LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % 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_M\METHODS\SOM-EPA-BM102417.M
Title : SVOA CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.157 | 5 | 12 | 20 | rBV | 1489754 | 1973138 | 13.95% | 1.249% |
| 2 | 3.263 | 26 | 30 | 38 | rVB | 172740 | 240880 | 1.70% | 0.152% |
| 3 | 3.357 | 42 | 46 | 57 | rVB3 | 93074 | 176363 | 1.25% | 0.112% |
| 4 | 3.798 | 87 | 121 | 124 | rBV2 | 1169955 | 6676796 | 47.19% | 4.225% |
| 5 | 4.175 | 148 | 185 | 188 | rBV | 1574860 | 9431573 | 66.66% | 5.968% |
| 6 | 4.575 | 249 | 253 | 259 | rVB | 184150 | 247236 | 1.75% | 0.156% |
| 7 | 4.781 | 281 | 288 | 292 | rBV | 176046 | 278767 | 1.97% | 0.176% |
| 8 | 4.910 | 302 | 310 | 317 | rBV | 138916 | 224466 | 1.59% | 0.142% |
| 9 | 5.357 | 380 | 386 | 395 | rVB3 | 129627 | 264766 | 1.87% | 0.168% |
| 10 | 5.516 | 408 | 413 | 427 | rVB | 76522 | 172376 | 1.22% | 0.109% |
| 11 | 5.716 | 441 | 447 | 461 | rVB | 2098847 | 2865396 | 20.25% | 1.813% |
| 12 | 5.875 | 469 | 474 | 480 | rBV2 | 206704 | 324151 | 2.29% | 0.205% |
| 13 | 6.134 | 512 | 518 | 526 | rBV5 | 69450 | 162566 | 1.15% | 0.103% |
| 14 | 6.398 | 557 | 563 | 565 | rBV4 | 120419 | 191437 | 1.35% | 0.121% |
| 15 | 6.534 | 578 | 586 | 596 | rVB4 | 172841 | 396978 | 2.81% | 0.251% |
| 16 | 6.987 | 630 | 663 | 665 | rBV | 1217228 | 5117276 | 36.17% | 3.238% |
| 17 | 7.028 | 665 | 670 | 686 | rVB2 | 1739236 | 3627165 | 25.64% | 2.295% |
| 18 | 7.186 | 692 | 697 | 703 | rVB | 1271515 | 1978576 | 13.98% | 1.252% |
| 19 | 7.392 | 721 | 732 | 745 | rVV | 1700692 | 3037198 | 21.47% | 1.922% |
| 20 | 7.510 | 749 | 752 | 760 | rVB3 | 96154 | 158078 | 1.12% | 0.100% |
| 21 | 7.786 | 788 | 799 | 802 | rBV6 | 172126 | 476373 | 3.37% | 0.301% |
| 22 | 7.857 | 802 | 811 | 818 | rVB2 | 1590973 | 3020301 | 21.35% | 1.911% |
| 23 | 7.928 | 818 | 823 | 828 | rVB3 | 128654 | 245273 | 1.73% | 0.155% |
| 24 | 7.992 | 828 | 834 | 838 | rBV5 | 98640 | 240966 | 1.70% | 0.152% |
| 25 | 8.045 | 838 | 843 | 847 | rVB4 | 212232 | 346800 | 2.45% | 0.219% |
| 26 | 8.116 | 847 | 855 | 857 | rBV3 | 113105 | 249331 | 1.76% | 0.158% |
| 27 | 8.416 | 899 | 906 | 916 | rVB5 | 169544 | 441485 | 3.12% | 0.279% |
| 28 | 8.563 | 922 | 931 | 939 | rBV | 1797106 | 2949278 | 20.85% | 1.866% |
| 29 | 8.733 | 956 | 960 | 970 | rVB5 | 114995 | 284566 | 2.01% | 0.180% |
| 30 | 8.957 | 988 | 998 | 1002 | rBV7 | 213746 | 558171 | 3.95% | 0.353% |
| 31 | 9.010 | 1002 | 1007 | 1013 | rBV | 1455532 | 2310571 | 16.33% | 1.462% |
| 32 | 9.169 | 1030 | 1034 | 1037 | rBV3 | 100387 | 163819 | 1.16% | 0.104% |
| 33 | 9.204 | 1037 | 1040 | 1046 | rVB4 | 112049 | 210435 | 1.49% | 0.133% |
| 34 | 9.322 | 1056 | 1060 | 1066 | rVB4 | 66767 | 150707 | 1.07% | 0.095% |

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % 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_M\METHODS\SOM-EPA-BM102417.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 9.498 | 1083 | 1090 | 1094 | rBV5 | 89129 | 169894 | 1.20% | 0.108% |
| 36 | 9.575 | 1097 | 1103 | 1107 | rVB5 | 104926 | 176151 | 1.25% | 0.111% |
| 37 | 9.733 | 1120 | 1130 | 1136 | rBV | 1341682 | 2307994 | 16.31% | 1.460% |
| 38 | 9.833 | 1141 | 1147 | 1155 | rVB5 | 155773 | 268723 | 1.90% | 0.170% |
| 39 | 10.022 | 1166 | 1179 | 1190 | rBV | 1143172 | 3010694 | 21.28% | 1.905% |
| 40 | 10.269 | 1211 | 1221 | 1230 | rVB | 2044322 | 3505342 | 24.78% | 2.218% |
| 41 | 10.651 | 1275 | 1286 | 1292 | rBV | 1958605 | 3527595 | 24.93% | 2.232% |
| 42 | 10.780 | 1302 | 1308 | 1321 | rBV2 | 686382 | 1418357 | 10.02% | 0.897% |
| 43 | 11.674 | 1455 | 1460 | 1464 | rBV2 | 106171 | 169678 | 1.20% | 0.107% |
| 44 | 11.963 | 1500 | 1509 | 1516 | rBV | 268057 | 461688 | 3.26% | 0.292% |
| 45 | 12.310 | 1562 | 1568 | 1575 | rVB3 | 65847 | 151492 | 1.07% | 0.096% |
| 46 | 13.027 | 1686 | 1690 | 1696 | rVB | 132346 | 175619 | 1.24% | 0.111% |
| 47 | 13.316 | 1735 | 1739 | 1746 | rVB2 | 183525 | 283217 | 2.00% | 0.179% |
| 48 | 13.657 | 1789 | 1797 | 1803 | rBV5 | 90275 | 253565 | 1.79% | 0.160% |
| 49 | 13.810 | 1818 | 1823 | 1828 | rBV2 | 146531 | 245437 | 1.73% | 0.155% |
| 50 | 13.898 | 1828 | 1838 | 1843 | rVV | 3426719 | 5025421 | 35.52% | 3.180% |
| 51 | 13.939 | 1843 | 1845 | 1849 | rVV | 321464 | 460859 | 3.26% | 0.292% |
| 52 | 13.974 | 1849 | 1851 | 1855 | rVB2 | 181297 | 187665 | 1.33% | 0.119% |
| 53 | 14.180 | 1880 | 1886 | 1898 | rVB | 3950099 | 5888283 | 41.62% | 3.726% |
| 54 | 14.321 | 1906 | 1910 | 1917 | rVB5 | 113942 | 193173 | 1.37% | 0.122% |
| 55 | 14.386 | 1917 | 1921 | 1927 | rBV | 282329 | 386611 | 2.73% | 0.245% |
| 56 | 14.486 | 1930 | 1938 | 1945 | rVV2 | 2859171 | 4585919 | 32.41% | 2.902% |
| 57 | 14.692 | 1967 | 1973 | 1983 | rBV | 1544392 | 2432185 | 17.19% | 1.539% |
| 58 | 14.839 | 1993 | 1998 | 2002 | rBV2 | 237224 | 378729 | 2.68% | 0.240% |
| 59 | 14.892 | 2003 | 2007 | 2013 | rVV4 | 212870 | 405429 | 2.87% | 0.257% |
| 60 | 14.962 | 2014 | 2019 | 2023 | rVV | 149068 | 229319 | 1.62% | 0.145% |
| 61 | 15.010 | 2024 | 2027 | 2036 | rVB8 | 106085 | 188473 | 1.33% | 0.119% |
| 62 | 15.110 | 2040 | 2044 | 2048 | rVV2 | 128547 | 218265 | 1.54% | 0.138% |
| 63 | 15.157 | 2049 | 2052 | 2056 | rVB | 125373 | 180126 | 1.27% | 0.114% |
| 64 | 15.339 | 2079 | 2083 | 2088 | rVB | 338286 | 410816 | 2.90% | 0.260% |
| 65 | 15.480 | 2101 | 2107 | 2113 | rVB | 5078645 | 7184261 | 50.78% | 4.546% |
| 66 | 15.604 | 2123 | 2128 | 2134 | rVV | 858510 | 1243281 | 8.79% | 0.787% |
| 67 | 15.657 | 2134 | 2137 | 2141 | rVB2 | 124370 | 174439 | 1.23% | 0.110% |
| 68 | 15.745 | 2149 | 2152 | 2159 | rVB | 262206 | 370194 | 2.62% | 0.234% |
| 69 | 15.845 | 2164 | 2169 | 2174 | rVB | 1078590 | 1521472 | 10.75% | 0.963% |
| 70 | 16.204 | 2225 | 2230 | 2232 | rBV2 | 490250 | 768016 | 5.43% | 0.486% |
| 71 | 16.598 | 2293 | 2297 | 2300 | rVB4 | 130086 | 182244 | 1.29% | 0.115% |

LSC Area Percent Report

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Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % 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_M\METHODS\SOM-EPA-BM102417.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 16.633 | 2300 | 2303 | 2311 | rVB4 | 247401 | 317335 | 2.24% | 0.201% |
| 73 | 16.992 | 2360 | 2364 | 2368 | rBV | 484665 | 570215 | 4.03% | 0.361% |
| 74 | 17.045 | 2369 | 2373 | 2383 | rVB2 | 433245 | 718468 | 5.08% | 0.455% |
| 75 | 17.233 | 2399 | 2405 | 2409 | rBV2 | 3740298 | 5276351 | 37.29% | 3.339% |
| 76 | 17.274 | 2409 | 2412 | 2416 | rVV | 888145 | 1214512 | 8.58% | 0.768% |
| 77 | 17.327 | 2416 | 2421 | 2431 | rVB2 | 5364944 | 8028374 | 56.74% | 5.080% |
| 78 | 17.727 | 2486 | 2489 | 2494 | rVB | 423171 | 485358 | 3.43% | 0.307% |
| 79 | 18.127 | 2550 | 2557 | 2571 | rBV2 | 2604972 | 4721821 | 33.37% | 2.988% |
| 80 | 18.421 | 2604 | 2607 | 2610 | rBV | 396940 | 484917 | 3.43% | 0.307% |
| 81 | 19.286 | 2750 | 2754 | 2761 | rVB | 2087820 | 3070704 | 21.70% | 1.943% |
| 82 | 19.397 | 2771 | 2773 | 2777 | rVB2 | 764594 | 787778 | 5.57% | 0.498% |
| 83 | 19.621 | 2806 | 2811 | 2815 | rBV | 7274649 | 10568922 | 74.70% | 6.688% |
| 84 | 20.539 | 2964 | 2967 | 2971 | rVB | 1639402 | 1850688 | 13.08% | 1.171% |
| 85 | 21.403 | 3109 | 3114 | 3118 | rBV | 5590638 | 7859790 | 55.55% | 4.973% |
| 86 | 23.568 | 3477 | 3482 | 3497 | rVB3 | 5998263 | 14148406 | 100.00% | 8.953% |

