LSC Area Percent Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110217\
Data File : BF100094.D
Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

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-BF102317.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 | 4.987 | 490 | 493 | 511 | rBV | 133731 | 220379 | 13.45% | 1.383% |
| 2 | 5.393 | 559 | 562 | 590 | rBV | 1211733 | 1447043 | 88.31% | 9.083% |
| 3 | 6.416 | 733 | 736 | 754 | rBV | 1368760 | 1455995 | 88.85% | 9.139% |
| 4 | 6.540 | 754 | 757 | 764 | rVB | 1559472 | 1521949 | 92.88% | 9.553% |
| 5 | 6.769 | 793 | 796 | 805 | rBV | 482207 | 438008 | 26.73% | 2.749% |
| 6 | 6.922 | 819 | 822 | 834 | rBV | 1323746 | 1200648 | 73.27% | 7.536% |
| 7 | 7.340 | 890 | 893 | 910 | rBV | 964239 | 901266 | 55.00% | 5.657% |
| 8 | 8.051 | 1011 | 1014 | 1025 | rBV | 713016 | 621998 | 37.96% | 3.904% |
| 9 | 9.122 | 1192 | 1196 | 1200 | rBV | 1924749 | 1638620 | 100.00% | 10.285% |
| 10 | 9.522 | 1261 | 1264 | 1271 | rVB | 276122 | 221428 | 13.51% | 1.390% |
| 11 | 9.669 | 1284 | 1289 | 1292 | rBV2 | 19393 | 23614 | 1.44% | 0.148% |
| 12 | 9.781 | 1305 | 1308 | 1310 | rBV | 471296 | 373857 | 22.82% | 2.347% |
| 13 | 9.804 | 1310 | 1312 | 1316 | rVB | 796713 | 624047 | 38.08% | 3.917% |
| 14 | 9.975 | 1338 | 1341 | 1344 | rVB | 164223 | 120410 | 7.35% | 0.756% |
| 15 | 10.057 | 1351 | 1355 | 1358 | rBV | 151263 | 133977 | 8.18% | 0.841% |
| 16 | 10.222 | 1380 | 1383 | 1386 | rVB | 24637 | 22306 | 1.36% | 0.140% |
| 17 | 10.522 | 1431 | 1434 | 1437 | rVB2 | 22839 | 21380 | 1.30% | 0.134% |
| 18 | 10.598 | 1444 | 1447 | 1457 | rBV | 915618 | 885384 | 54.03% | 5.557% |
| 19 | 10.692 | 1460 | 1463 | 1470 | rBV3 | 59051 | 62610 | 3.82% | 0.393% |
| 20 | 10.904 | 1493 | 1499 | 1502 | rBV6 | 12735 | 22311 | 1.36% | 0.140% |
| 21 | 11.145 | 1537 | 1540 | 1543 | rBV2 | 28759 | 30439 | 1.86% | 0.191% |
| 22 | 11.298 | 1562 | 1566 | 1568 | rBV | 656040 | 577536 | 35.25% | 3.625% |
| 23 | 11.322 | 1568 | 1570 | 1574 | rVB | 94956 | 82991 | 5.06% | 0.521% |
| 24 | 11.380 | 1577 | 1580 | 1583 | rBV | 25454 | 26934 | 1.64% | 0.169% |
| 25 | 11.810 | 1649 | 1653 | 1655 | rBV3 | 54761 | 67756 | 4.13% | 0.425% |
| 26 | 11.916 | 1668 | 1671 | 1673 | rVB2 | 35895 | 33555 | 2.05% | 0.211% |
| 27 | 12.351 | 1742 | 1745 | 1750 | rBV3 | 21915 | 25530 | 1.56% | 0.160% |
| 28 | 12.516 | 1770 | 1773 | 1779 | rBV | 179409 | 170127 | 10.38% | 1.068% |
| 29 | 12.745 | 1807 | 1812 | 1817 | rVB | 144711 | 165623 | 10.11% | 1.040% |
| 30 | 12.880 | 1831 | 1835 | 1838 | rBV | 1233893 | 1121715 | 68.45% | 7.041% |
| 31 | 13.080 | 1867 | 1869 | 1874 | rVB2 | 29000 | 31519 | 1.92% | 0.198% |
| 32 | 13.551 | 1946 | 1949 | 1953 | rBV2 | 61598 | 56776 | 3.46% | 0.356% |
| 33 | 13.757 | 1981 | 1984 | 1990 | rVB3 | 72917 | 90582 | 5.53% | 0.569% |
| 34 | 13.945 | 2012 | 2016 | 2020 | rBV | 443171 | 497961 | 30.39% | 3.126% |

LSC Area Percent Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110217\
Data File : BF100094.D
Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

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-BF102317.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 14.057 | 2033 | 2035 | 2037 | rBV3 | 20812 | 19800 | 1.21% | 0.124% |
| 36 | 14.204 | 2057 | 2060 | 2063 | rVB2 | 127725 | 104762 | 6.39% | 0.658% |
| 37 | 14.821 | 2162 | 2165 | 2168 | rVB3 | 47002 | 45718 | 2.79% | 0.287% |
| 38 | 14.998 | 2191 | 2195 | 2204 | rVB2 | 74011 | 146652 | 8.95% | 0.920% |
| 39 | 15.357 | 2253 | 2256 | 2261 | rVV2 | 55129 | 64744 | 3.95% | 0.406% |
| 40 | 15.410 | 2261 | 2265 | 2274 | rVB | 336542 | 420065 | 25.64% | 2.637% |
| 41 | 15.563 | 2288 | 2291 | 2297 | rVB8 | 27255 | 35291 | 2.15% | 0.222% |
| 42 | 16.439 | 2436 | 2440 | 2447 | rVB3 | 47154 | 81262 | 4.96% | 0.510% |
| 43 | 16.810 | 2501 | 2503 | 2510 | rVB6 | 21888 | 38791 | 2.37% | 0.243% |
| 44 | 17.862 | 2676 | 2682 | 2685 | rBV8 | 17212 | 38843 | 2.37% | 0.244% |

