE-WALL(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_G\METHODS\8270-BG060616.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Sum of corrected areas: 14587559

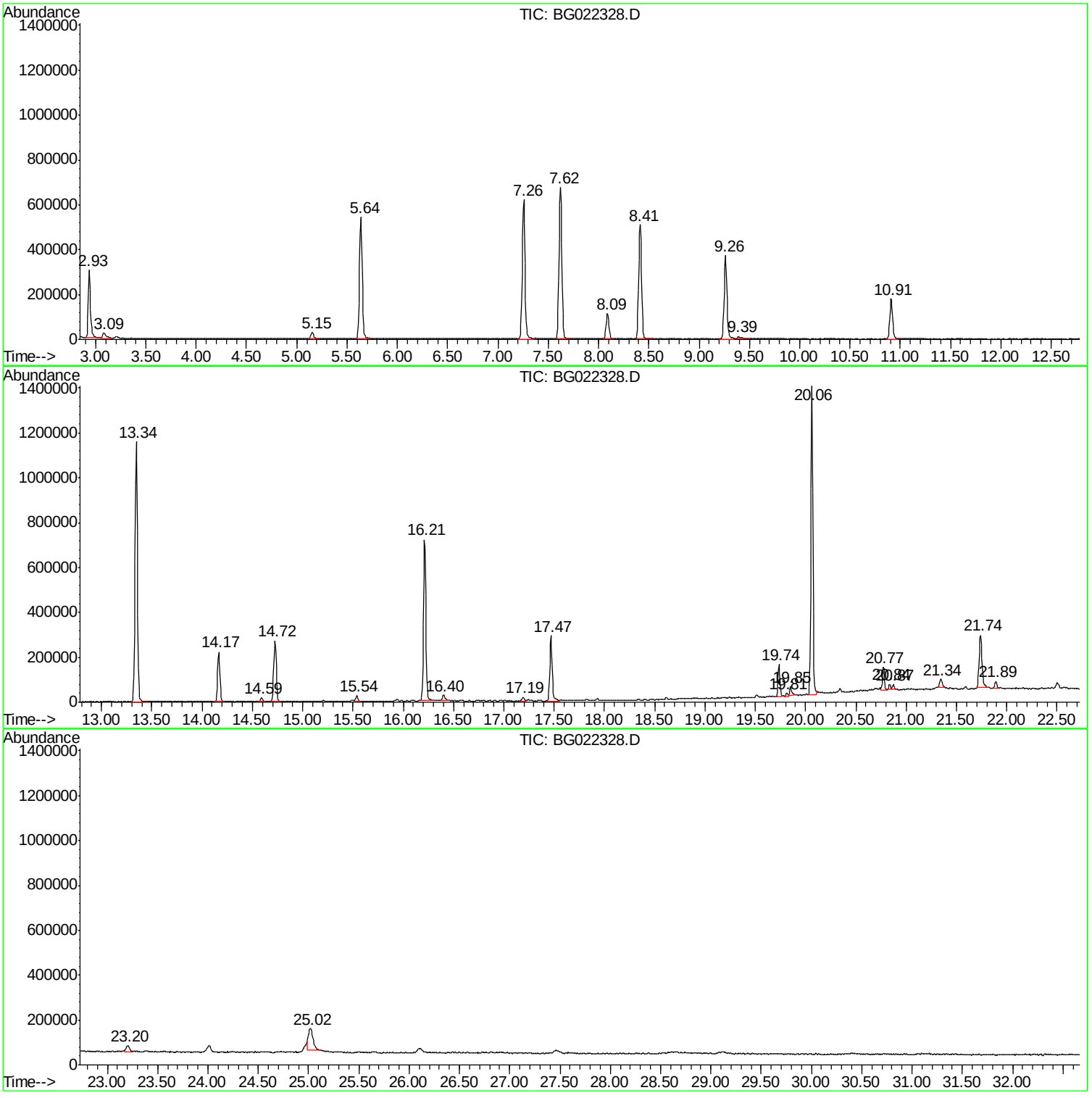
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

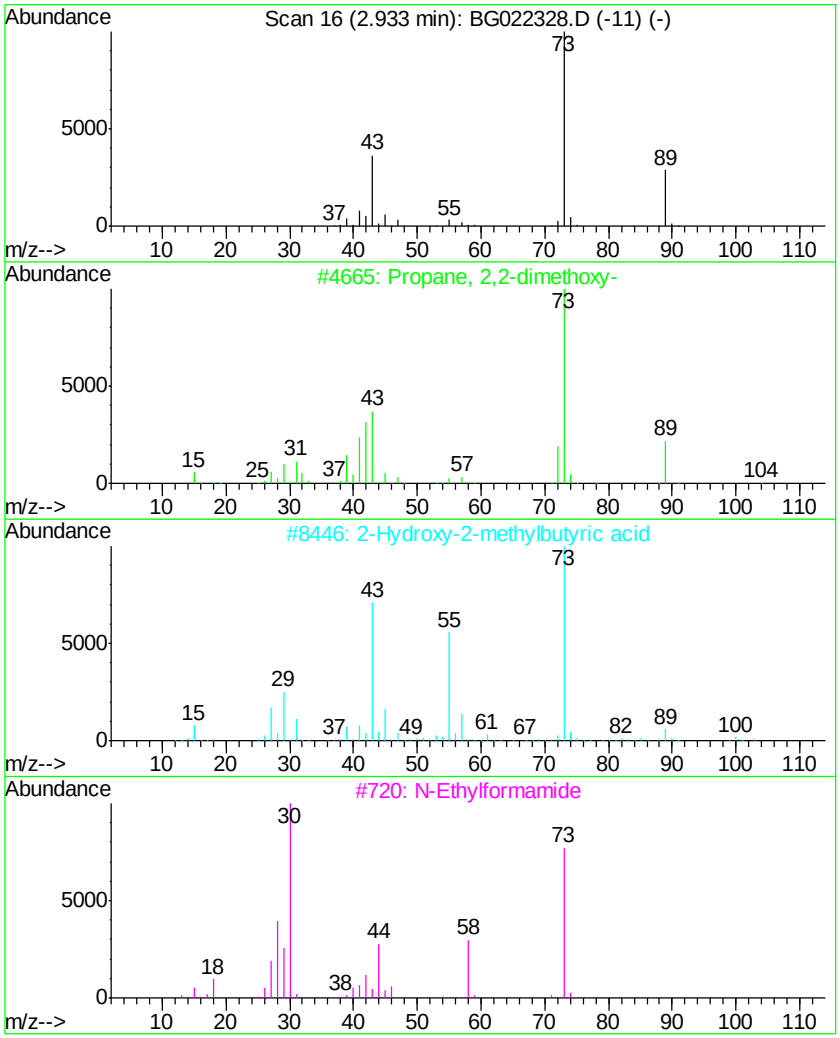
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