Sum of corrected areas: 158037523

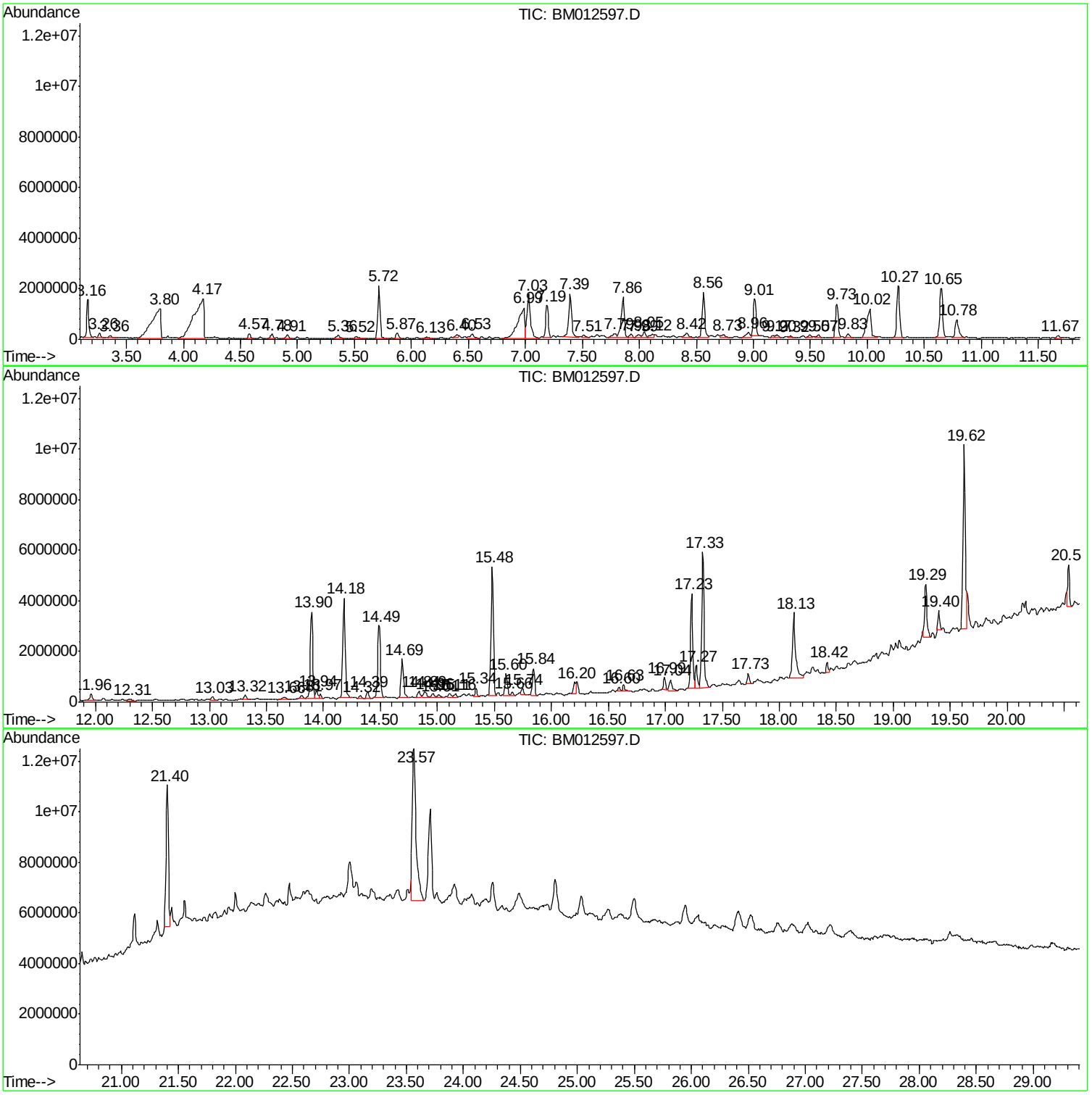
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

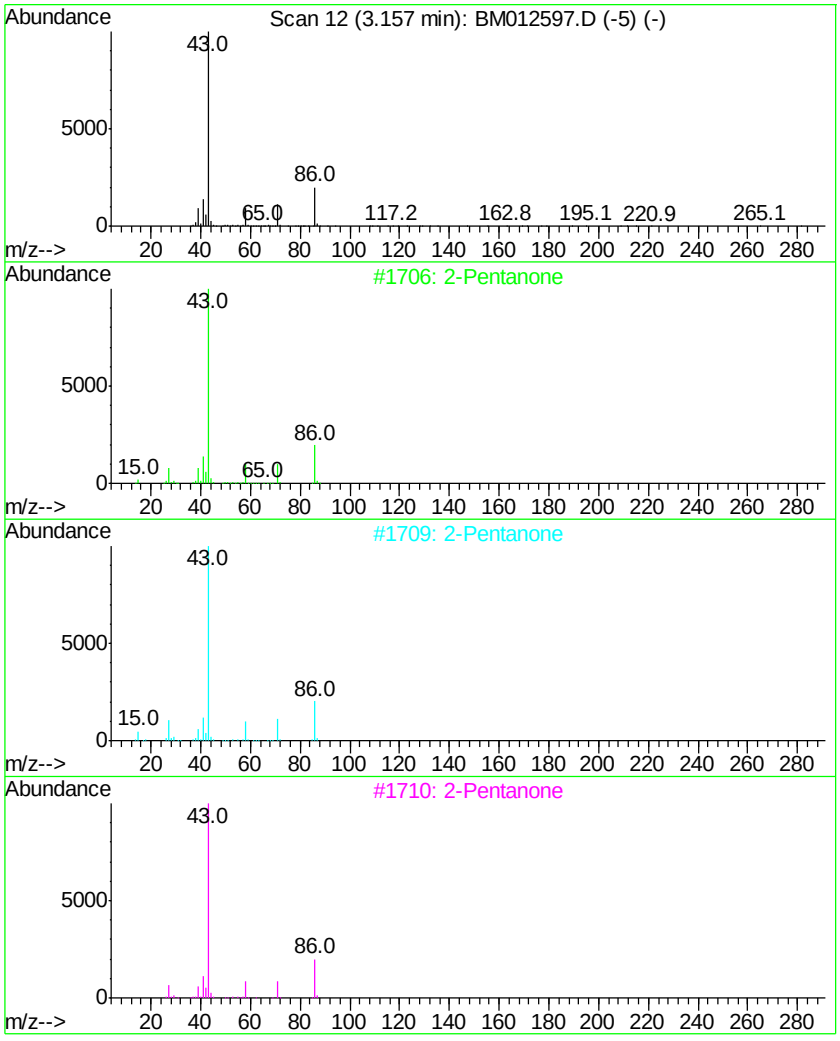
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

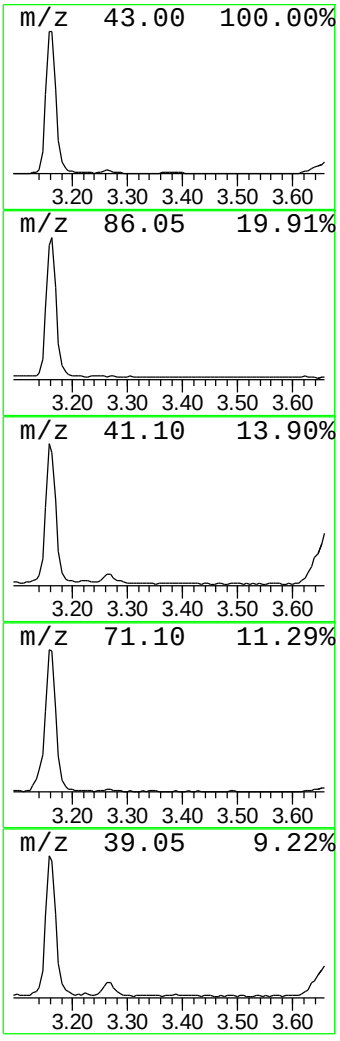
**********

Peak Number 1 2-Pentanone Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.16 | 13.07 ng/ul | 1973140 | 1,4-Dichlorobenzene-d4 | 7.86 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone | 86 | C5H10O | 000107-87-9 | 91 |
| 2 | 2-Pentanone | 86 | C5H10O | 000107-87-9 | 90 |
| 3 | 2-Pentanone | 86 | C5H10O | 000107-87-9 | 86 |
| 4 | 2-Butanone, 3-methyl- | 86 | C5H10O | 000563-80-4 | 59 |
| 5 | 2,3-Butanedione | 86 | C4H6O2 | 000431-03-8 | 52 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

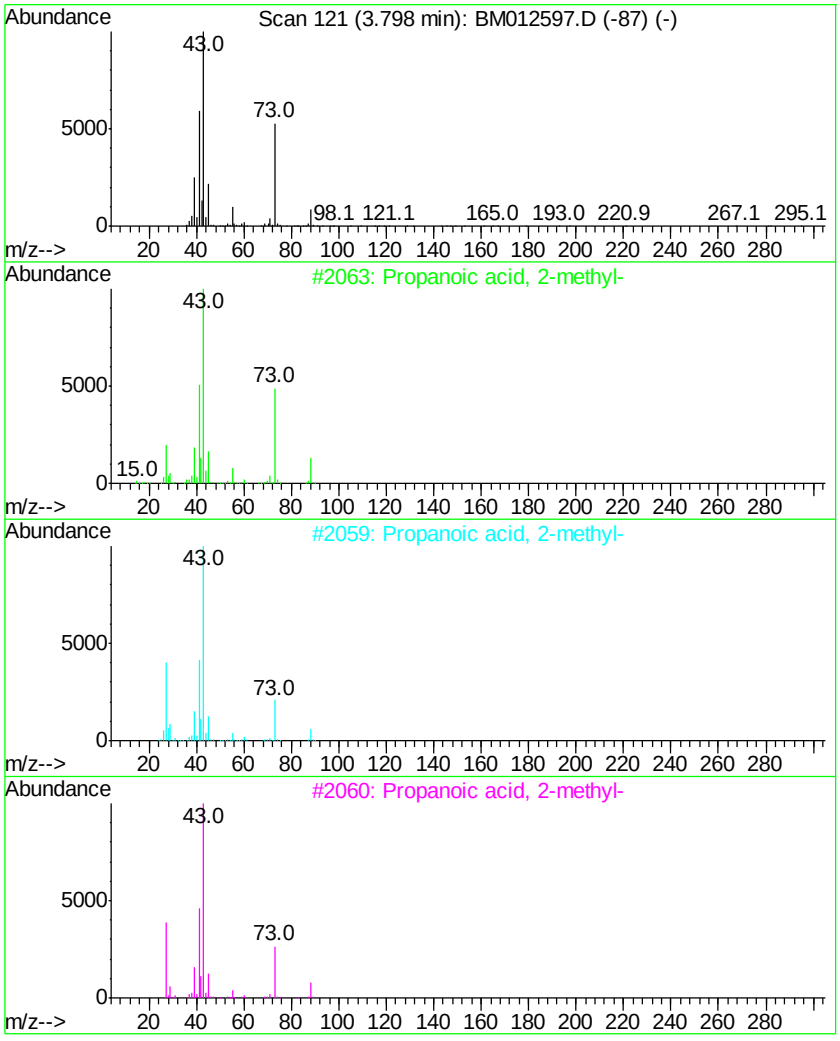
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

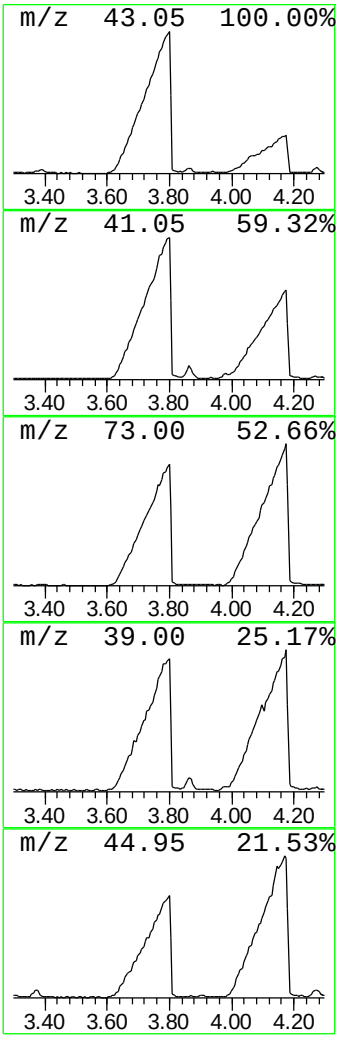
**********

Peak Number 2 Propanoic acid, 2-methyl- Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.80 | 44.21 ng/ul | 6676800 | 1,4-Dichlorobenzene-d4 | 7.86 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propanoic acid, 2-methyl- | 88 | C4H8O2 | 000079-31-2 | 90 |
| 2 | Propanoic acid, 2-methyl- | 88 | C4H8O2 | 000079-31-2 | 86 |
| 3 | Propanoic acid, 2-methyl- | 88 | C4H8O2 | 000079-31-2 | 86 |
| 4 | Propanoic acid, 2-methyl- | 88 | C4H8O2 | 000079-31-2 | 80 |
| 5 | Propanoic acid, 2-methyl- | 88 | C4H8O2 | 000079-31-2 | 80 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
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Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