Sum of corrected areas: 15932202

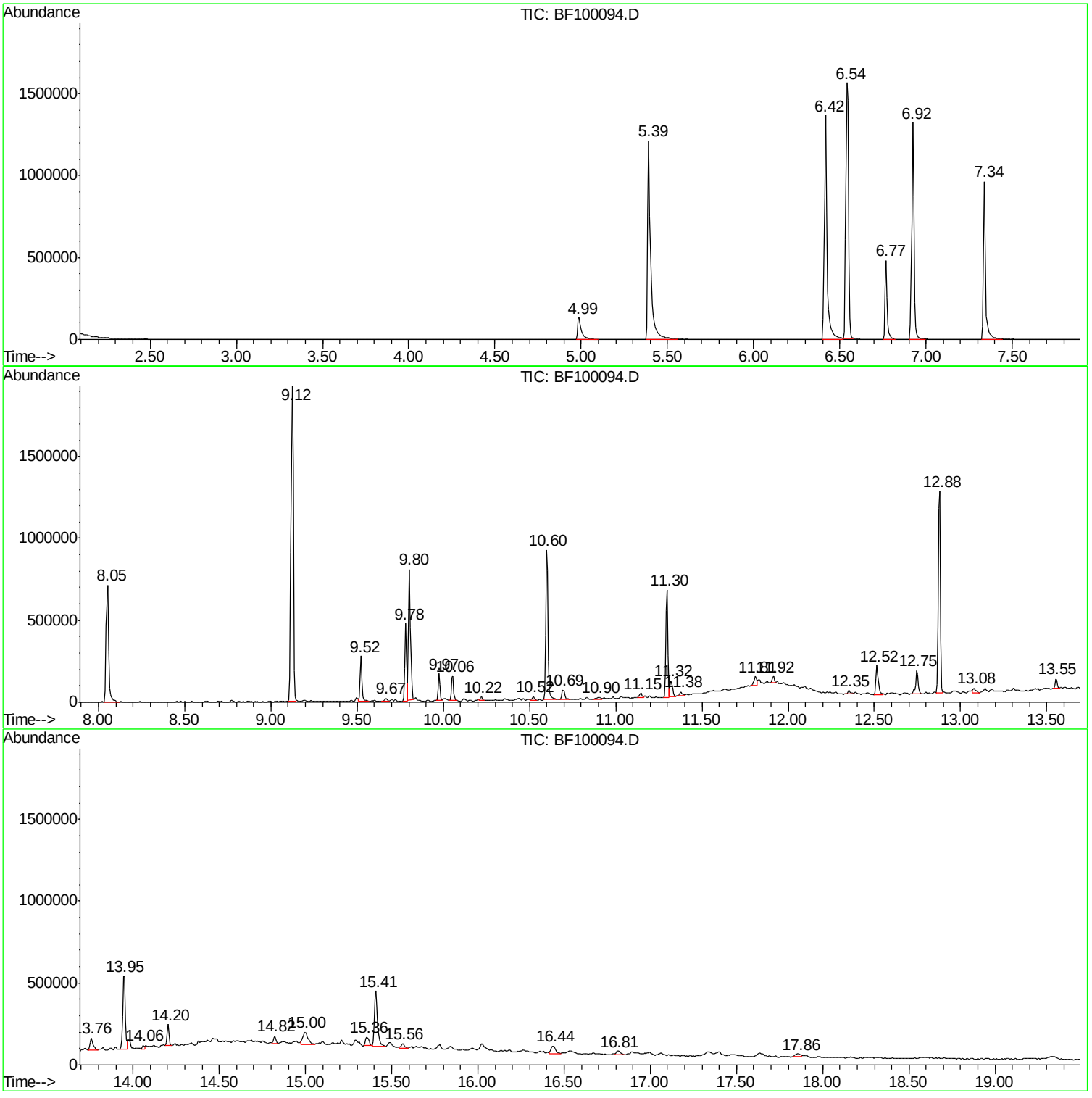
LSC Report - Integrated Chromatogram

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110217\
Data File : BF100094.D
Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110217\
Data File : BF100094.D
Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

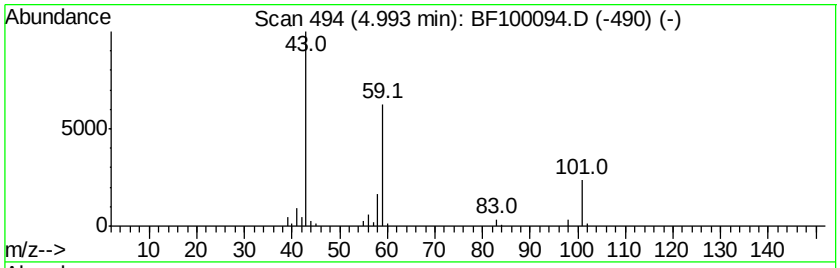
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

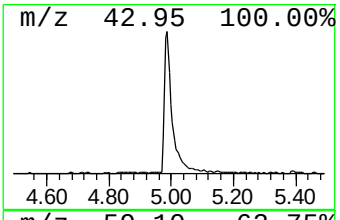
**********

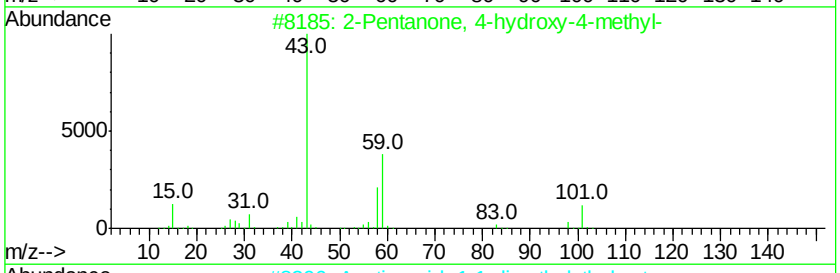
Peak Number 1 unknown4.99 Concentration Rank 3

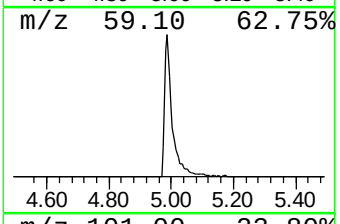
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.99 | 10.06 ng | 220379 | 1,4-Dichlorobenzene-d4 | 6.77 |

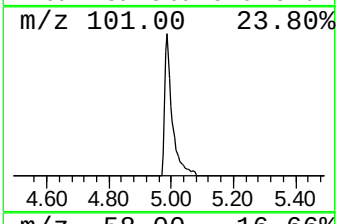
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 42 |
| 2 | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5 | 38 |
| 3 | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 25 |
| 4 | 5-Hexen-2-one | 98 | C6H10O | 000109-49-9 | 9 |
| 5 | Acetone | 58 | C3H6O | 000067-64-1 | 9 |