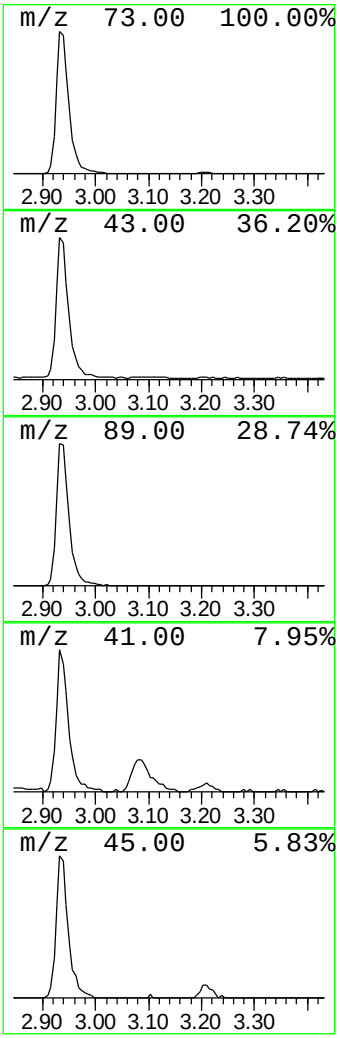
**********

Peak Number 1 unknown2.93 Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.93 | 47.52 ng | 502553 | 1,4-Dichlorobenzene-d4 | 8.09 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 2,2-dimethoxy- | 104 | C5H12O2 | 000077-76-9 | 45 |
| 2 | 2-Hydroxy-2-methylbutyric acid | 118 | C5H10O3 | 003739-30-8 | 9 |
| 3 | N-Ethylformamide | 73 | C3H7NO | 000627-45-2 | 5 |
| 4 | Silane, tetramethyl- | 88 | C4H12Si | 000075-76-3 | 4 |
| 5 | 1-Propene-1-thiol | 74 | C3H6S | 000925-89-3 | 4 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** 15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

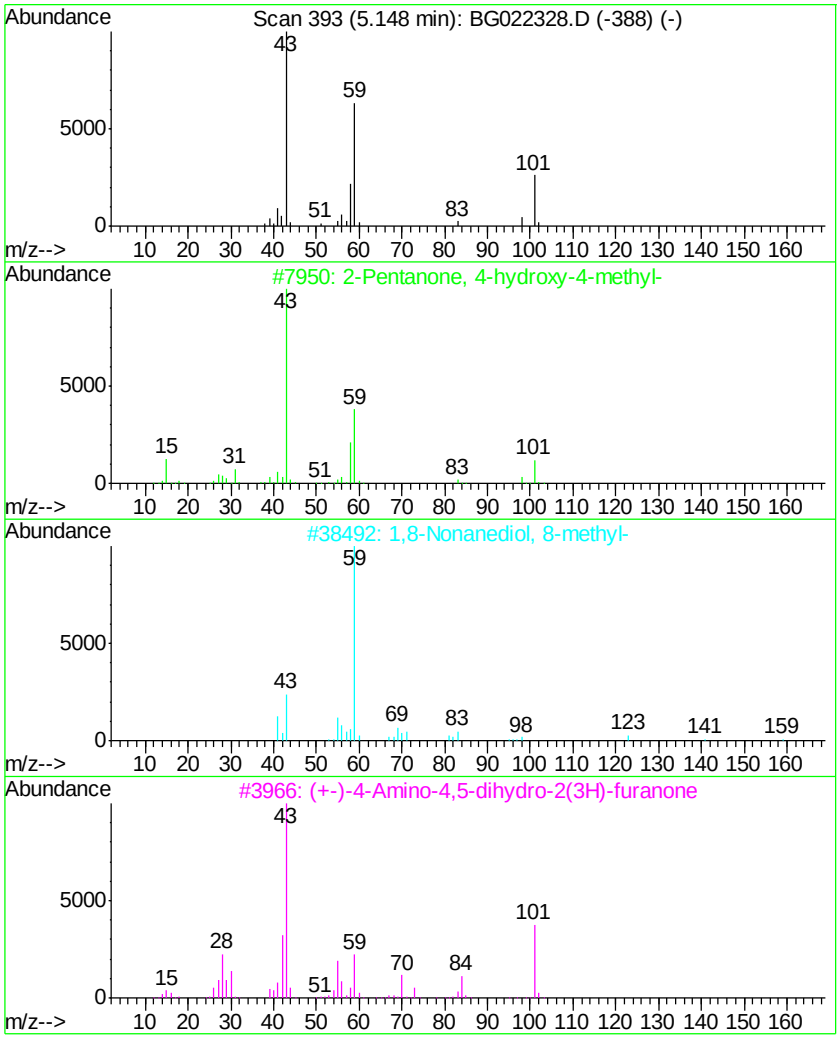
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

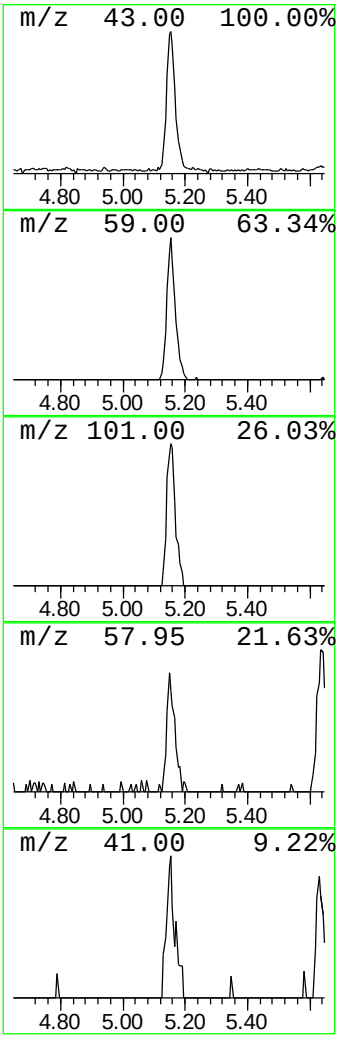
**********

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.15 | 5.27 ng | 55758 | 1,4-Dichlorobenzene-d4 | 8.09 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 56 |
| 2 | 1,8-Nonanediol, 8-methyl- | 174 | C10H22O2 | 054725-73-4 | 17 |
| 3 | (+-)-4-Amino-4,5-dihydro-2(3H)-f... | 101 | C4H7NO2 | 016504-58-8 | 9 |
| 4 | 2,3-Butanedione, monooxime | 101 | C4H7NO2 | 000057-71-6 | 9 |
| 5 | Butane, 1-ethoxy- | 102 | C6H14O | 000628-81-9 | 9 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** 15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