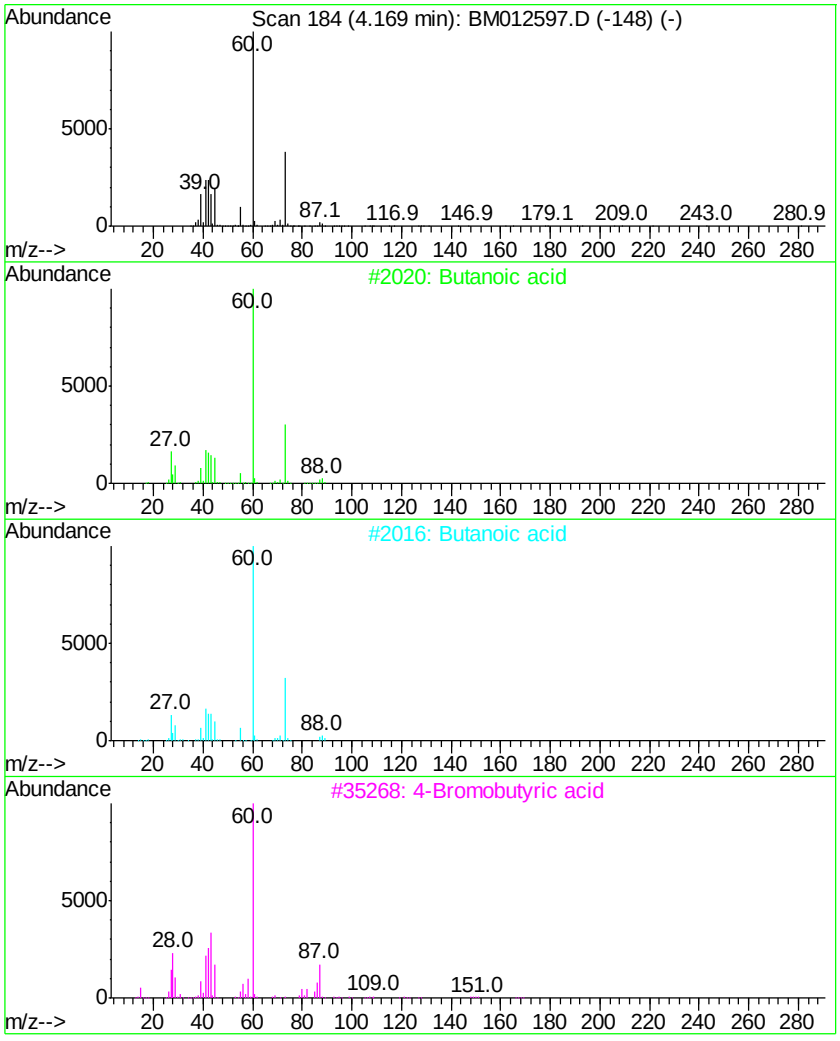
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

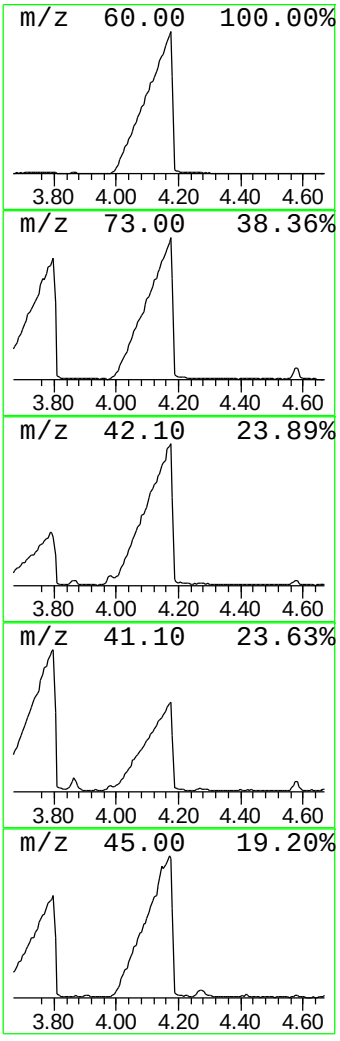
**********

Peak Number 3 Butanoic acid Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.17 | 62.45 ng/ul | 9431570 | 1,4-Dichlorobenzene-d4 | 7.86 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butanoic acid | 88 | C4H8O2 | 000107-92-6 | 72 |
| 2 | Butanoic acid | 88 | C4H8O2 | 000107-92-6 | 50 |
| 3 | 4-Bromobutyric acid | 166 | C4H7BrO2 | 002623-87-2 | 9 |
| 4 | Pentanoic acid | 102 | C5H10O2 | 000109-52-4 | 9 |
| 5 | Hexanoic acid | 116 | C6H12O2 | 000142-62-1 | 9 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

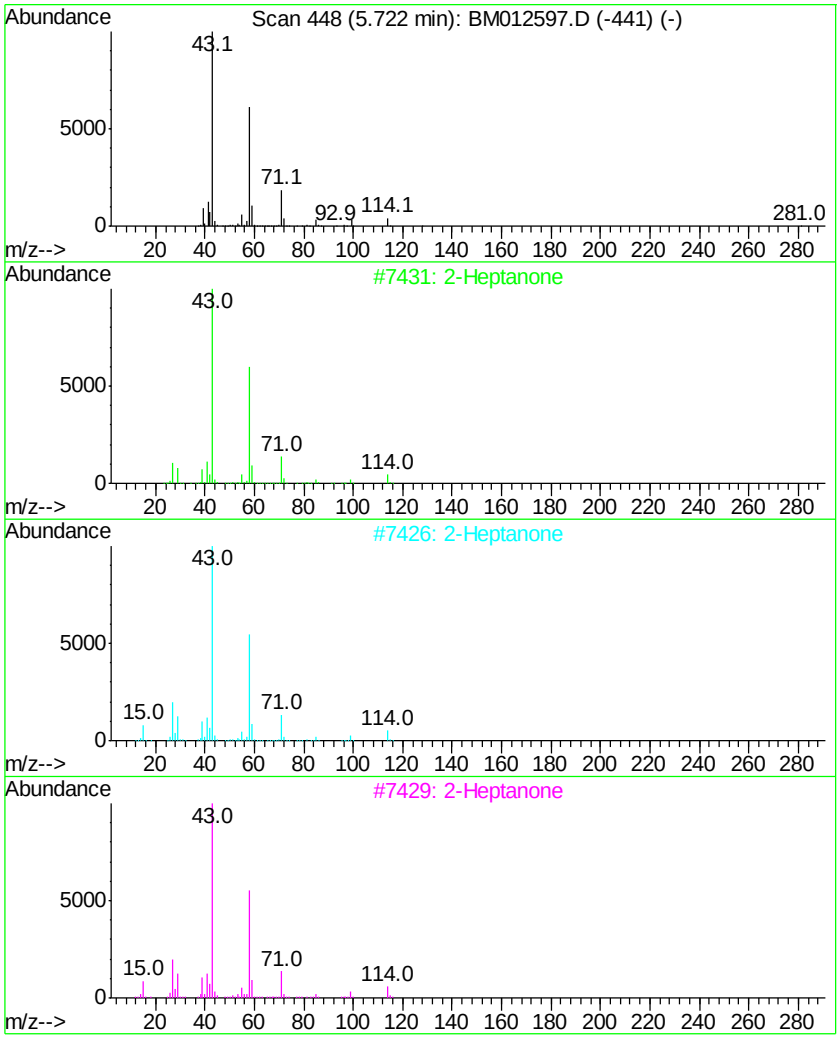
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

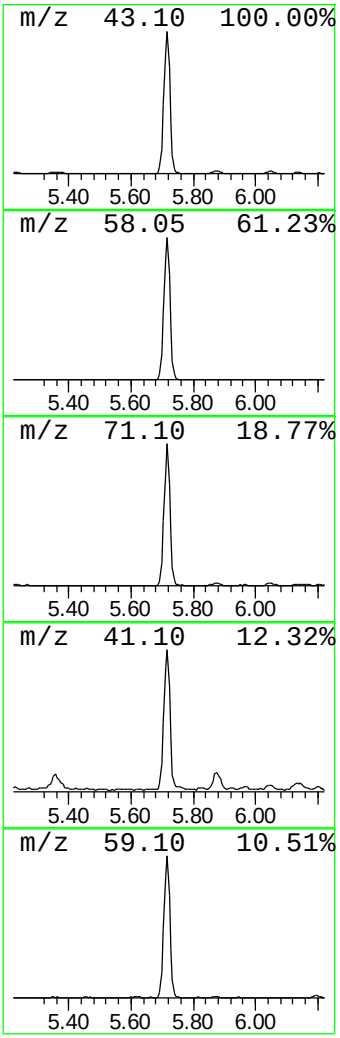
**********

Peak Number 4 2-Heptanone Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.72 | 18.97 ng/ul | 2865400 | 1,4-Dichlorobenzene-d4 | 7.86 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Heptanone | 114 | C7H14O | 000110-43-0 | 90 |
| 2 | 2-Heptanone | 114 | C7H14O | 000110-43-0 | 86 |
| 3 | 2-Heptanone | 114 | C7H14O | 000110-43-0 | 83 |
| 4 | 2-Hexanone, 5-methyl- | 114 | C7H14O | 000110-12-3 | 72 |
| 5 | 2-Hexanone, 5-methyl- | 114 | C7H14O | 000110-12-3 | 72 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

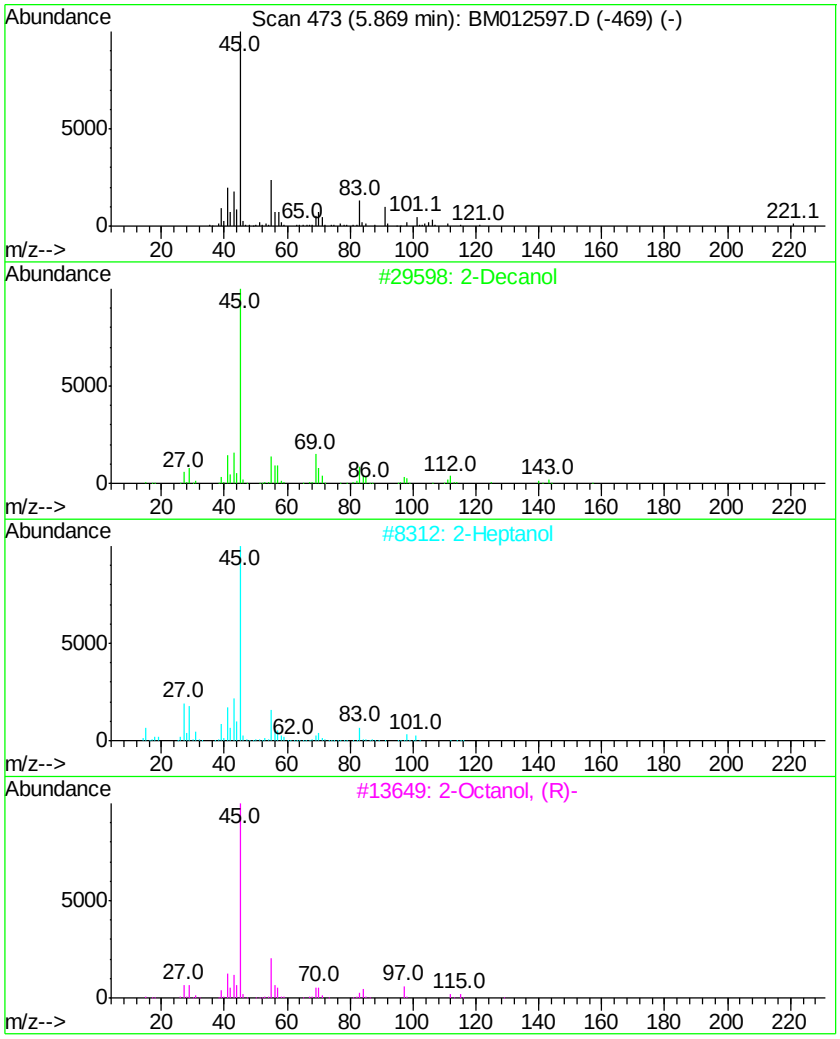
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

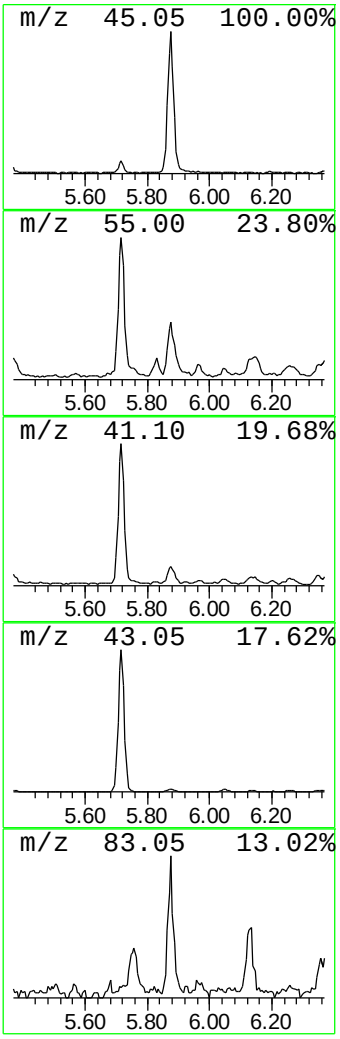
**********

Peak Number 5 2-Decanol Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.87 | 2.15 ng/ul | 324151 | 1,4-Dichlorobenzene-d4 | 7.86 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Decanol | 158 | C10H22O | 001120-06-5 | 59 |
| 2 | 2-Heptanol | 116 | C7H16O | 000543-49-7 | 59 |
| 3 | 2-Octanol, (R)- | 130 | C8H18O | 005978-70-1 | 53 |
| 4 | 2-Heptanol | 116 | C7H16O | 000543-49-7 | 50 |
| 5 | 2-Heptanol | 116 | C7H16O | 000543-49-7 | 50 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