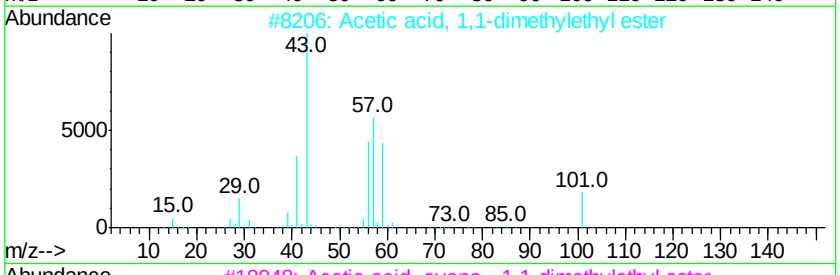


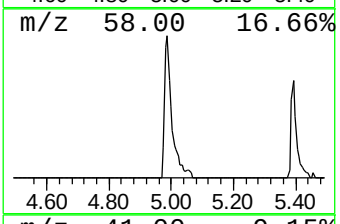



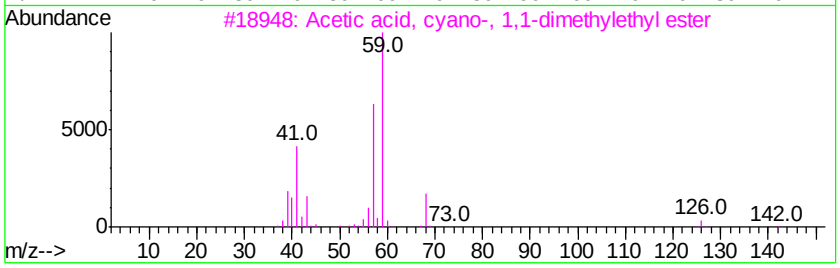



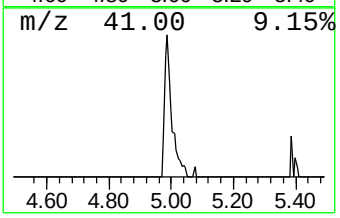



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110217\
Data File : BF100094.D
Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

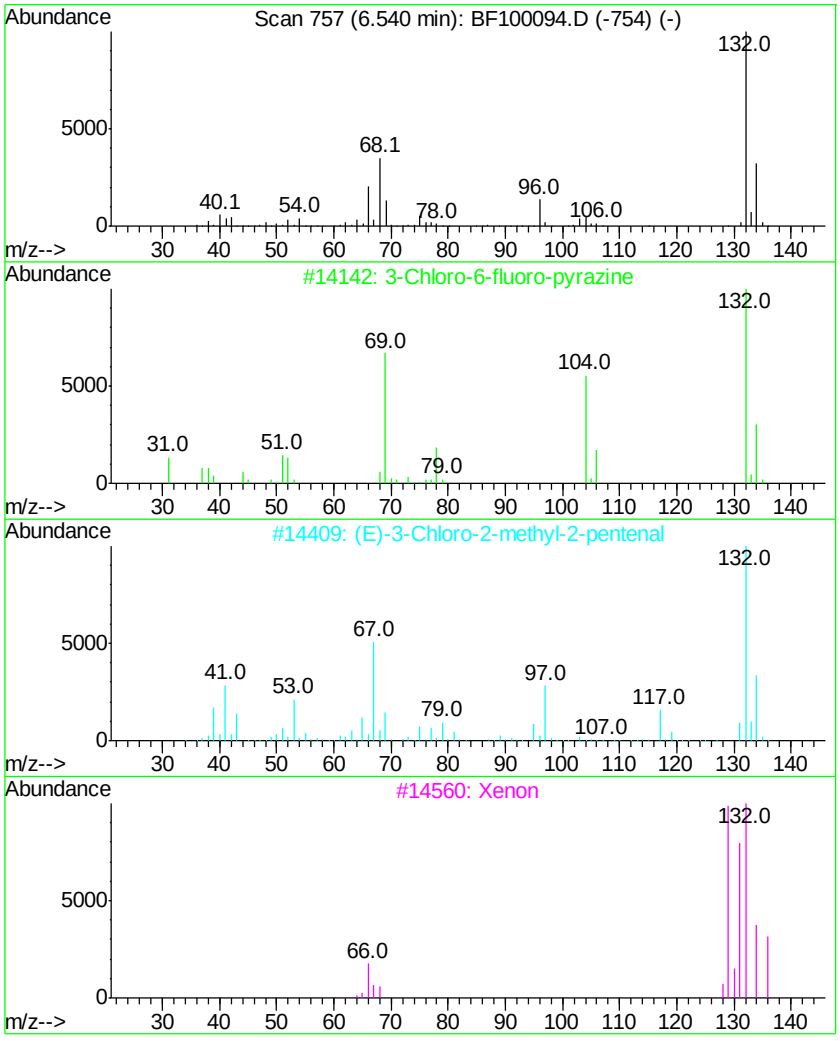
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

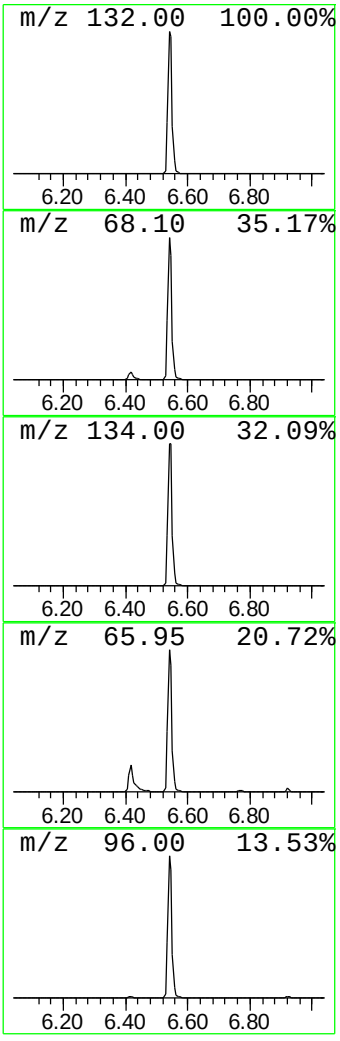
**********

Peak Number 2 unknown6.54 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.54 | 69.49 ng | 1521950 | 1,4-Dichlorobenzene-d4 | 6.77 |

| 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 | 17 |
| 3 | Xenon | 132 | Xe | 007440-63-3 | 9 |
| 4 | 4-Chloro-6-fluoro-pyrimidine | 132 | C4H2ClFN2 | 051422-01-6 | 9 |
| 5 | 5-Fluoro-2-chloropyrimidine | 132 | C4H2ClFN2 | 062802-42-0 | 9 |



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110217\
Data File : BF100094.D
Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