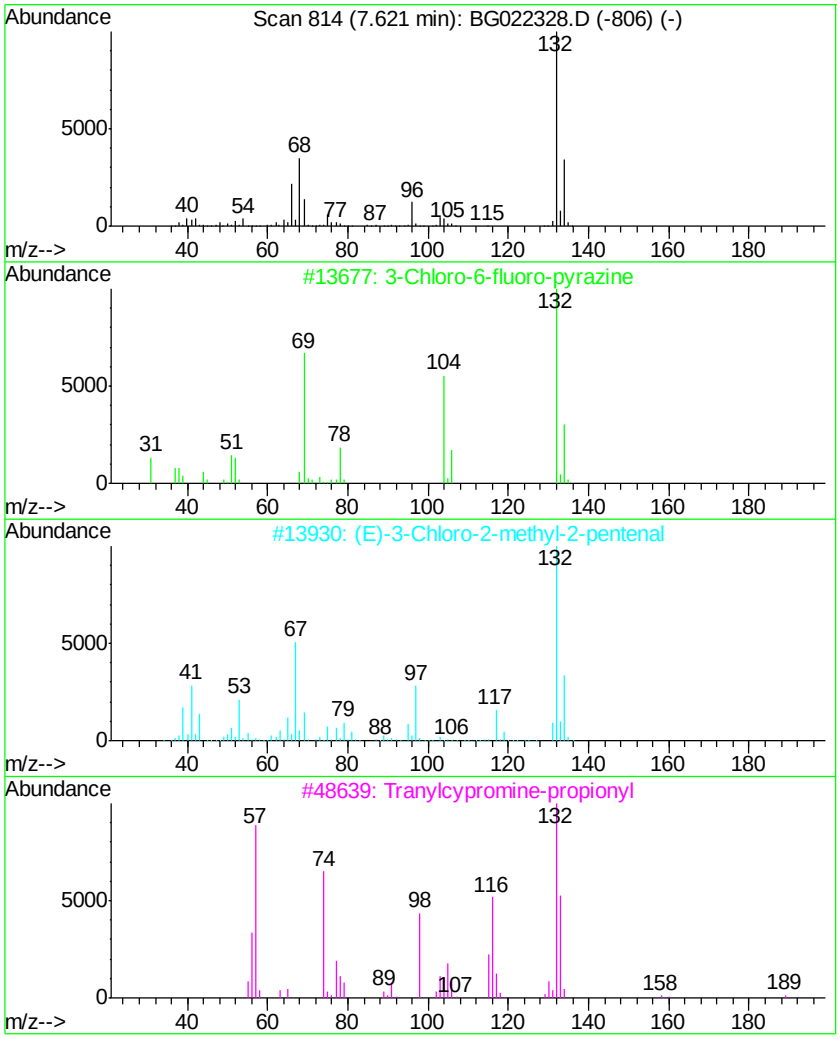
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

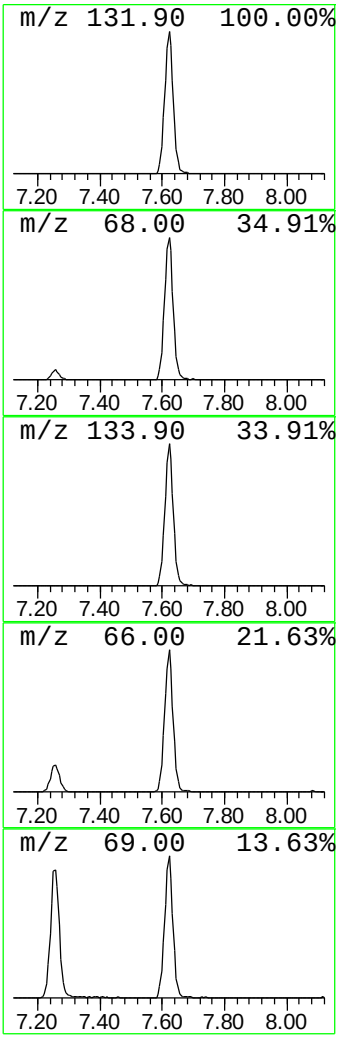
**********

Peak Number 4 unknown7.62 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.62 | 121.38 ng | 1283820 | 1,4-Dichlorobenzene-d4 | 8.09 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Chloro-6-fluoro-pyrazine | 132 | C4H2ClFN2 | 1000146-10-7 | 38 |
| 2 | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO | 031357-76-3 | 35 |
| 3 | Tranylcypromine-propionyl | 189 | C12H15NO | 1000123-86-3 | 12 |
| 4 | Amphetaminil | 250 | C17H18N2 | 017590-01-1 | 12 |
| 5 | Xenon | 132 | Xe | 007440-63-3 | 9 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** 15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

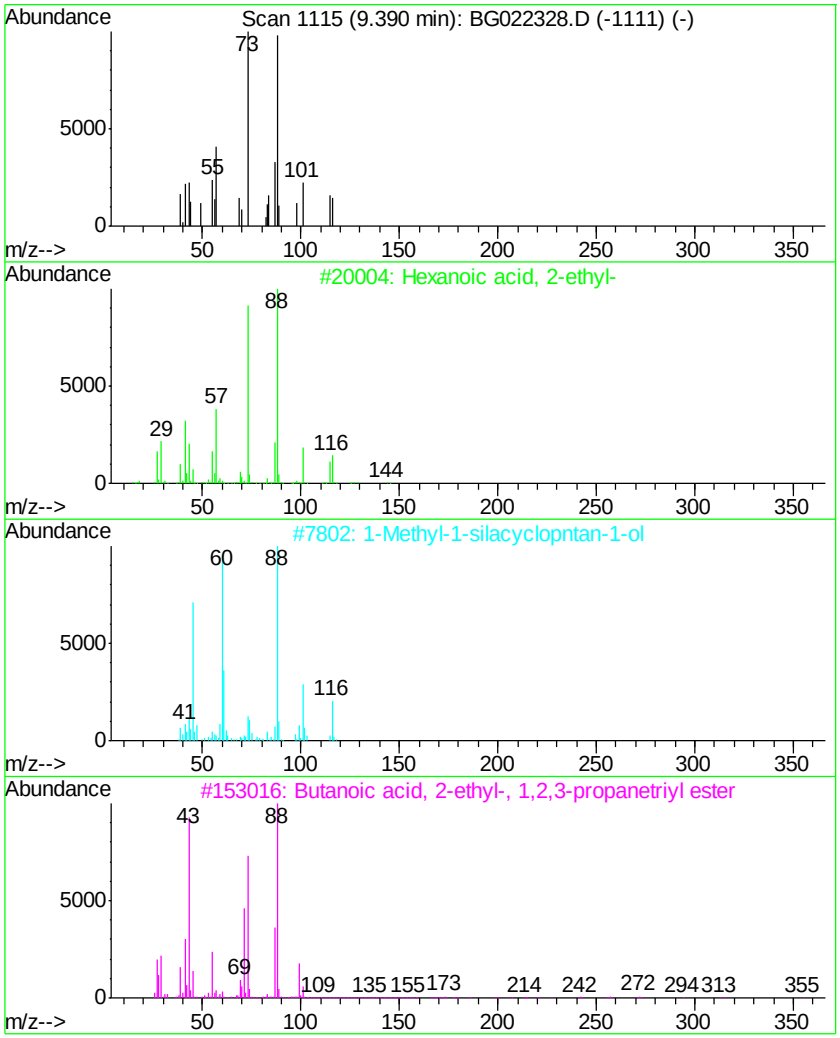
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