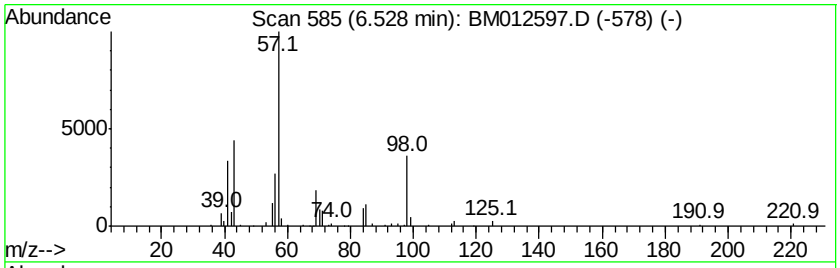
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

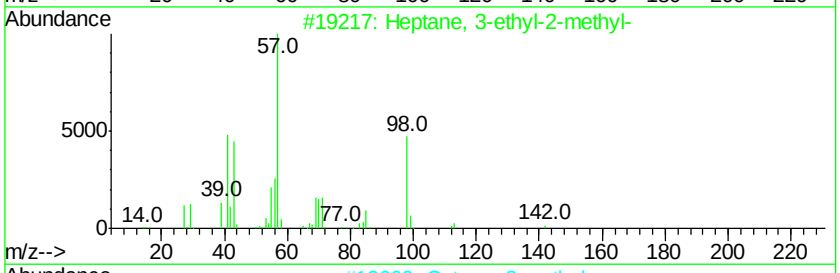
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

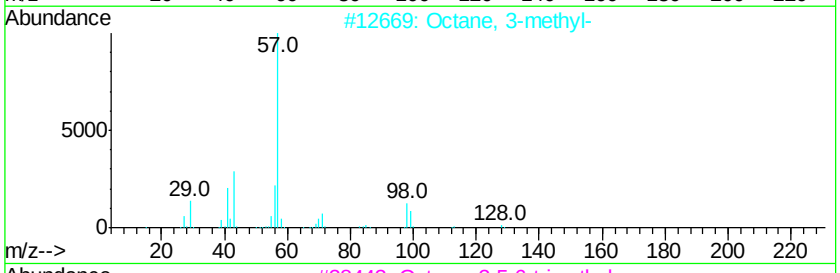
**********
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 14

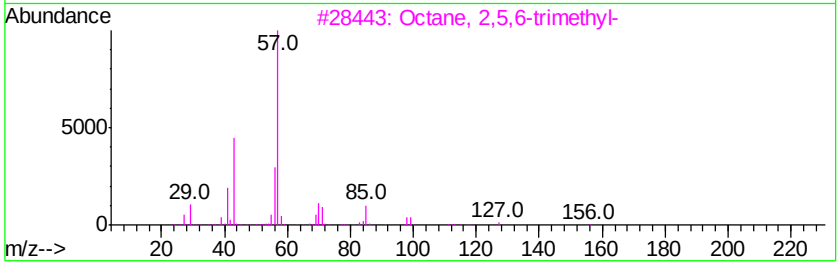
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.53 | 2.63 ng/ul | 396978 | 1,4-Dichlorobenzene-d4 | 7.86 |

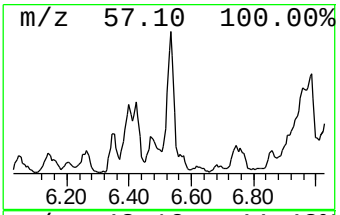
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptane, 3-ethyl-2-methyl- | 142 | C10H22 | 014676-29-0 | 72 |
| 2 | Octane, 3-methyl- | 128 | C9H20 | 002216-33-3 | 47 |
| 3 | Octane, 2,5,6-trimethyl- | 156 | C11H24 | 062016-14-2 | 47 |
| 4 | Propanoic acid, 2,2-dimethyl-, o... | 214 | C13H26O2 | 027751-88-8 | 38 |
| 5 | Heptane, 2,5-dimethyl- | 128 | C9H20 | 002216-30-0 | 38 |

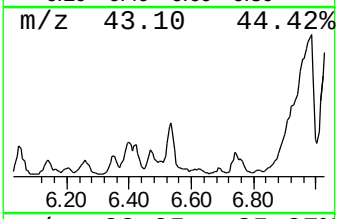


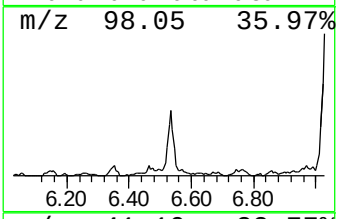



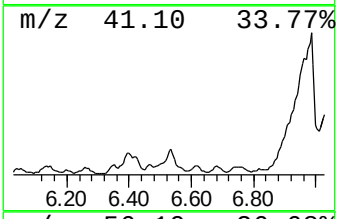



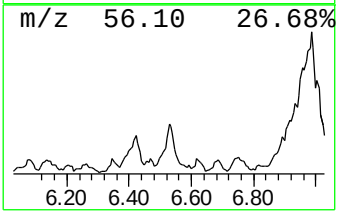



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

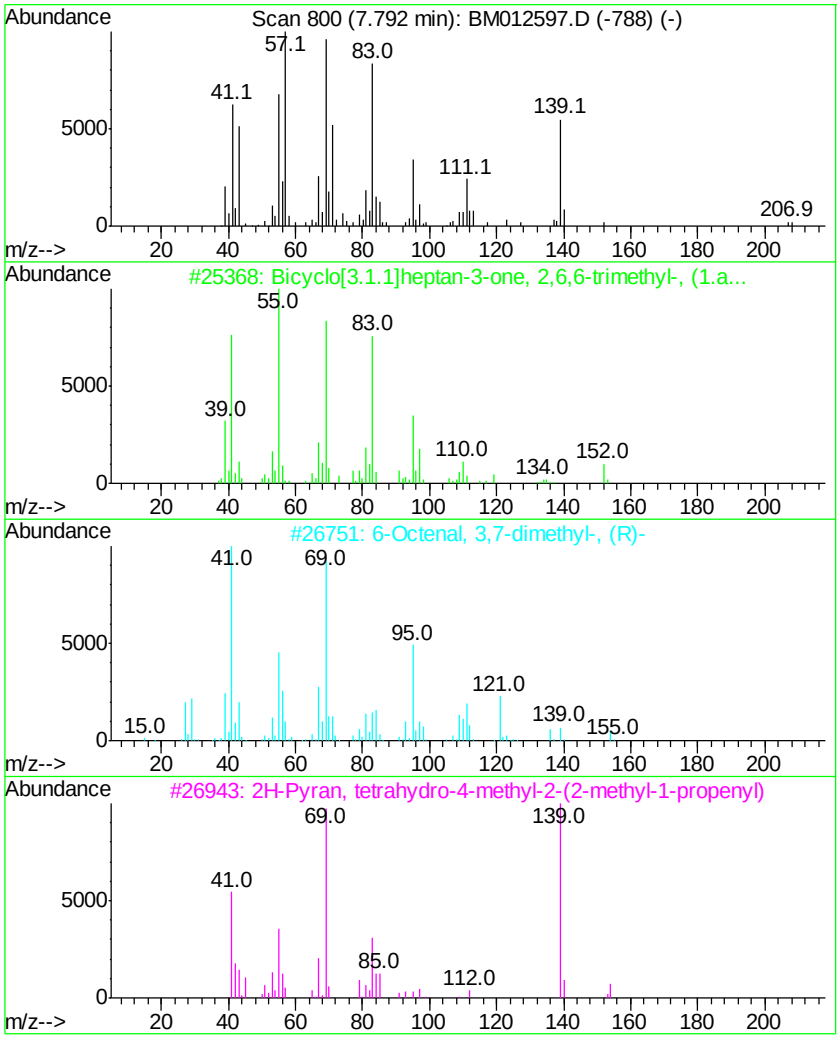
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

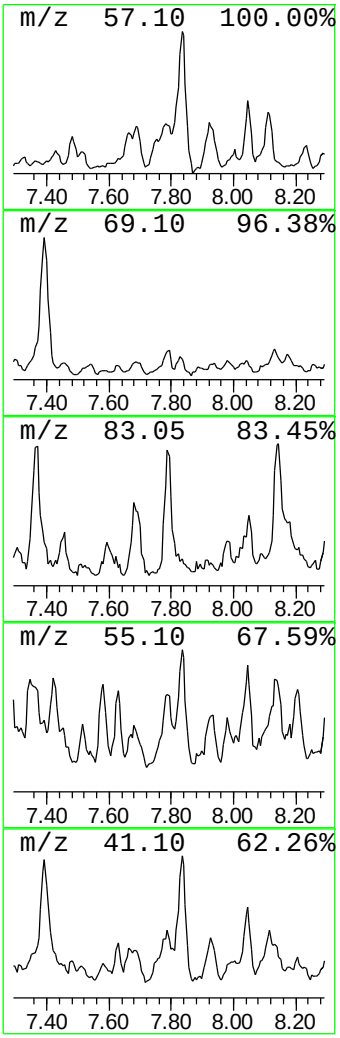
**********

Peak Number 7 unknown-01 Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.79 | 3.15 ng/ul | 476373 | 1,4-Dichlorobenzene-d4 | 7.86 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Bicyclo[3.1.1]heptan-3-one, 2,6,... | 152 | C10H16O | 015358-88-0 | 49 |
| 2 | 6-Octenal, 3,7-dimethyl-, (R)- | 154 | C10H18O | 002385-77-5 | 35 |
| 3 | 2H-Pyran, tetrahydro-4-methyl-2-... | 154 | C10H18O | 016409-43-1 | 35 |
| 4 | 2-Hydroxy-2-methyl-but-3-enyl 2-... | 184 | C10H16O3 | 080758-67-4 | 27 |
| 5 | Cyclopentane, 1-butyl-2-ethyl- | 154 | C11H22 | 072993-32-9 | 14 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

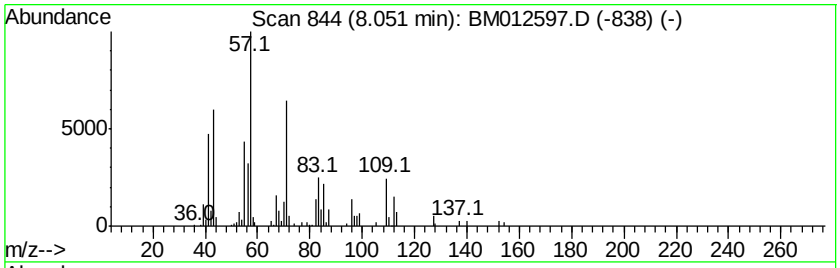
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

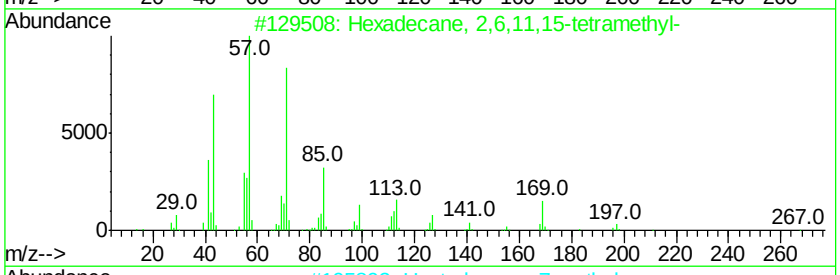
**********

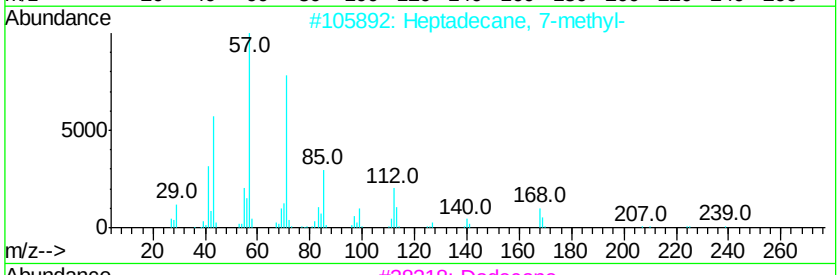
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 16

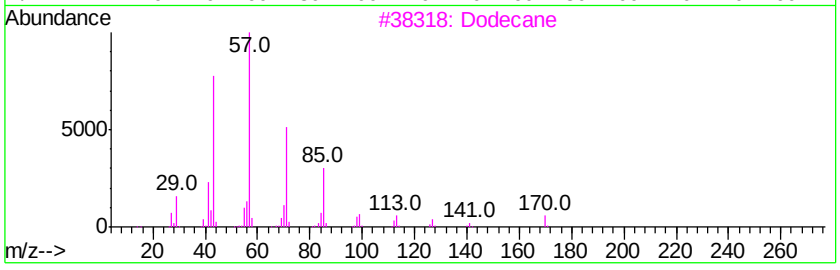
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.05 | 2.30 ng/ul | 346800 | 1,4-Dichlorobenzene-d4 | 7.86 |