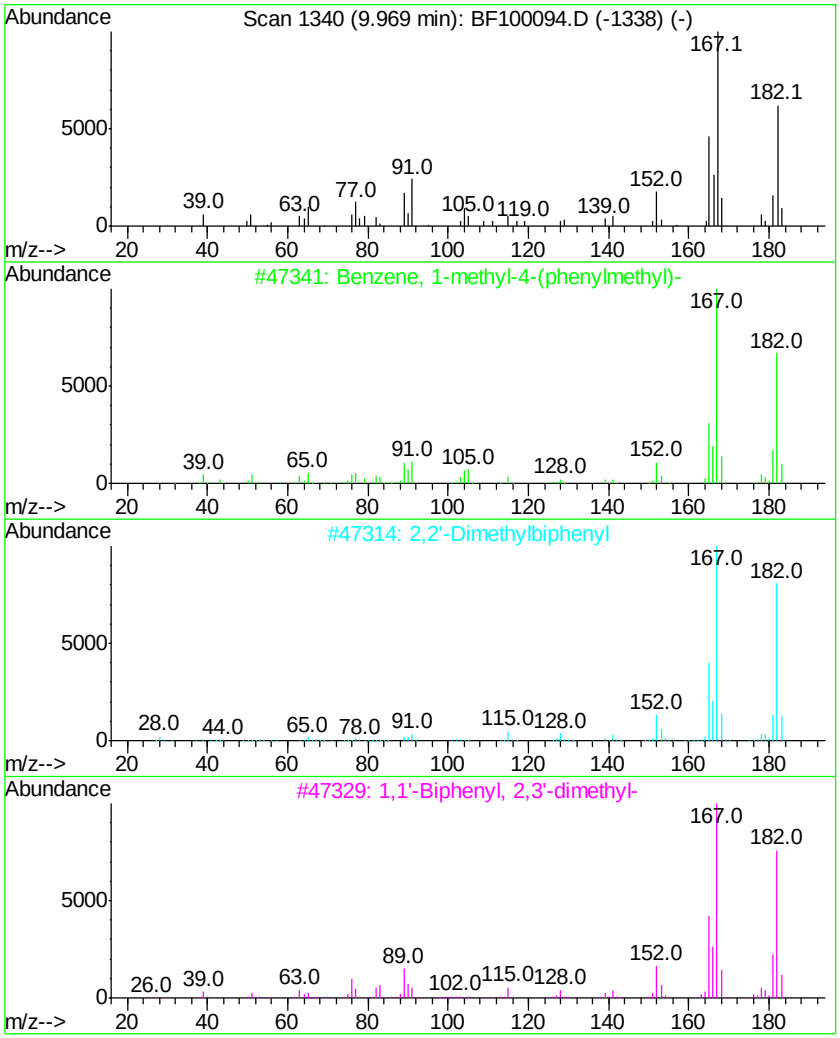
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

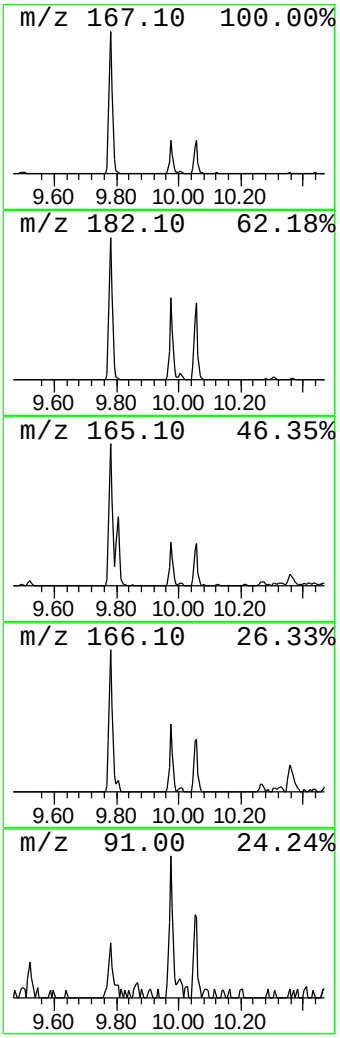
**********

Peak Number 4 Benzene, 1-methyl-4-(phenyl... Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.97 | 3.86 ng | 120410 | Acenaphthene-d10 | 9.80 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-methyl-4-(phenylmethyl)- | 182 | C14H14 | 000620-83-7 | 96 |
| 2 | 2,2'-Dimethylbiphenyl | 182 | C14H14 | 000605-39-0 | 94 |
| 3 | 1,1'-Biphenyl, 2,3'-dimethyl- | 182 | C14H14 | 000611-43-8 | 94 |
| 4 | Benzene, 1-methyl-2-(phenylmethyl)- | 182 | C14H14 | 000713-36-0 | 91 |
| 5 | Benzene, 1-methyl-3-(phenylmethyl)- | 182 | C14H14 | 000620-47-3 | 91 |



# Library Search Compound Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110217\
Data File : BF100094.D
Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

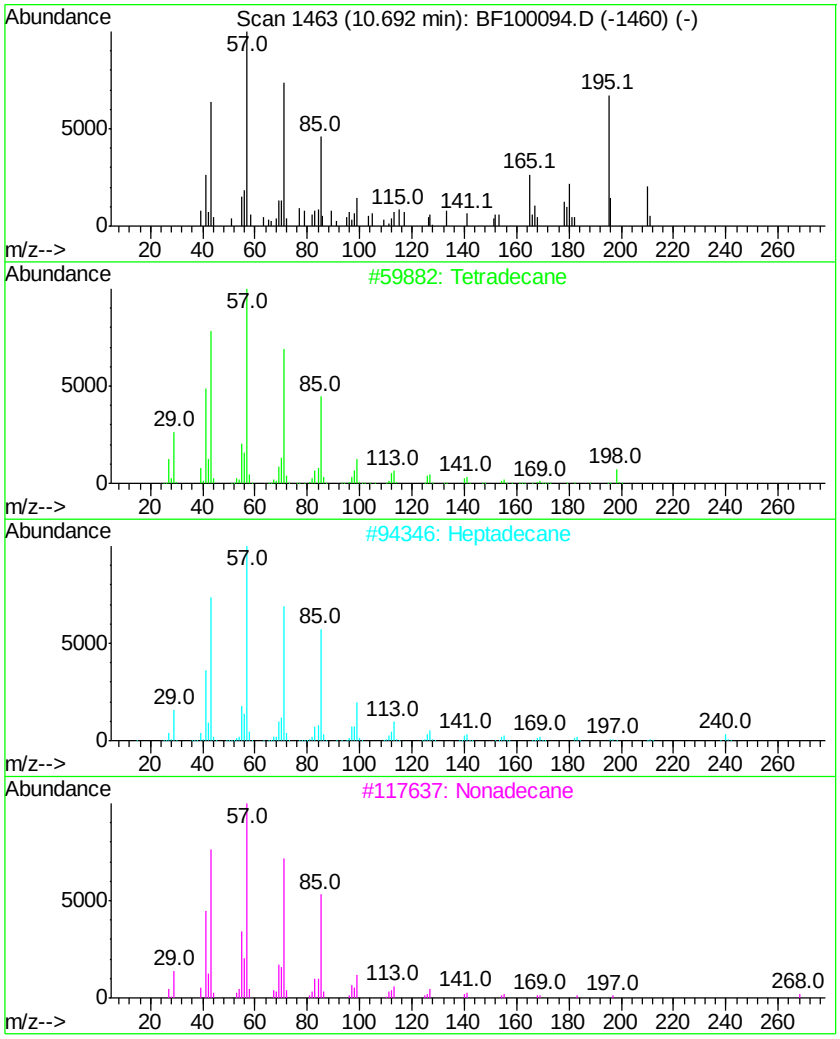
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