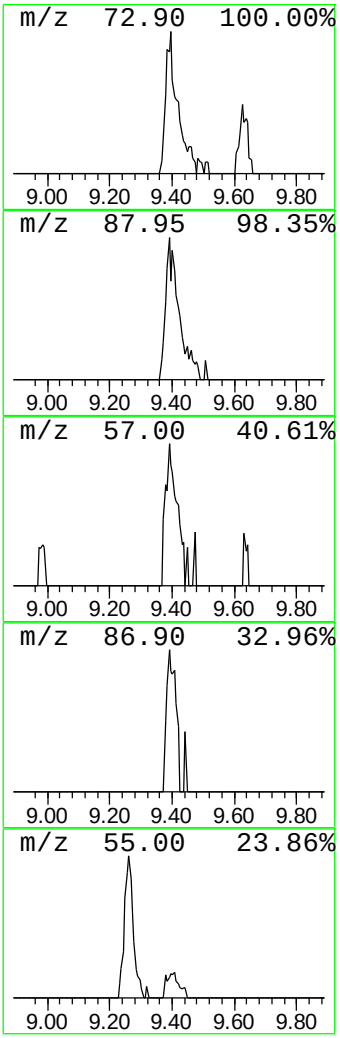
**********

Peak Number 6 Hexanoic acid, 2-ethyl- Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.39 | 2.47 ng | 26087 | 1,4-Dichlorobenzene-d4 | 8.09 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexanoic acid, 2-ethyl- | 144 | C8H16O2 | 000149-57-5 | 64 |
| 2 | 1-Methyl-1-silacyclopntan-1-ol | 116 | C5H12OSi | 1000216-84-4 | 50 |
| 3 | Butanoic acid, 2-ethyl-, 1,2,3-p... | 386 | C21H38O6 | 056554-54-2 | 33 |
| 4 | Pentanoic acid, 2-methyl-, methy... | 130 | C7H14O2 | 002177-77-7 | 28 |
| 5 | Hexanamide, 6-(heptyloxy)- | 229 | C13H27NO2 | 055191-08-7 | 25 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** 15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

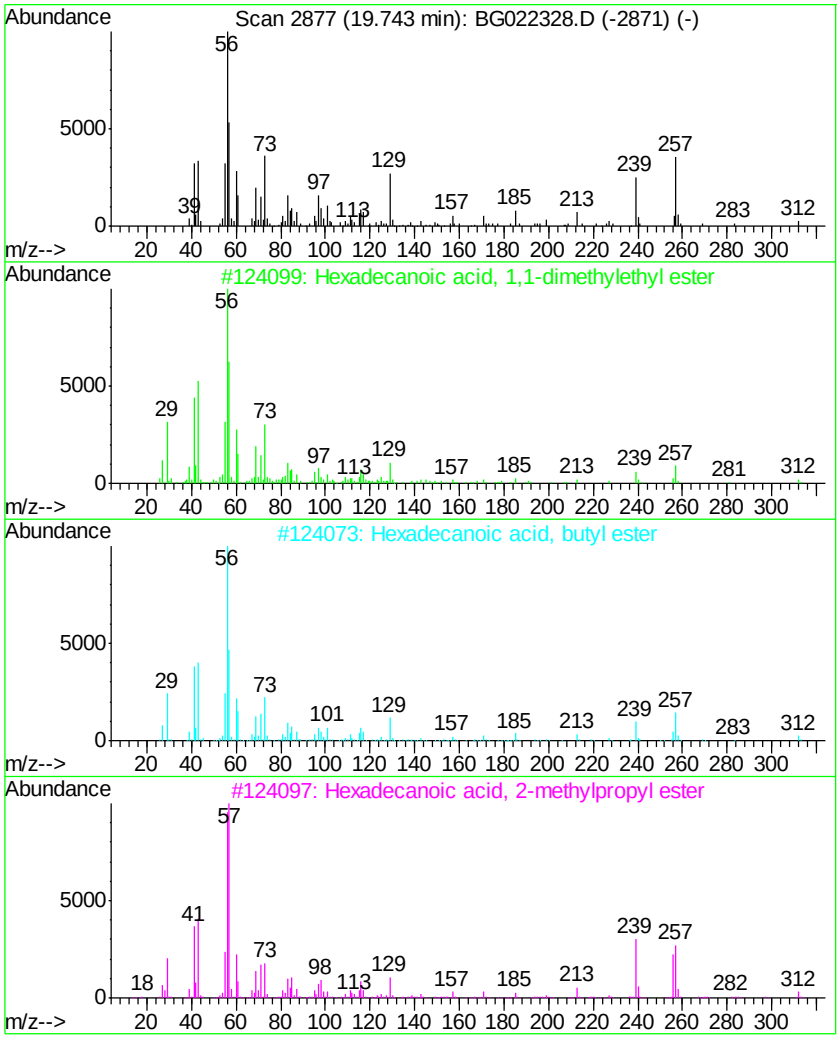
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