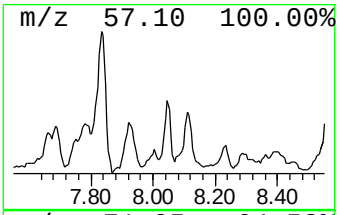
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexadecane, 2,6,11,15-tetramethyl- | 282 | C20H42 | 000504-44-9 | 64 |
| 2 | Heptadecane, 7-methyl- | 254 | C18H38 | 020959-33-5 | 58 |
| 3 | Dodecane | 170 | C12H26 | 000112-40-3 | 50 |
| 4 | Tridecane, 5-propyl- | 226 | C16H34 | 055045-11-9 | 50 |
| 5 | Tetradecane | 198 | C14H30 | 000629-59-4 | 50 |

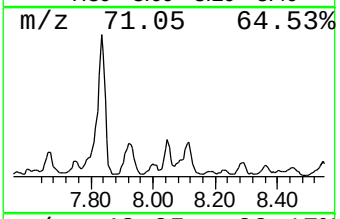


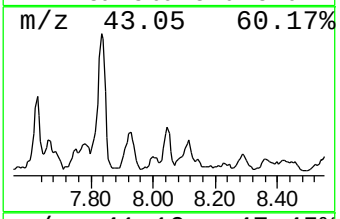



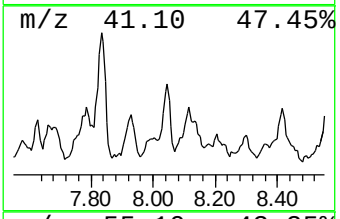



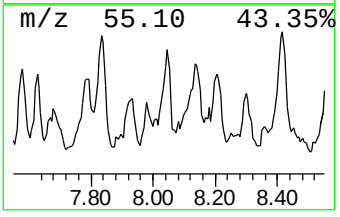



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

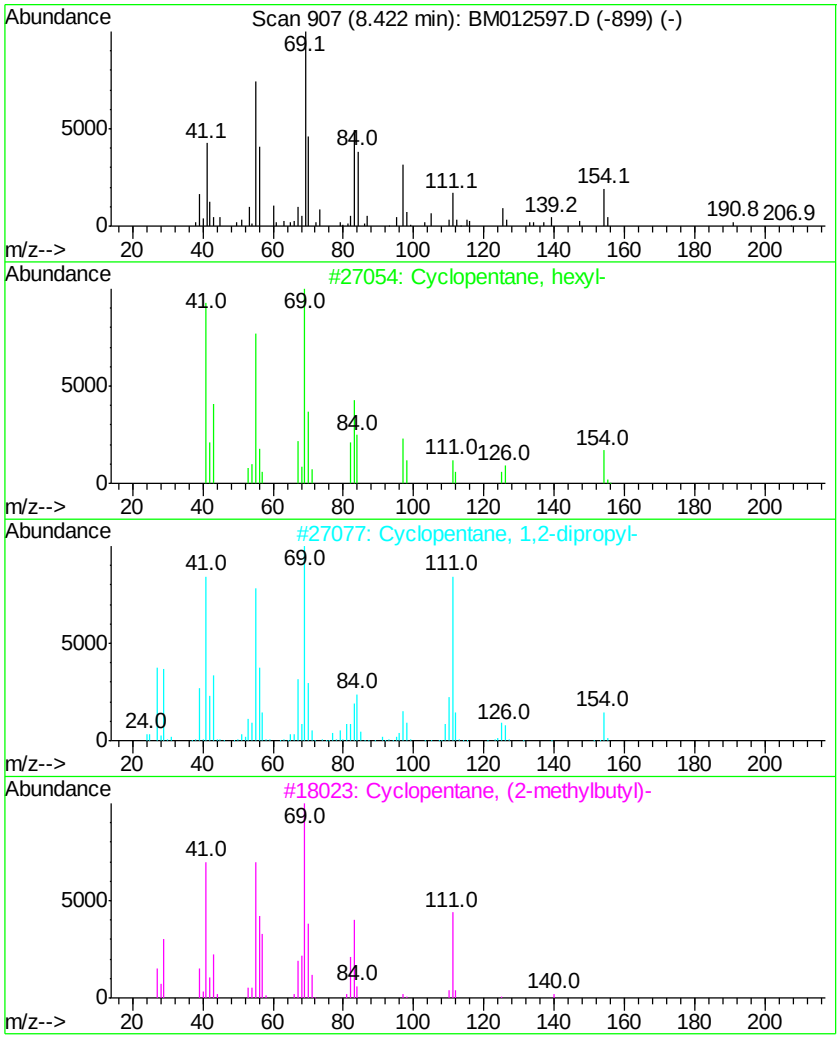
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

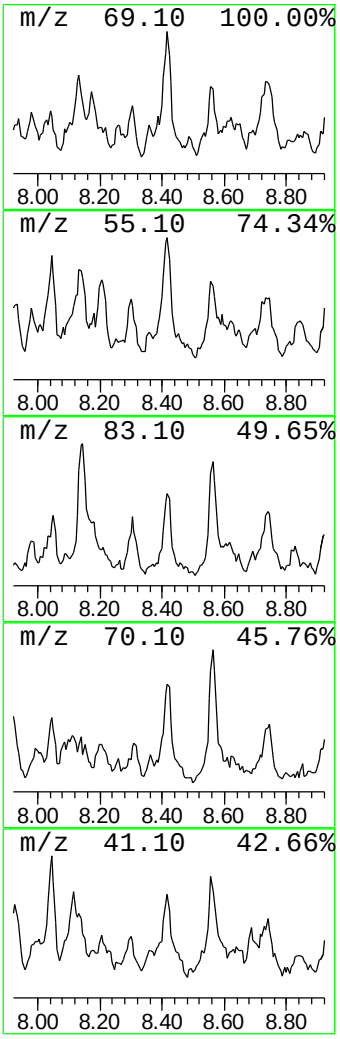
**********

Peak Number 9 (DEL) Alkane: Cyclic8.42 Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.42 | 2.92 ng/ul | 441485 | 1,4-Dichlorobenzene-d4 | 7.86 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclopentane, hexyl- | 154 | C11H22 | 004457-00-5 | 64 |
| 2 | Cyclopentane, 1,2-dipropyl- | 154 | C11H22 | 091242-57-8 | 52 |
| 3 | Cyclopentane, (2-methylbutyl)- | 140 | C10H20 | 053366-38-4 | 50 |
| 4 | Nonane, 2-methyl-3-methylene- | 154 | C11H22 | 055499-08-6 | 50 |
| 5 | 2-Decene, 4-methyl-, (Z)- | 154 | C11H22 | 074630-30-1 | 50 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

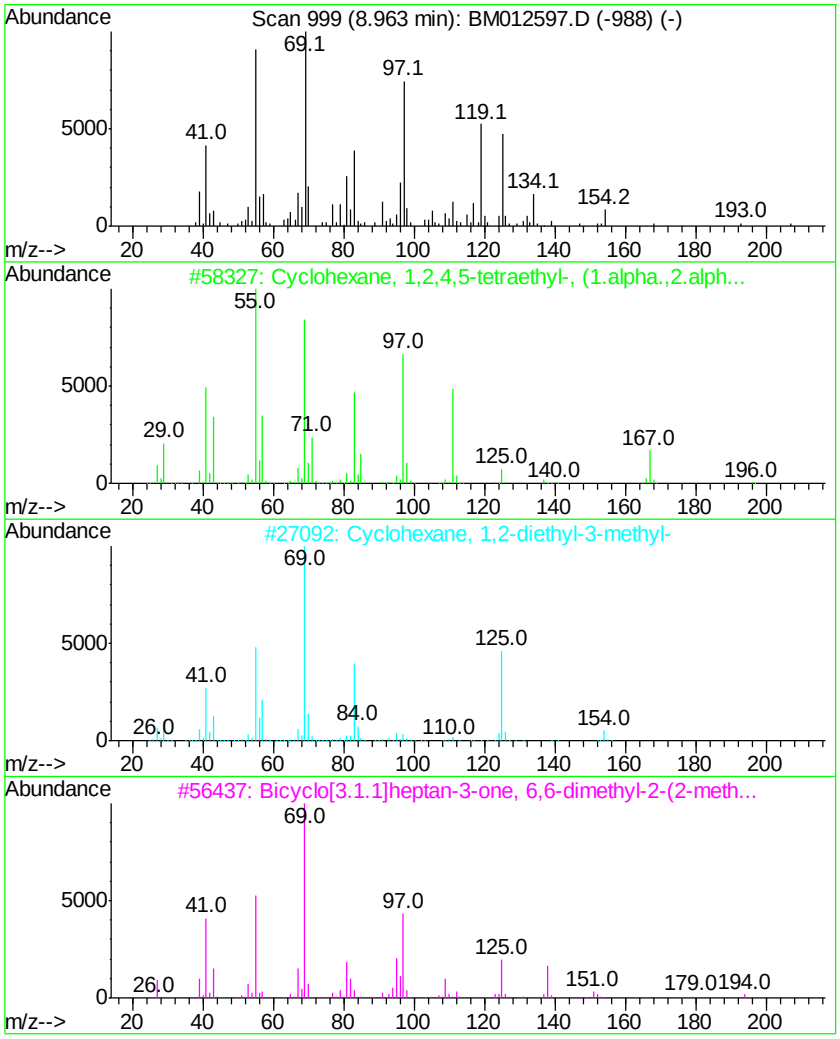
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

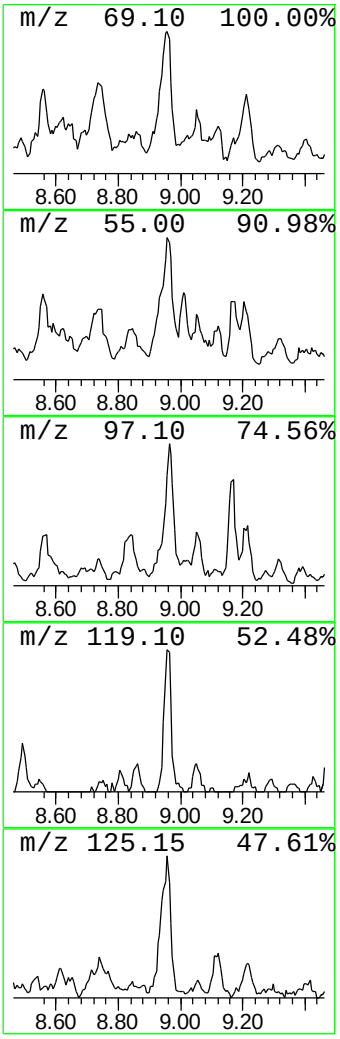
**********

Peak Number 10 (DEL) Alkane: Cyclic8.96 Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.96 | 3.70 ng/ul | 558171 | 1,4-Dichlorobenzene-d4 | 7.86 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclohexane, 1,2,4,5-tetraethyl-... | 196 | C14H28 | 061142-24-3 | 43 |
| 2 | Cyclohexane, 1,2-diethyl-3-methyl- | 154 | C11H22 | 061141-80-8 | 43 |
| 3 | Bicyclo[3.1.1]heptan-3-one, 6,6-... | 194 | C13H22O | 1000163-97-9 | 38 |
| 4 | 4-Undecene, 4-methyl-, (Z)- | 168 | C12H24 | 074630-57-2 | 37 |
| 5 | Cyclohexane, 1,1'-(1-methylethyl... | 208 | C15H28 | 054934-90-6 | 32 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