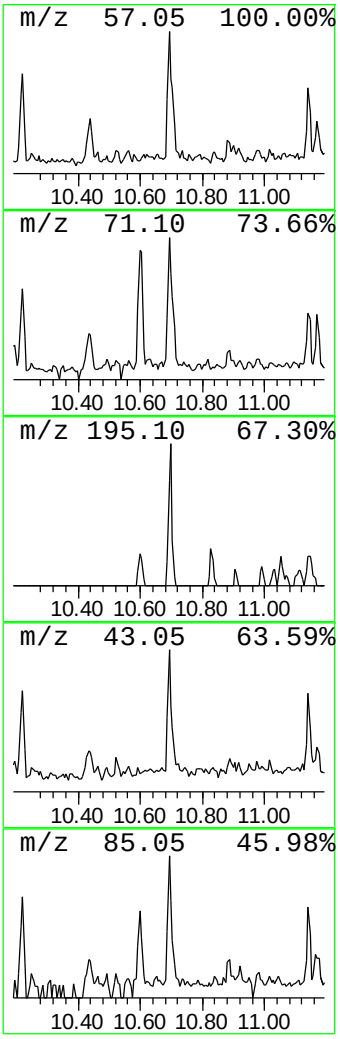
**********

Peak Number 6 Tetradecane Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.69 | 2.17 ng | 62610 | Phenanthrene-d10 | 11.30 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tetradecane | 198 | C14H30 | 000629-59-4 | 41 |
| 2 | Heptadecane | 240 | C17H36 | 000629-78-7 | 41 |
| 3 | Nonadecane | 268 | C19H40 | 000629-92-5 | 41 |
| 4 | 1-Octadecanesulphonyl chloride | 352 | C18H37ClO2S | 1000342-70-4 | 38 |
| 5 | Heneicosane | 296 | C21H44 | 000629-94-7 | 38 |



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Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

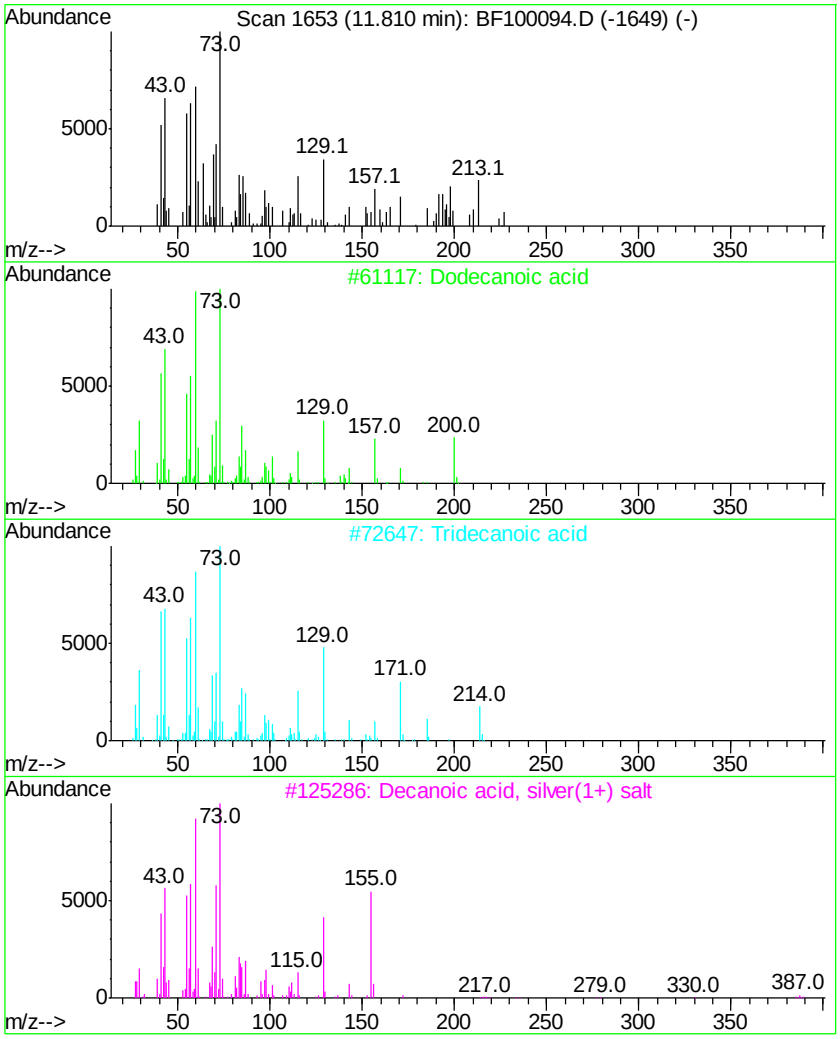
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

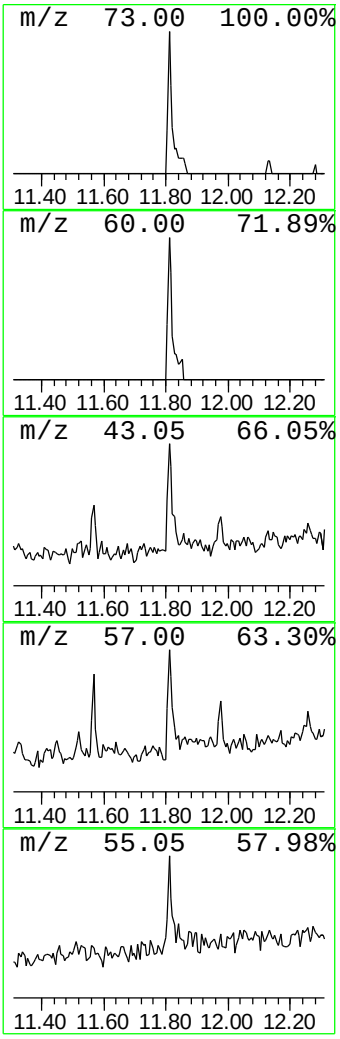
**********

Peak Number 7 unknown11.81 Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.81 | 2.35 ng | 67756 | Phenanthrene-d10 | 11.30 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dodecanoic acid | 200 | C12H24O2 | 000143-07-7 | 47 |
| 2 | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 43 |
| 3 | Decanoic acid, silver(1+) salt | 278 | C10H19AgO2 | 013126-67-5 | 35 |
| 4 | Undecanoic acid | 186 | C11H22O2 | 000112-37-8 | 32 |
| 5 | Trehalose | 342 | C12H22O11 | 000099-20-7 | 27 |