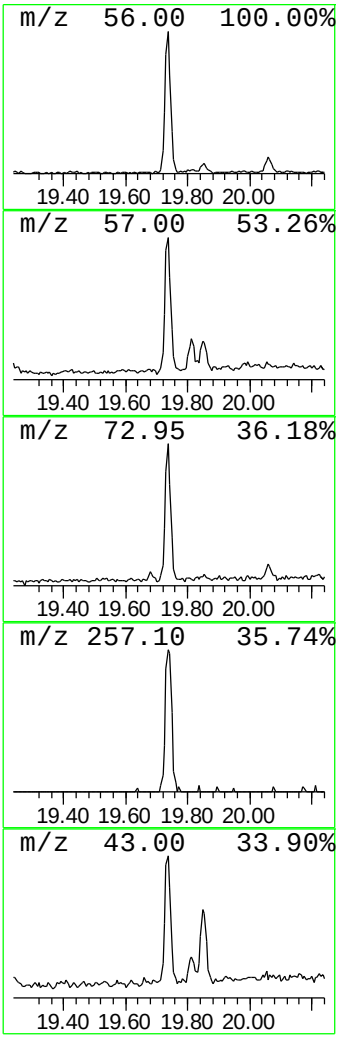
**********

Peak Number 7 Hexadecanoic acid, 1,1-dime... Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.74 | 8.35 ng | 188946 | Chrysene-d12 | 21.74 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexadecanoic acid, 1,1-dimethyle... | 312 | C20H40O2 | 031158-91-5 | 98 |
| 2 | Hexadecanoic acid, butyl ester | 312 | C20H40O2 | 000111-06-8 | 96 |
| 3 | Hexadecanoic acid, 2-methylpropy... | 312 | C20H40O2 | 000110-34-9 | 43 |
| 4 | Cyclobutane, 1,1-dimethyl-2-octyl- | 196 | C14H28 | 062338-30-1 | 35 |
| 5 | Oxirane, 2,3-bis(1-methylethyl)-... | 128 | C8H16O | 054644-32-5 | 35 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** 15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

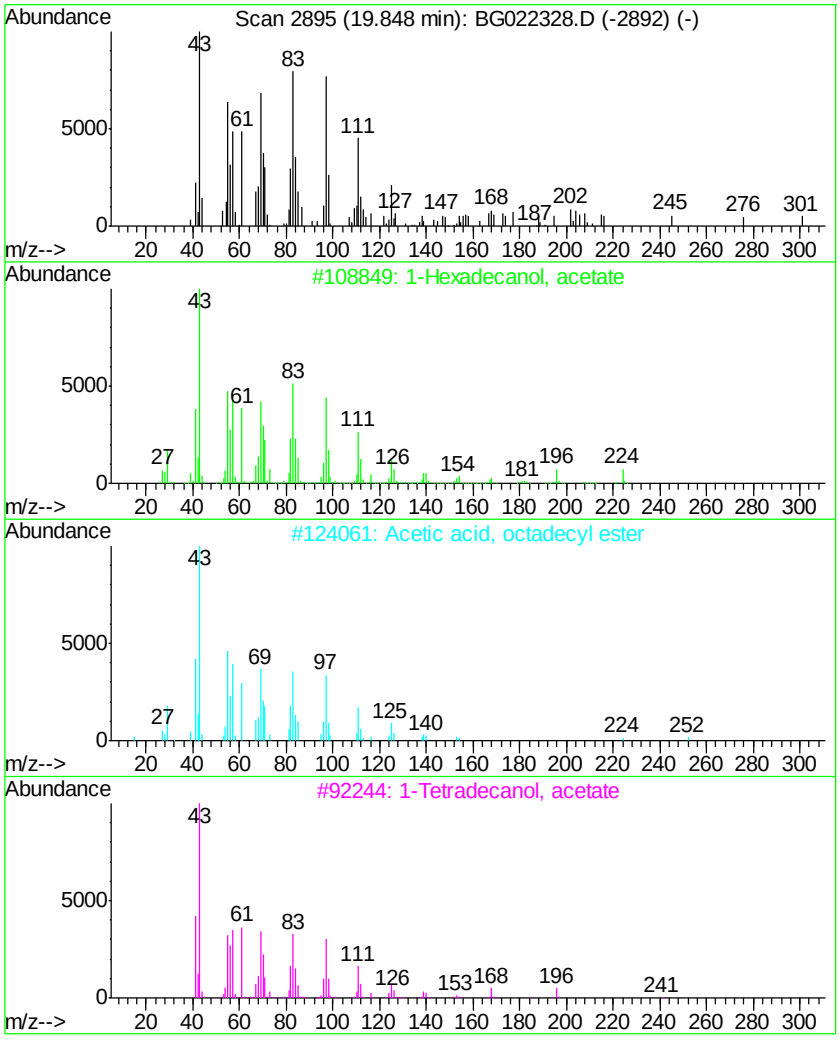
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

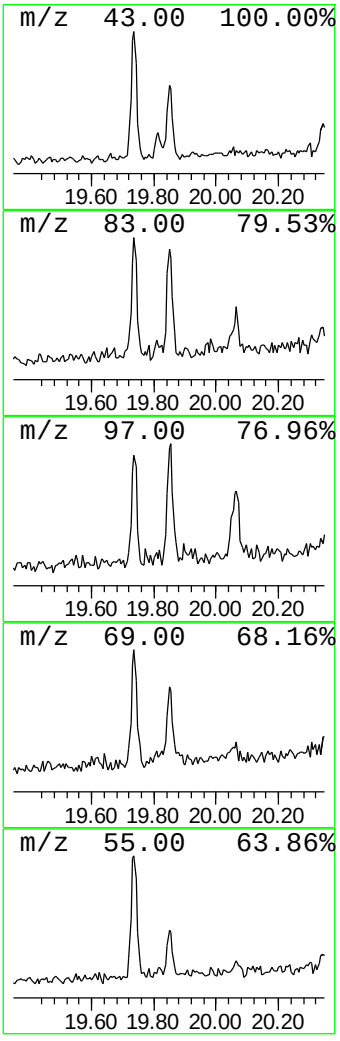
**********

Peak Number 8 1-Hexadecanol, acetate Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.85 | 2.60 ng | 58907 | Chrysene-d12 | 21.74 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Hexadecanol, acetate | 284 | C18H36O2 | 000629-70-9 | 90 |
| 2 | Acetic acid, octadecyl ester | 312 | C20H40O2 | 000822-23-1 | 87 |
| 3 | 1-Tetradecanol, acetate | 256 | C16H32O2 | 000638-59-5 | 86 |
| 4 | 1-Nonadecene | 266 | C19H38 | 018435-45-5 | 81 |
| 5 | 1-Hexadecanol | 242 | C16H34O | 036653-82-4 | 81 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