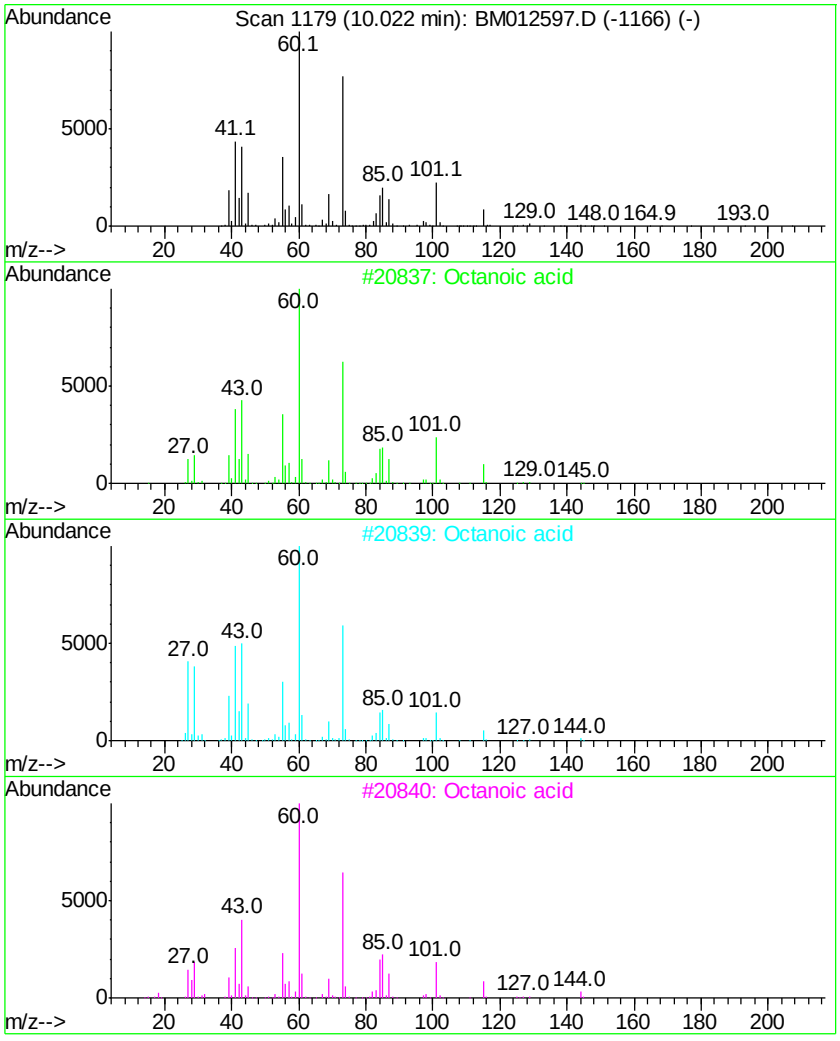
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

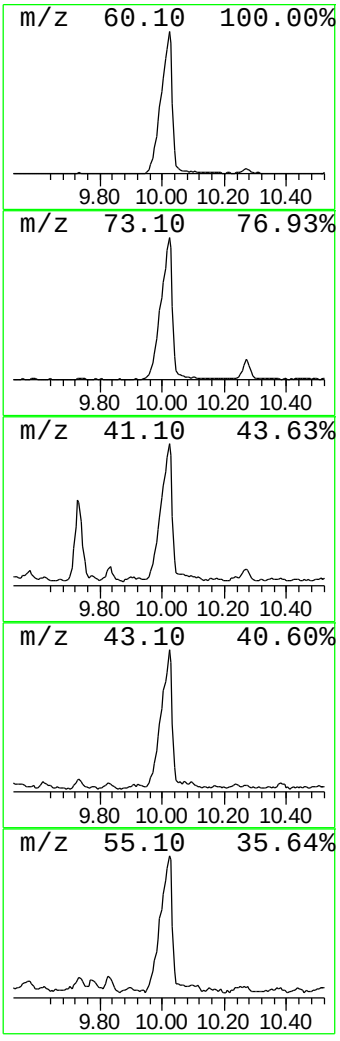
**********

Peak Number 11 Octanoic acid Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.02 | 17.07 ng/ul | 3010690 | Naphthalene-d8 | 10.65 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octanoic acid | 144 | C8H16O2 | 000124-07-2 | 91 |
| 2 | Octanoic acid | 144 | C8H16O2 | 000124-07-2 | 86 |
| 3 | Octanoic acid | 144 | C8H16O2 | 000124-07-2 | 80 |
| 4 | Hexanoic acid | 116 | C6H12O2 | 000142-62-1 | 53 |
| 5 | Hexanoic acid | 116 | C6H12O2 | 000142-62-1 | 53 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

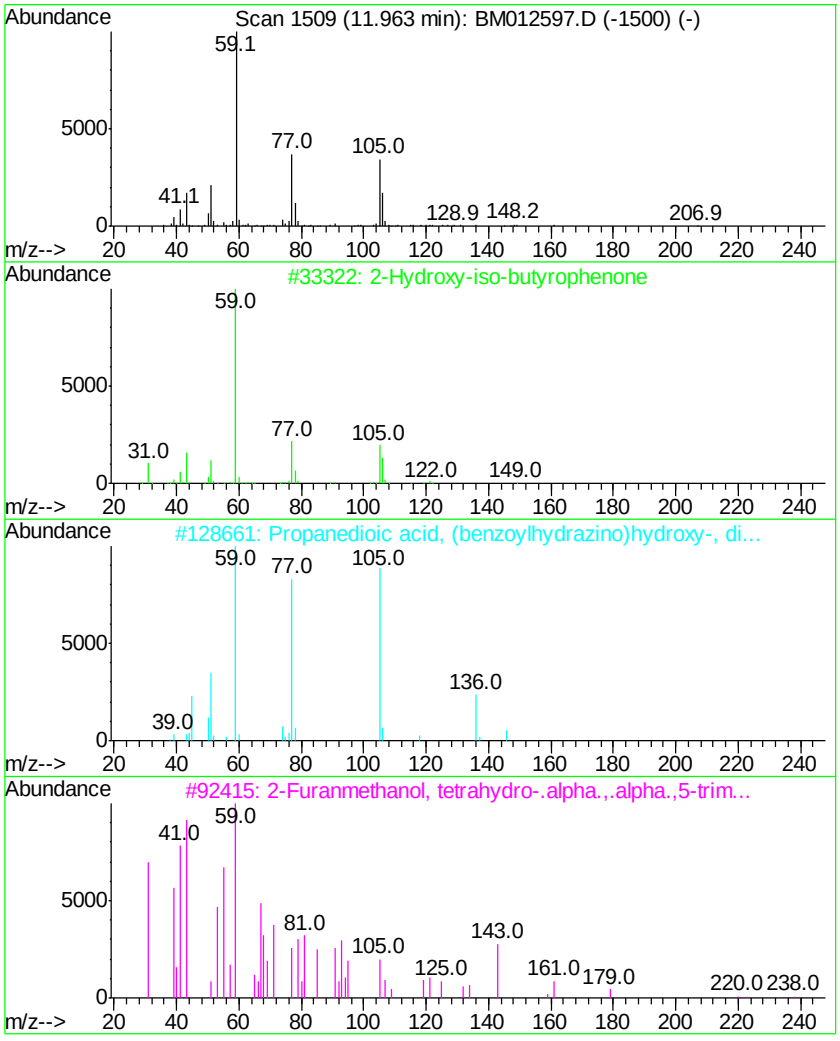
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

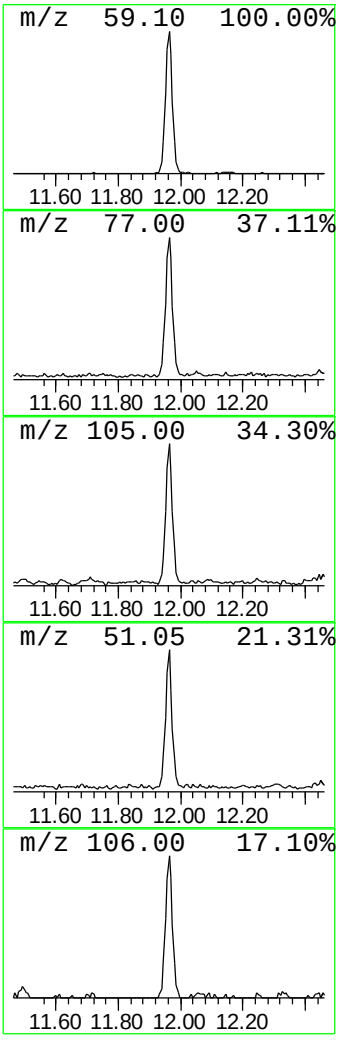
**********

Peak Number 12 2-Hydroxy-iso-butyrophenone Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.96 | 2.62 ng/ul | 461688 | Naphthalene-d8 | 10.65 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Hydroxy-iso-butyrophenone | 164 | C10H12O2 | 007473-98-5 | 78 |
| 2 | Propanedioic acid, (benzoylhydra... | 282 | C12H14N2O6 | 1000142-99-9 | 25 |
| 3 | 2-Furanmethanol, tetrahydro-.alp... | 238 | C15H26O2 | 026184-88-3 | 25 |
| 4 | Propionic acid, 2-isopropoxy-, m... | 146 | C7H14O3 | 1000151-15-1 | 23 |
| 5 | Propanoic acid, 2-hydroxy-2-meth... | 118 | C5H10O3 | 002110-78-3 | 23 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

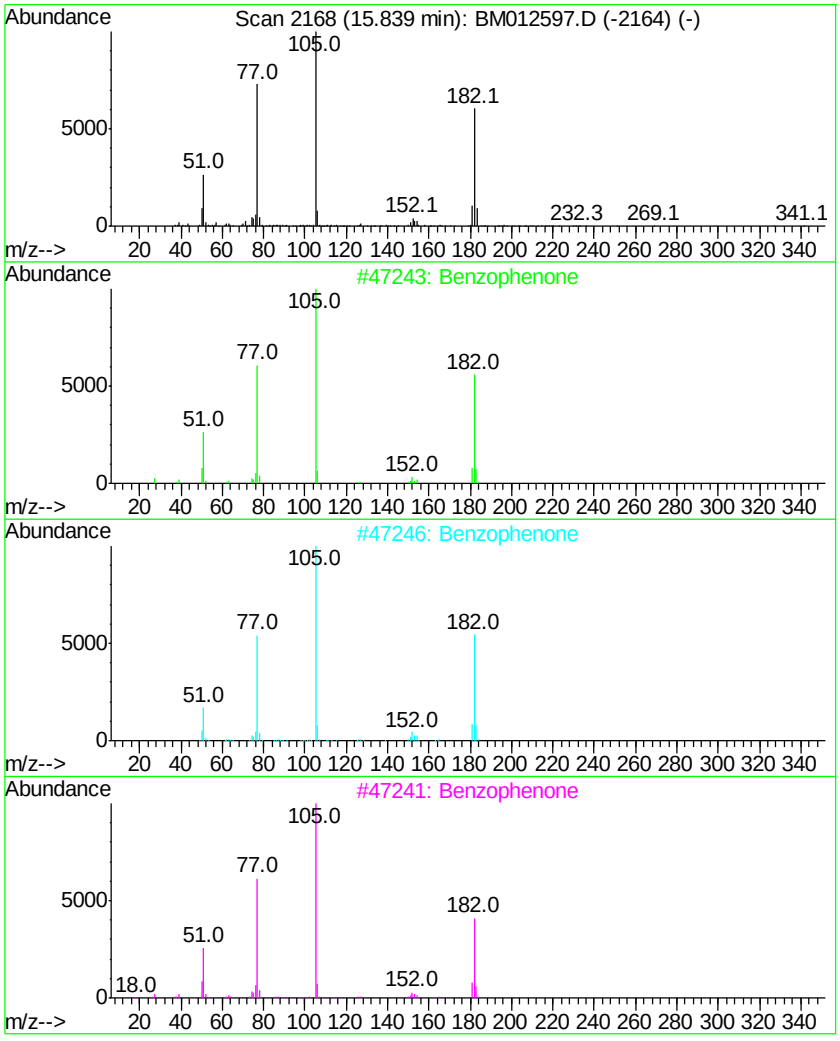
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

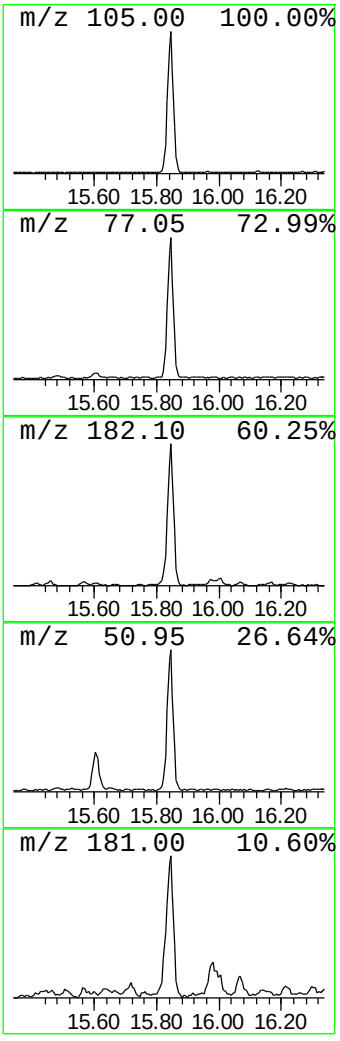
**********

Peak Number 13 Benzophenone Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.84 | 6.64 ng/ul | 1521470 | Acenaphthene-d10 | 14.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzophenone | 182 | C13H10O | 000119-61-9 | 96 |
| 2 | Benzophenone | 182 | C13H10O | 000119-61-9 | 95 |
| 3 | Benzophenone | 182 | C13H10O | 000119-61-9 | 95 |
| 4 | Benzophenone | 182 | C13H10O | 000119-61-9 | 91 |
| 5 | Benzophenone | 182 | C13H10O | 000119-61-9 | 81 |