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Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

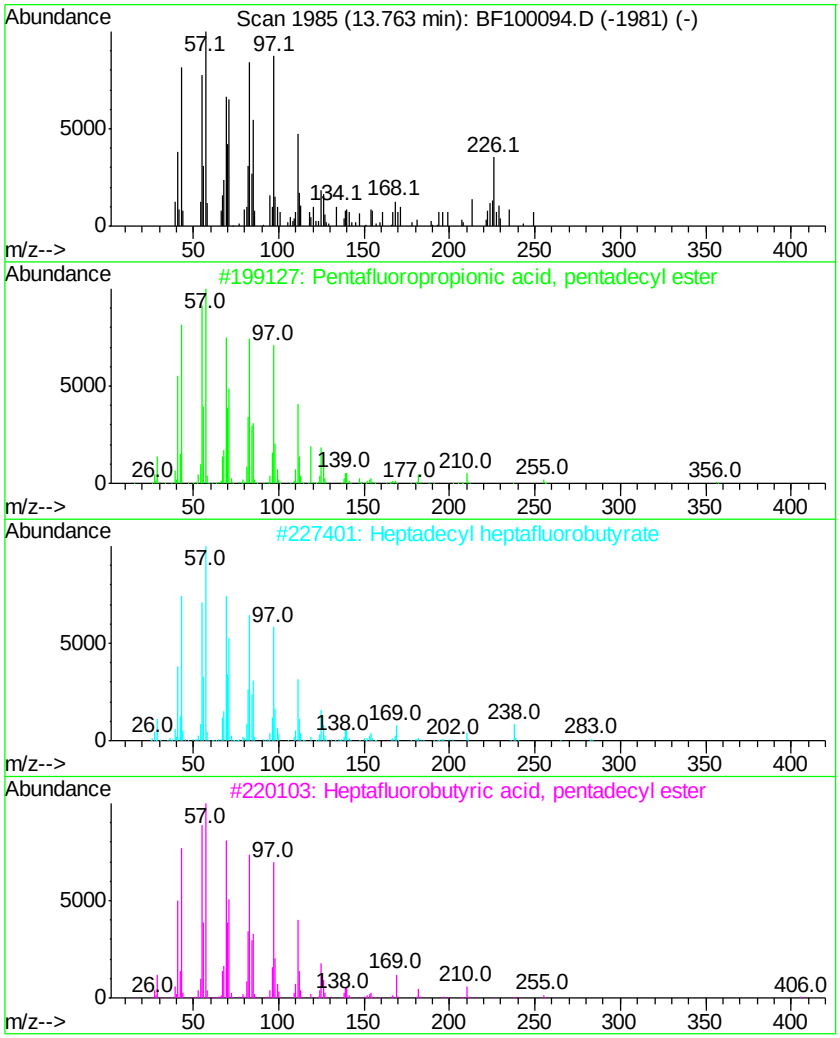
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

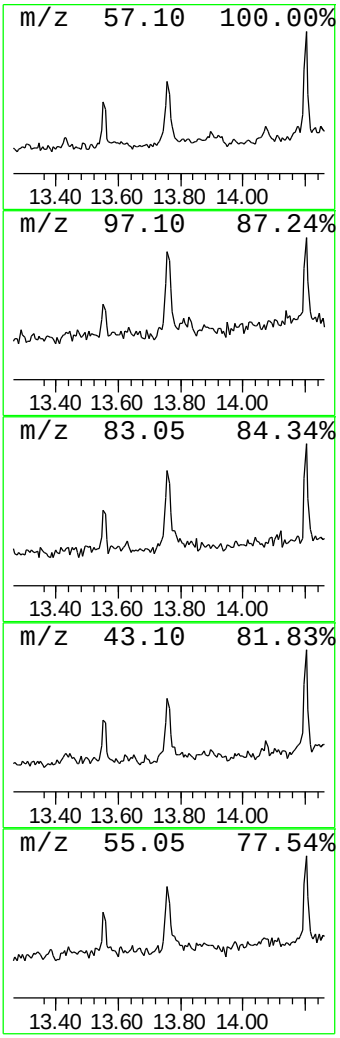
**********

Peak Number 9 Pentafluoropropionic acid, ... Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.76 | 3.64 ng | 90582 | Chrysene-d12 | 13.95 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentafluoropropionic acid, penta... | 374 | C18H31F5O2 | 959092-08-1 | 90 |
| 2 | Heptadecyl heptafluorobutyrate | 452 | C21H35F7O2 | 959085-66-6 | 90 |
| 3 | Heptafluorobutyric acid, pentade... | 424 | C19H31F7O2 | 959261-23-5 | 90 |
| 4 | Pentafluoropropionic acid, hexad... | 388 | C19H33F5O2 | 006222-07-7 | 83 |
| 5 | n-Tetracosanol-1 | 354 | C24H50O | 000506-51-4 | 81 |