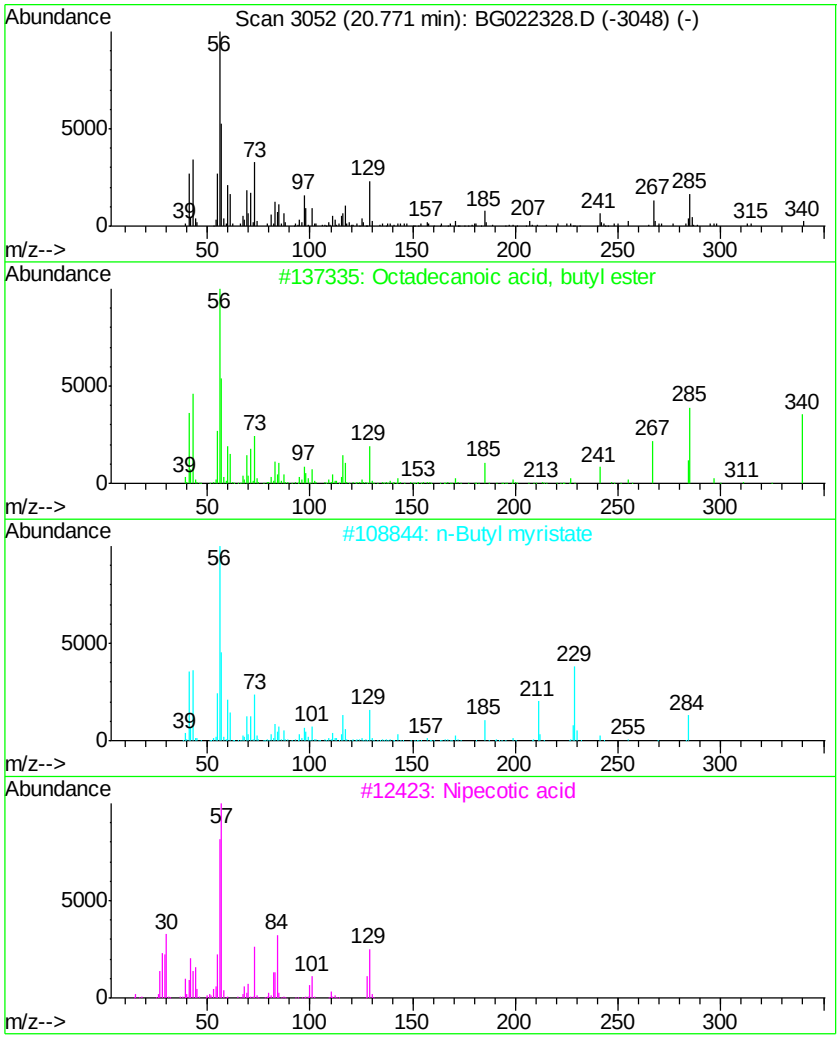
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

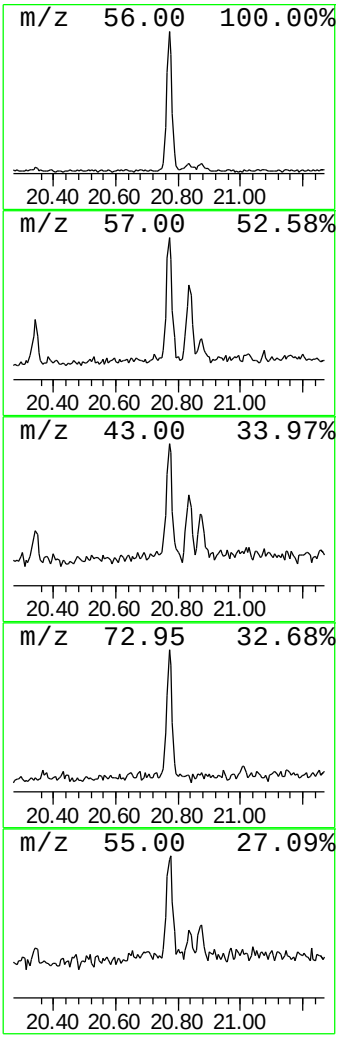
**********

Peak Number 9 Octadecanoic acid, butyl ester Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.77 | 6.16 ng | 139310 | Chrysene-d12 | 21.74 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecanoic acid, butyl ester | 340 | C22H44O2 | 000123-95-5 | 95 |
| 2 | n-Butyl myristate | 284 | C18H36O2 | 000110-36-1 | 59 |
| 3 | Nipecotic acid | 129 | C6H11NO2 | 000498-95-3 | 50 |
| 4 | 2,2-Dimethyl-1,3-butanediol | 118 | C6H14O2 | 000076-35-7 | 35 |
| 5 | 1-Hexene, 2,5-dimethyl- | 112 | C8H16 | 006975-92-4 | 30 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** 15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

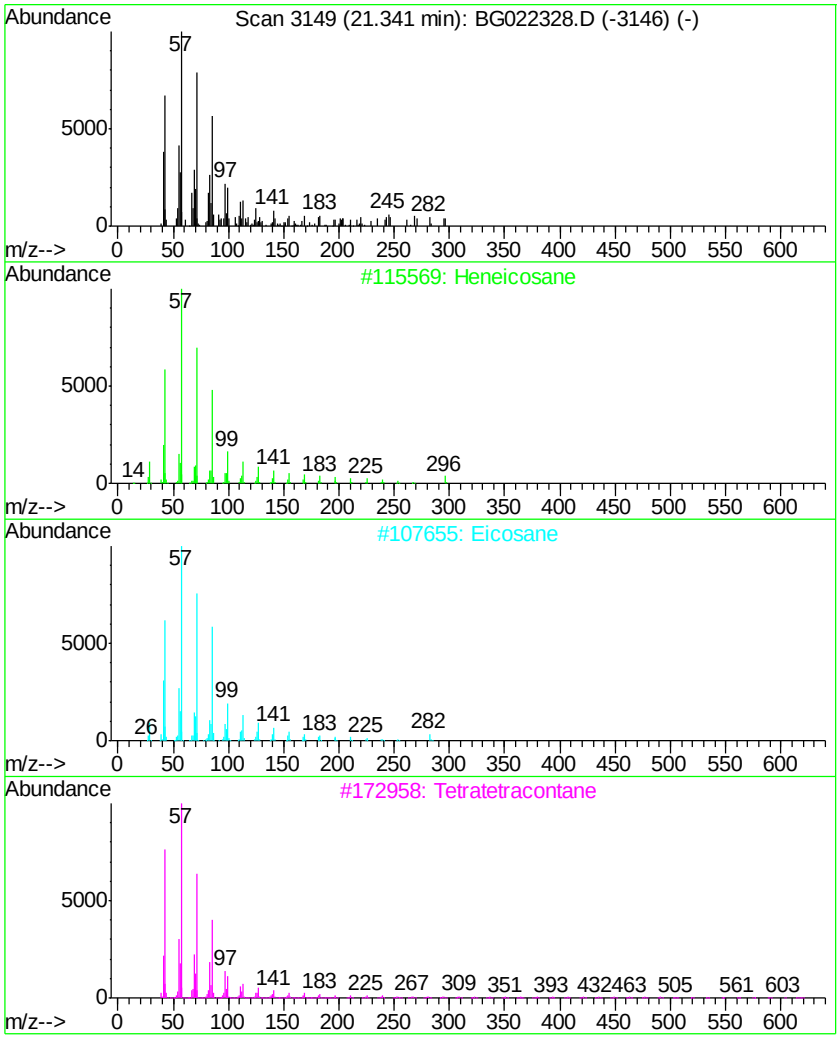
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