# Library Search Compound Report

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Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

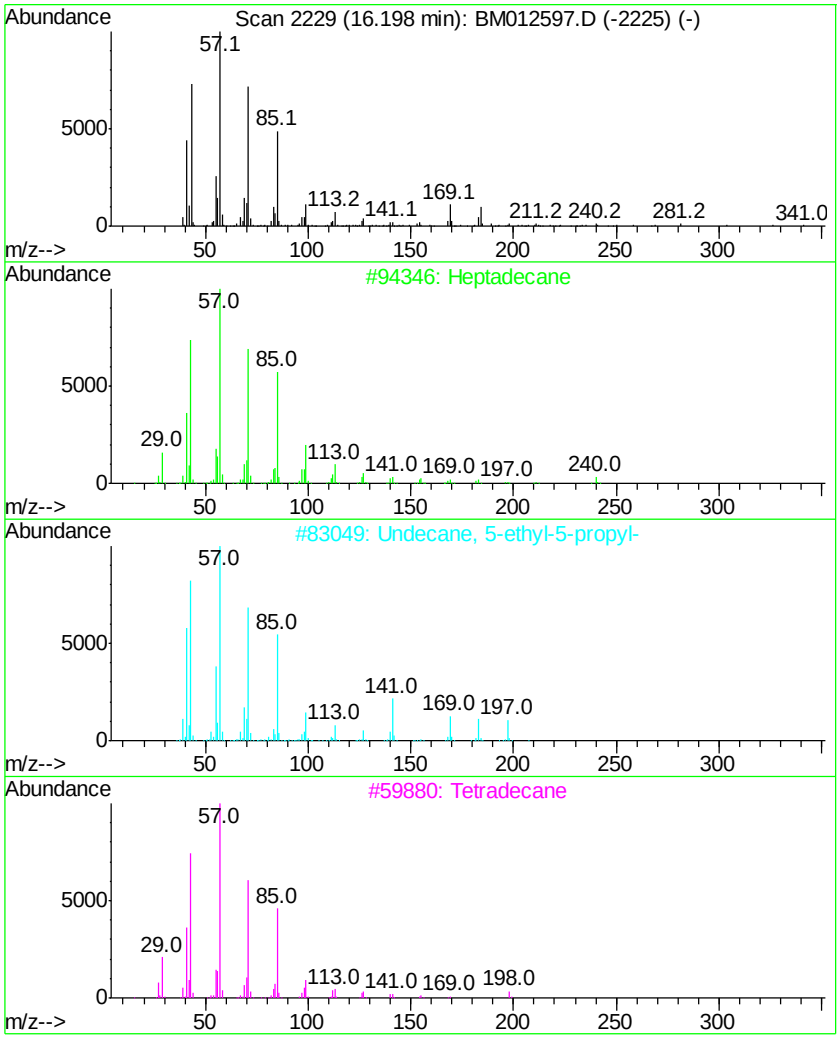
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

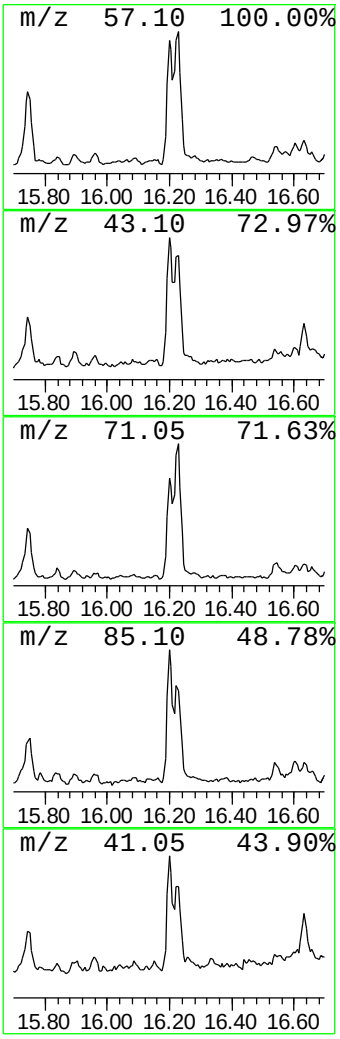
**********

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.20 | 2.91 ng/ul | 768016 | Phenanthrene-d10 | 17.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptadecane | 240 | C17H36 | 000629-78-7 | 95 |
| 2 | Undecane, 5-ethyl-5-propyl- | 226 | C16H34 | 002755-07-9 | 83 |
| 3 | Tetradecane | 198 | C14H30 | 000629-59-4 | 70 |
| 4 | Dodecane | 170 | C12H26 | 000112-40-3 | 70 |
| 5 | Eicosane | 282 | C20H42 | 000112-95-8 | 64 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

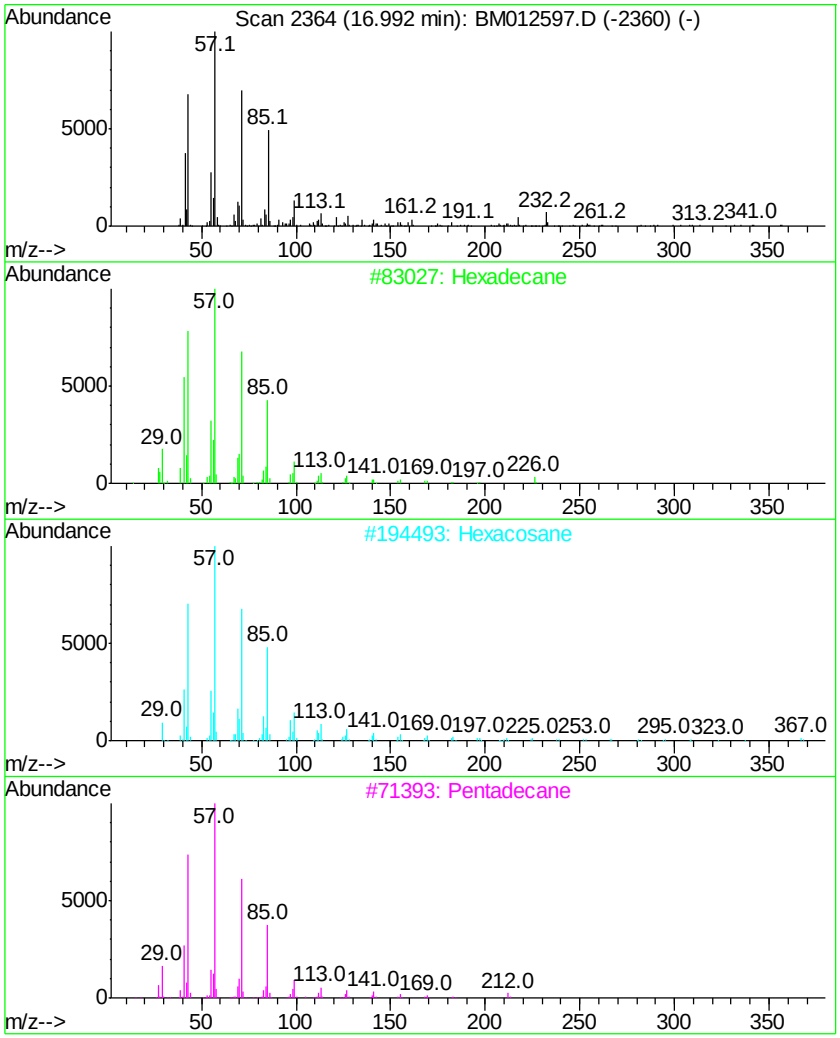
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

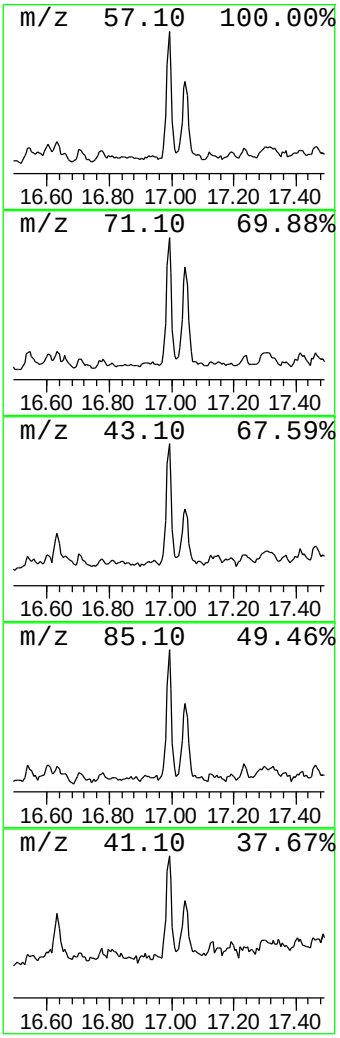
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.99 | 2.16 ng/ul | 570215 | Phenanthrene-d10 | 17.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexadecane | 226 | C16H34 | 000544-76-3 | 78 |
| 2 | Hexacosane | 366 | C26H54 | 000630-01-3 | 74 |
| 3 | Pentadecane | 212 | C15H32 | 000629-62-9 | 74 |
| 4 | Nonadecane | 268 | C19H40 | 000629-92-5 | 74 |
| 5 | Tridecane, 1-iodo- | 310 | C13H27I | 035599-77-0 | 74 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

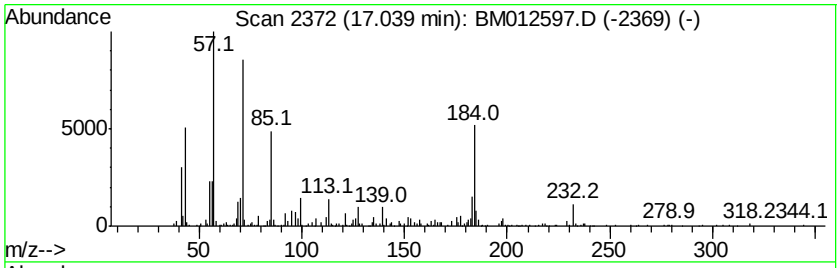
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

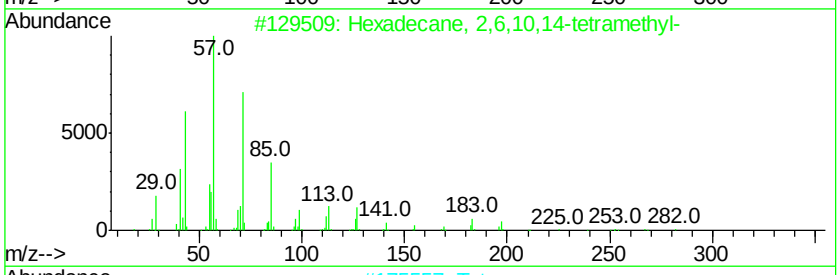
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

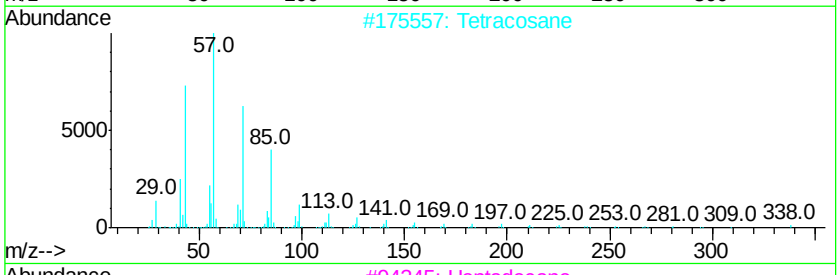
**********
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 13

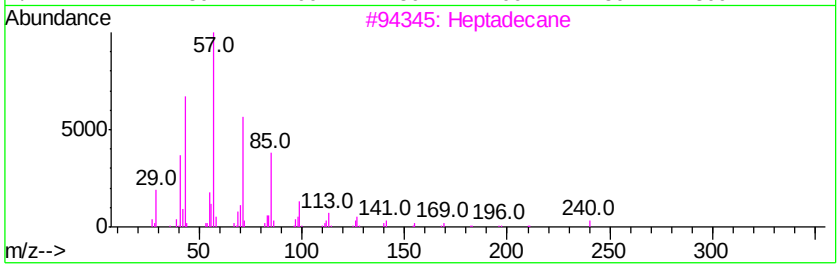
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.04 | 2.72 ng/ul | 718468 | Phenanthrene-d10 | 17.23 |