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Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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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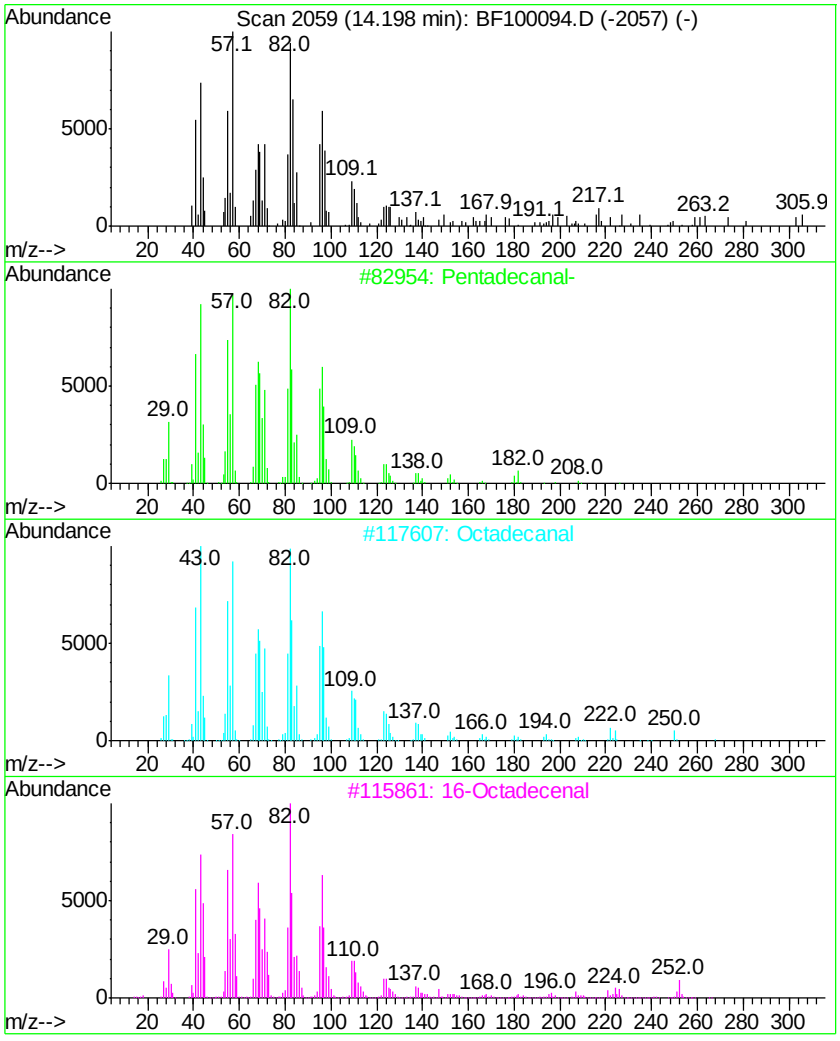
Peak Number 10 Pentadecanal- Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.20 | 4.21 ng | 104762 | Chrysene-d12 | 13.95 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentadecanal- | 226 | C15H30O | 002765-11-9 | 94 |
| 2 | Octadecanal | 268 | C18H36O | 000638-66-4 | 91 |
| 3 | 16-Octadecenal | 266 | C18H34O | 056554-87-1 | 91 |
| 4 | 17-Octadecenal | 266 | C18H34O | 056554-86-0 | 90 |
| 5 | 13-Octadecenal | 266 | C18H34O | 056554-90-6 | 86 |




m/z 57.10 100.00%
m/z 82.05 94.23%
m/z 43.05 73.69%
m/z 83.10 65.16%
m/z 55.00 59.39%

# Library Search Compound Report

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110217\
Data File : BF100094.D
Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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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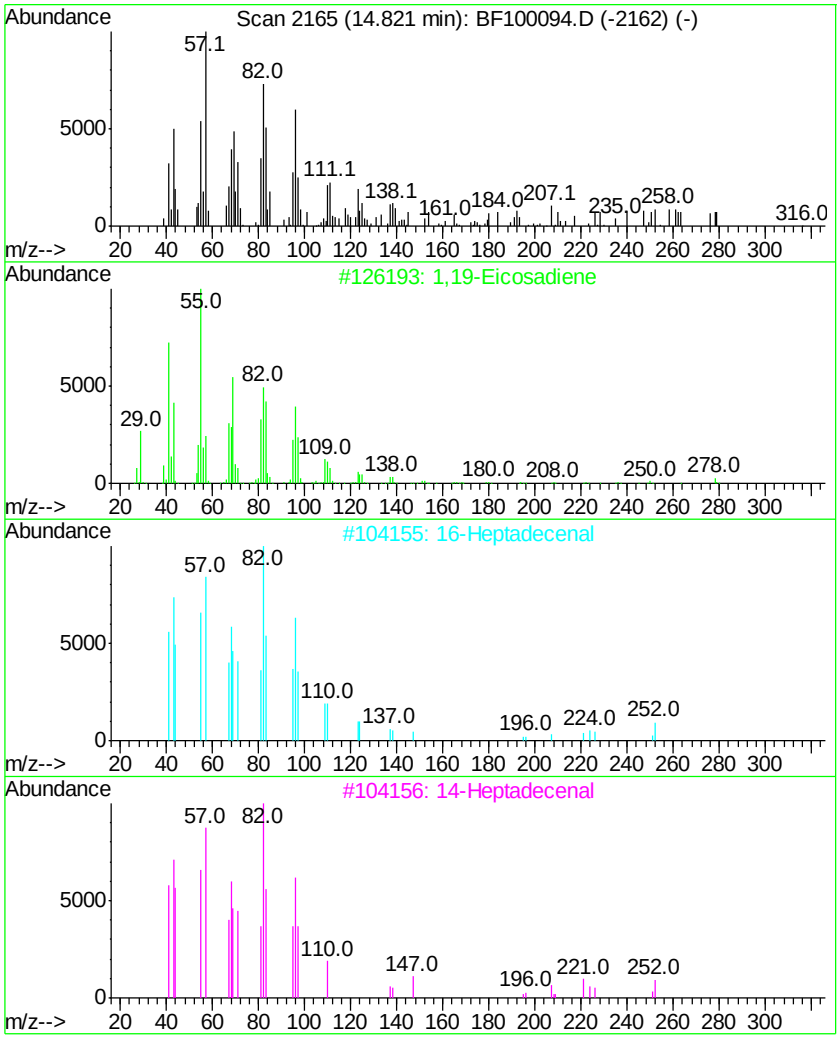
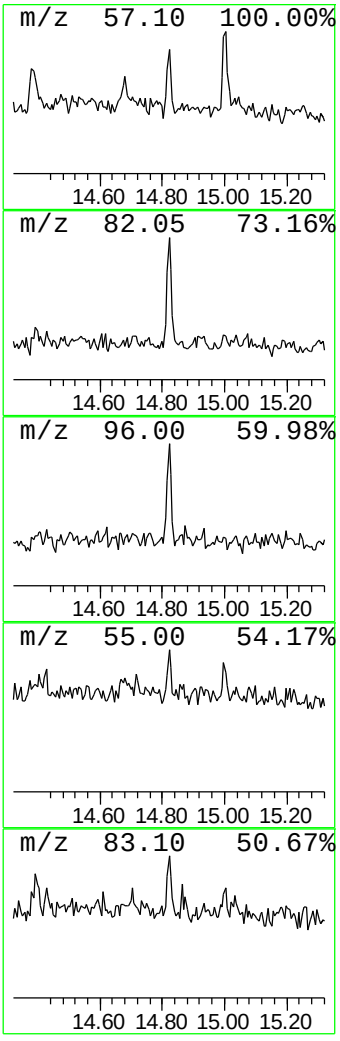
Peak Number 11 1,19-Eicosadiene Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.82 | 2.18 ng | 45718 | Perylene-d12 | 15.41 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,19-Eicosadiene | 278 | C20H38 | 014811-95-1 | 86 |
| 2 | 16-Heptadecenal | 252 | C17H32O | 1000144-57-9 | 83 |
| 3 | 14-Heptadecenal | 252 | C17H32O | 1000144-58-0 | 62 |
| 4 | Bicyclo[4.1.0]heptane, 7-butyl- | 152 | C11H20 | 018645-10-8 | 46 |
| 5 | 1-Cyclohexylheptene | 180 | C13H24 | 114614-83-4 | 43 |