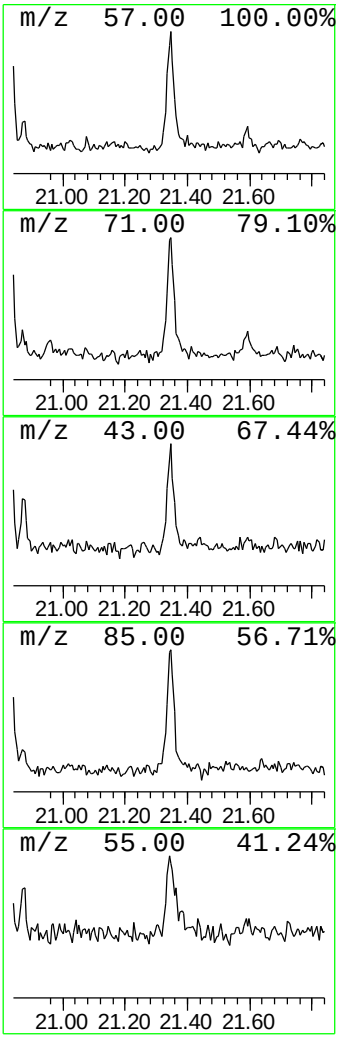
**********

Peak Number 10 Heneicosane Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.34 | 2.83 ng | 63907 | Chrysene-d12 | 21.74 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heneicosane | 296 | C21H44 | 000629-94-7 | 93 |
| 2 | Eicosane | 282 | C20H42 | 000112-95-8 | 89 |
| 3 | Tetratetracontane | 619 | C44H90 | 007098-22-8 | 80 |
| 4 | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44 | 054833-48-6 | 68 |
| 5 | Tetratriacontane | 479 | C34H70 | 014167-59-0 | 64 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** 15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

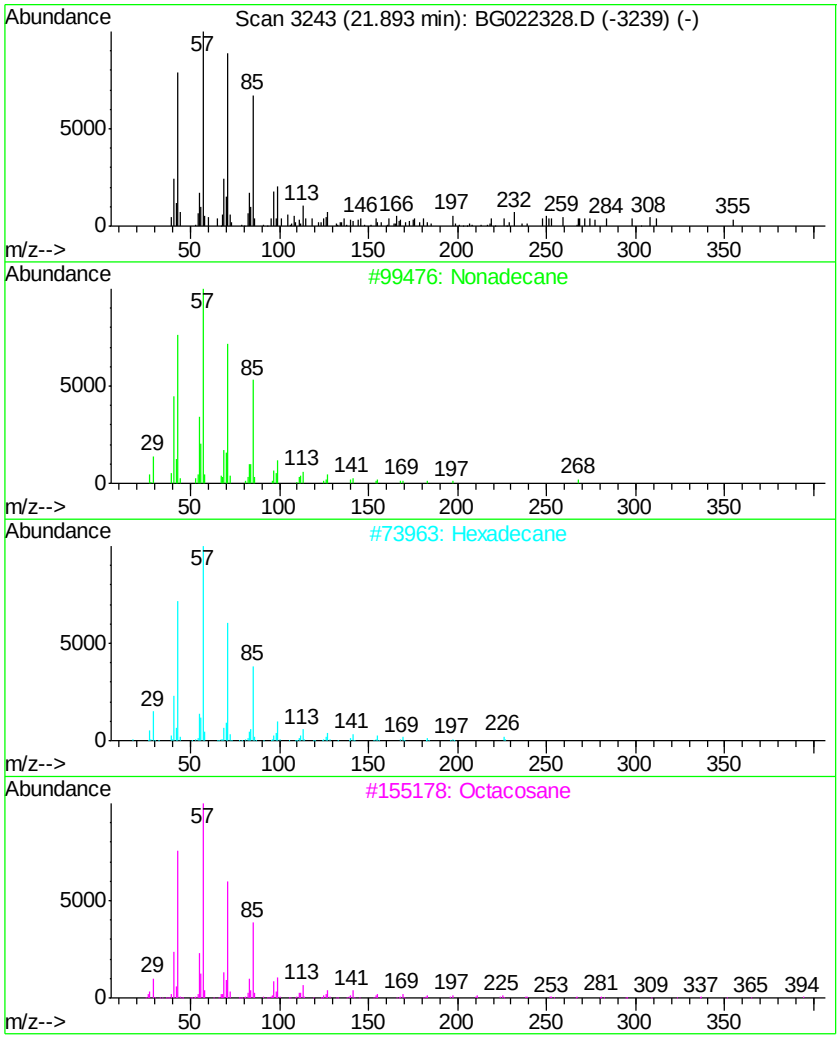
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

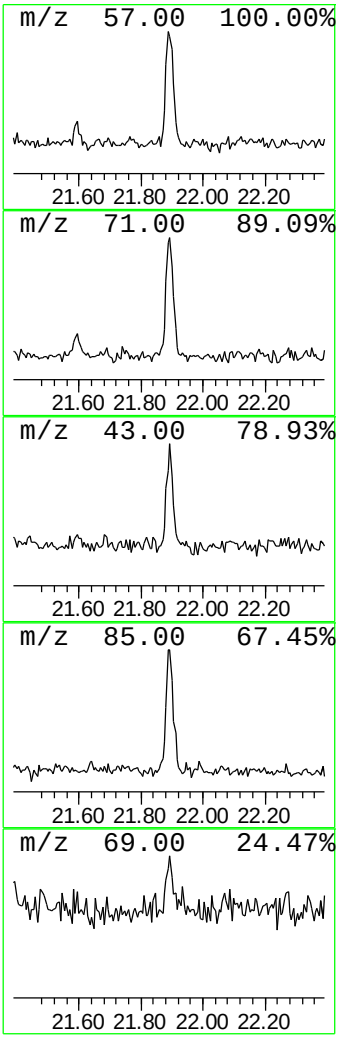
**********

Peak Number 11 Nonadecane Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.89 | 2.16 ng | 48739 | Chrysene-d12 | 21.74 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Nonadecane | 268 | C19H40 | 000629-92-5 | 89 |
| 2 | Hexadecane | 226 | C16H34 | 000544-76-3 | 81 |
| 3 | Octacosane | 394 | C28H58 | 000630-02-4 | 64 |
| 4 | Pentadecane | 212 | C15H32 | 000629-62-9 | 64 |
| 5 | Dodecane, 1-iodo- | 296 | C12H25I | 004292-19-7 | 59 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** 15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