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110217\
Data File : BF100094.D
Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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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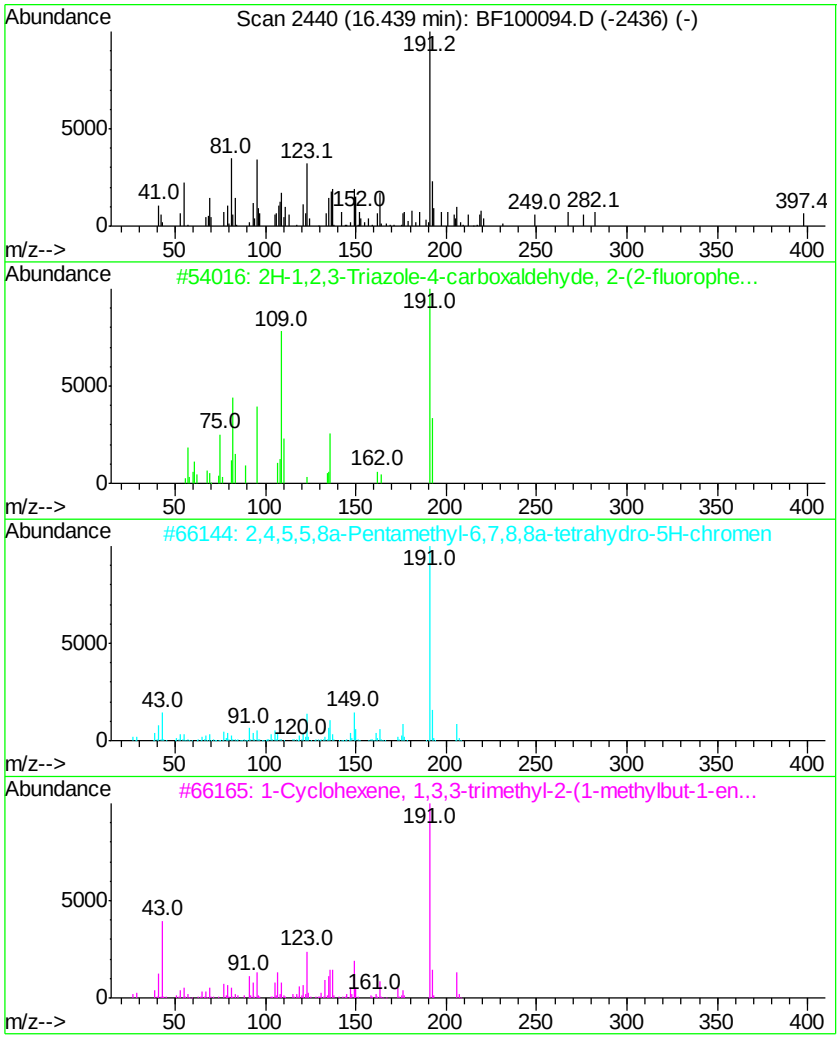
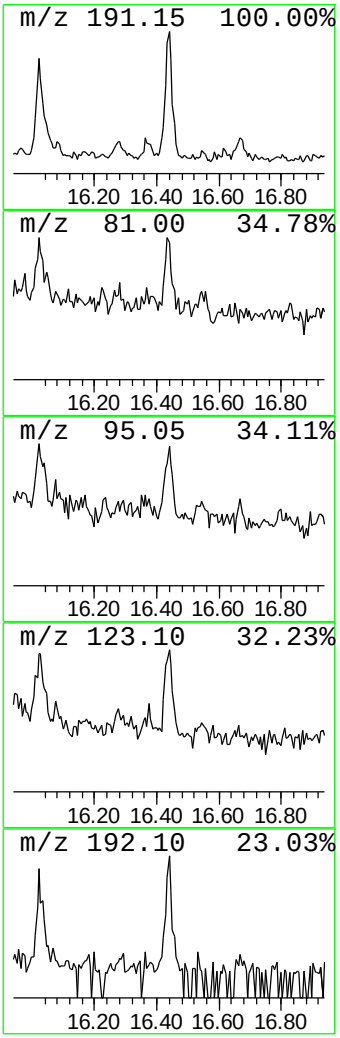
Peak Number 12 2H-1,2,3-Triazole-4-carboxa... Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.44 | 3.87 ng | 81262 | Perylene-d12 | 15.41 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2H-1,2,3-Triazole-4-carboxaldehv... | 191 | C9H6FN3O | 051306-43-5 | 70 |
| 2 | 2,4,5,5,8a-Pentamethyl-6,7,8,8a-... | 206 | C14H22O | 1000195-40-9 | 68 |
| 3 | 1-Cyclohexene, 1,3,3-trimethyl-2... | 206 | C14H22O | 1000197-08-4 | 59 |
| 4 | Amiphenazole | 191 | C9H9N3S | 000490-55-1 | 47 |
| 5 | 1-Penten-3-one, 1-(2,6,6-trimeth... | 206 | C14H22O | 000127-43-5 | 47 |



Tentatively Identified Compound (LSC) summary

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA_F\DATA\BF110217\
Data File : BF100094.D
Acq On : 2 Nov 2017 19:28
Operator : SJ/JU
Sample : I6249-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** COMP-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| | | | | | Internal Standard | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| unknown4.99 | 4.99 | 10.1 | ng | 220379 | 1 | 6.77 | 438008 | 20.0 |
| unknown6.54 | 6.54 | 69.5 | ng | 1521950 | 1 | 6.77 | 438008 | 20.0 |
| Benzene, 1-methyl... | 9.97 | 3.9 | ng | 120410 | 3 | 9.80 | 624047 | 20.0 |
| Tetradecane | 10.69 | 2.2 | ng | 62610 | 4 | 11.30 | 577536 | 20.0 |
| unknown11.81 | 11.81 | 2.4 | ng | 67756 | 4 | 11.30 | 577536 | 20.0 |
| Pentafluoropropio... | 13.76 | 3.6 | ng | 90582 | 5 | 13.95 | 497961 | 20.0 |
| Pentadecanal- | 14.20 | 4.2 | ng | 104762 | 5 | 13.95 | 497961 | 20.0 |
| 1,19-Eicosadiene | 14.82 | 2.2 | ng | 45718 | 6 | 15.41 | 420065 | 20.0 |
| 2H-1,2,3-Triazole... | 16.44 | 3.9 | ng | 81262 | 6 | 15.41 | 420065 | 20.0 |