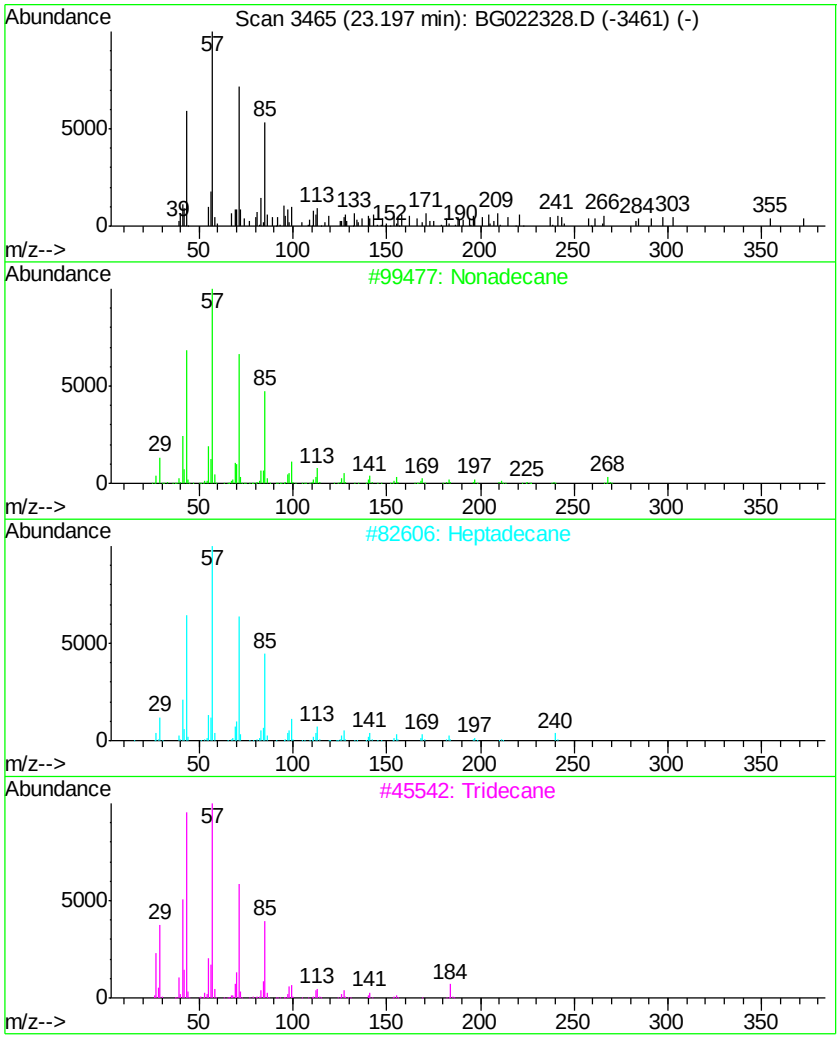
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

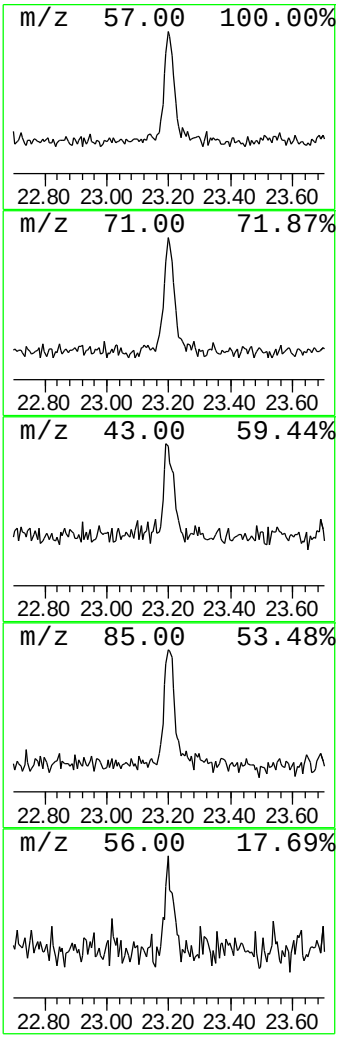
**********

Peak Number 12 Tridecane Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 23.20 | 2.70 ng | 61148 | Chrysene-d12 | 21.74 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Nonadecane | 268 | C19H40 | 000629-92-5 | 64 |
| 2 | Heptadecane | 240 | C17H36 | 000629-78-7 | 58 |
| 3 | Tridecane | 184 | C13H28 | 000629-50-5 | 58 |
| 4 | Pentadecane, 8-heptyl- | 310 | C22H46 | 071005-15-7 | 52 |
| 5 | Decane, 1-iodo- | 268 | C10H21I | 002050-77-3 | 50 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061316\
Data File : BG022328.D
Acq On : 13 Jun 2016 17:37
Operator : UM/SJ
Sample : H3448-29
Misc :
ALS Vial : 39 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** 15-VE-PILE-WALL(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.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 |
| unknown2.93 | 2.93 | 47.5 | ng | 502553 | 1 | 8.09 | 211529 | 20.0 |
| 2-Pentanone, 4-hy... | 5.15 | 5.3 | ng | 55758 | 1 | 8.09 | 211529 | 20.0 |
| unknown7.62 | 7.62 | 121.4 | ng | 1283820 | 1 | 8.09 | 211529 | 20.0 |
| Hexanoic acid, 2-... | 9.39 | 2.5 | ng | 26087 | 1 | 8.09 | 211529 | 20.0 |
| Hexadecanoic acid... | 19.74 | 8.3 | ng | 188946 | 5 | 21.74 | 452310 | 20.0 |
| 1-Hexadecanol, ac... | 19.85 | 2.6 | ng | 58907 | 5 | 21.74 | 452310 | 20.0 |
| Octadecanoic acid... | 20.77 | 6.2 | ng | 139310 | 5 | 21.74 | 452310 | 20.0 |
| Heneicosane | 21.34 | 2.8 | ng | 63907 | 5 | 21.74 | 452310 | 20.0 |
| Nonadecane | 21.89 | 2.2 | ng | 48739 | 5 | 21.74 | 452310 | 20.0 |
| Tridecane | 23.20 | 2.7 | ng | 61148 | 5 | 21.74 | 452310 | 20.0 |