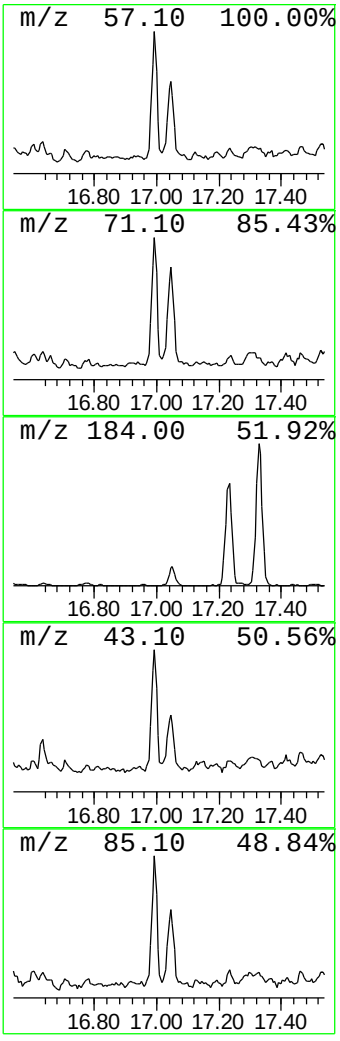
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42 | 000638-36-8 | 70 |
| 2 | Tetracosane | 338 | C24H50 | 000646-31-1 | 49 |
| 3 | Heptadecane | 240 | C17H36 | 000629-78-7 | 49 |
| 4 | Heptacosane | 380 | C27H56 | 000593-49-7 | 49 |
| 5 | Hexadecane | 226 | C16H34 | 000544-76-3 | 49 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

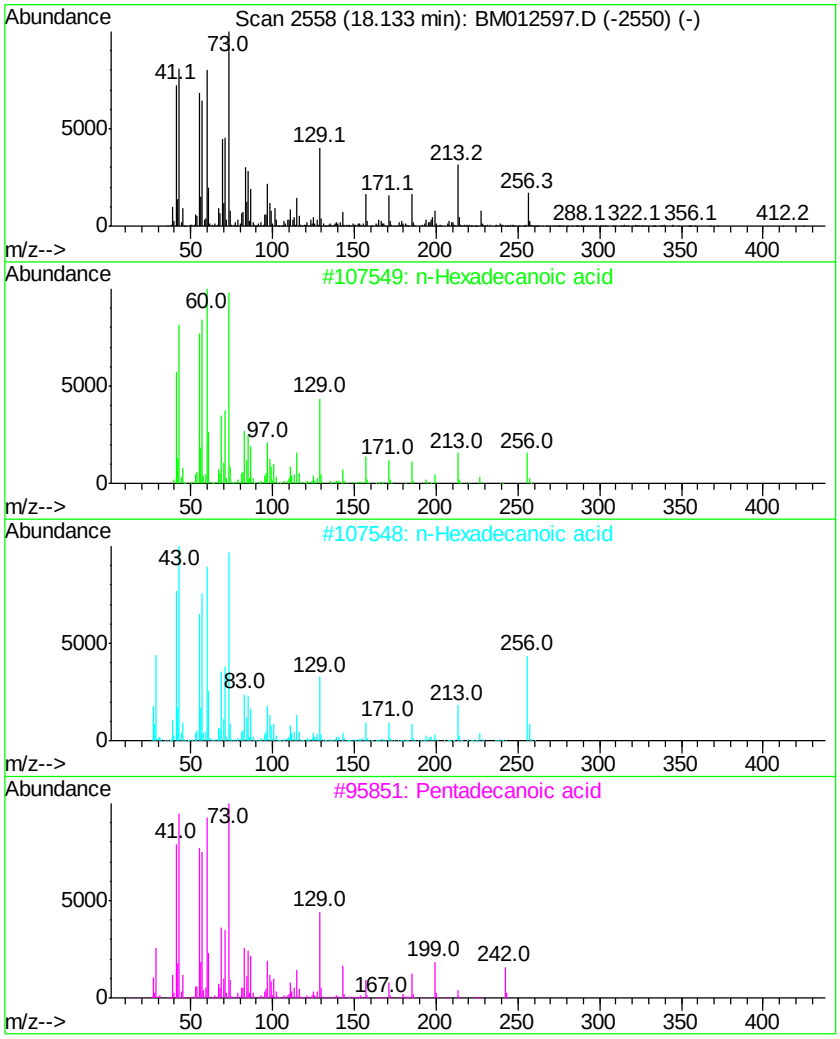
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

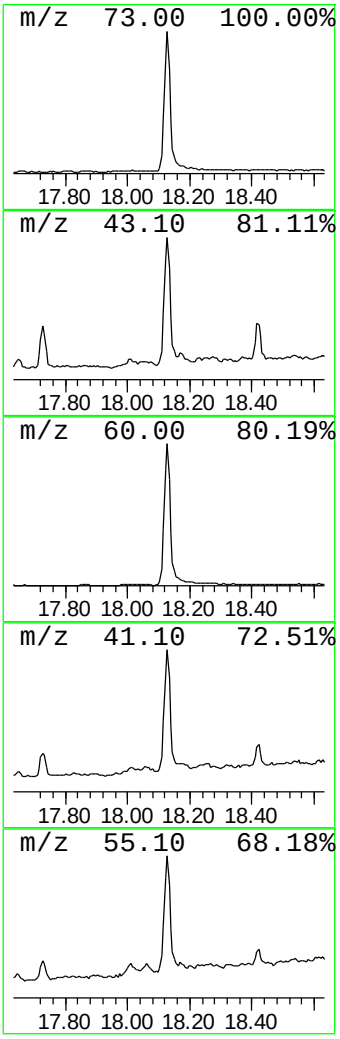
**********

Peak Number 17 n-Hexadecanoic acid Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.13 | 17.90 ng/ul | 4721820 | Phenanthrene-d10 | 17.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99 |
| 2 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98 |
| 3 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 91 |
| 4 | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 91 |
| 5 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 90 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

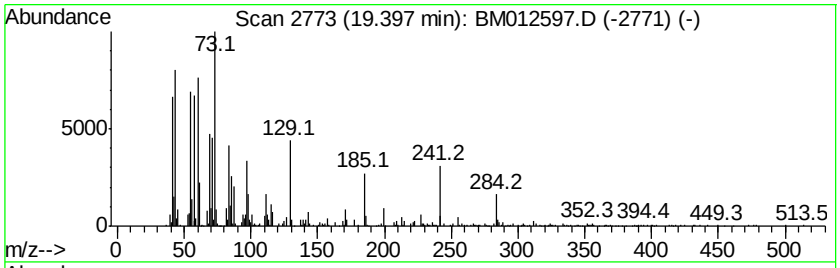
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

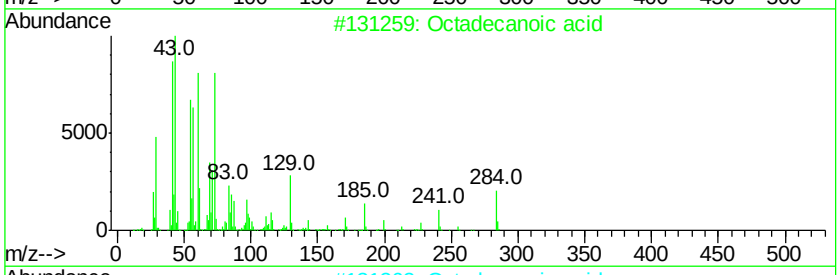
**********

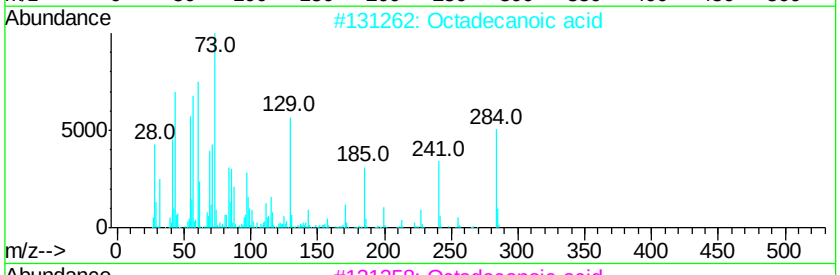
Peak Number 18 Octadecanoic acid Concentration Rank 19

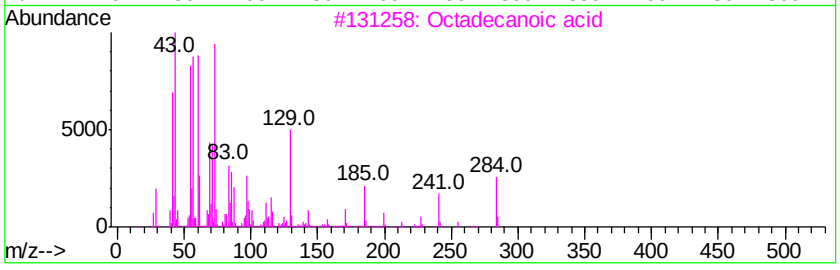
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.40 | 2.00 ng/ul | 787778 | Chrysene-d12 | 21.40 |

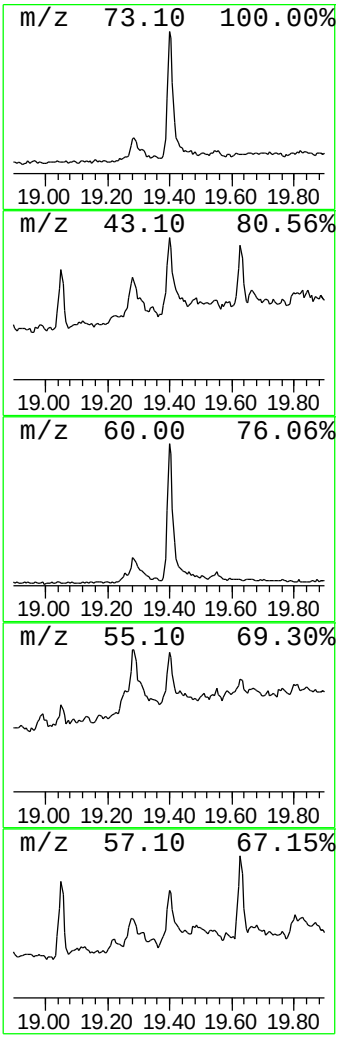
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 96 |
| 2 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 95 |
| 3 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 94 |
| 4 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 93 |
| 5 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 93 |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

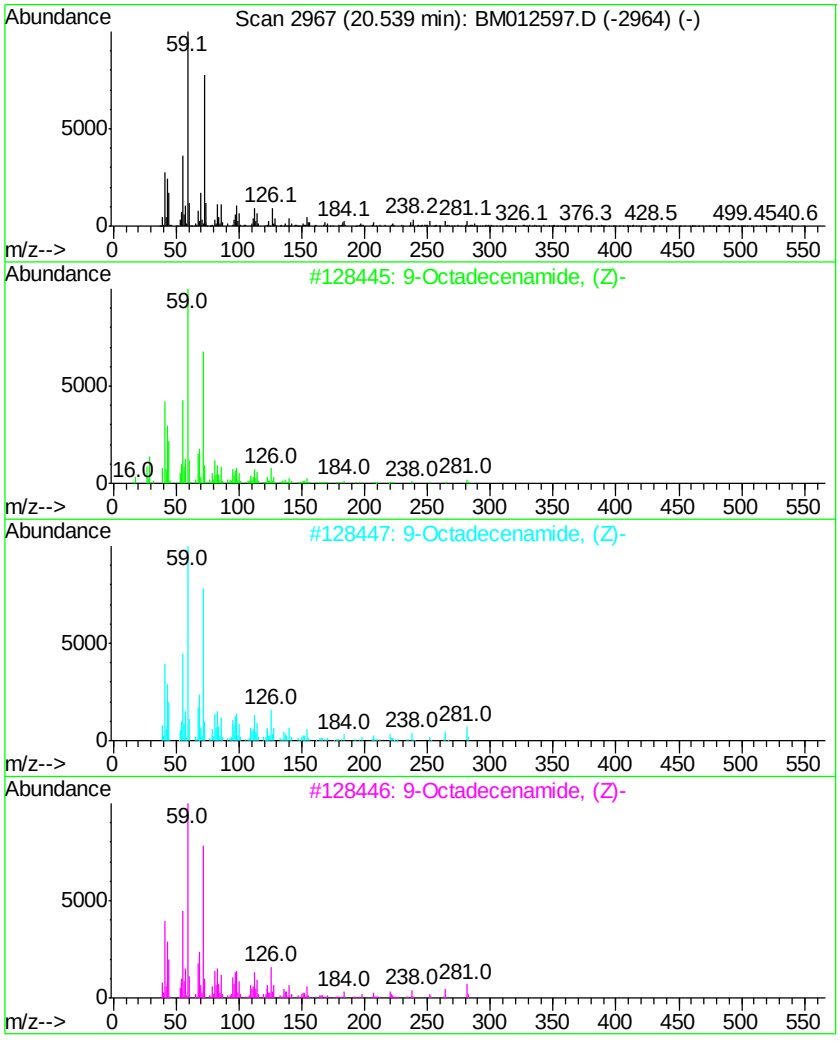
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

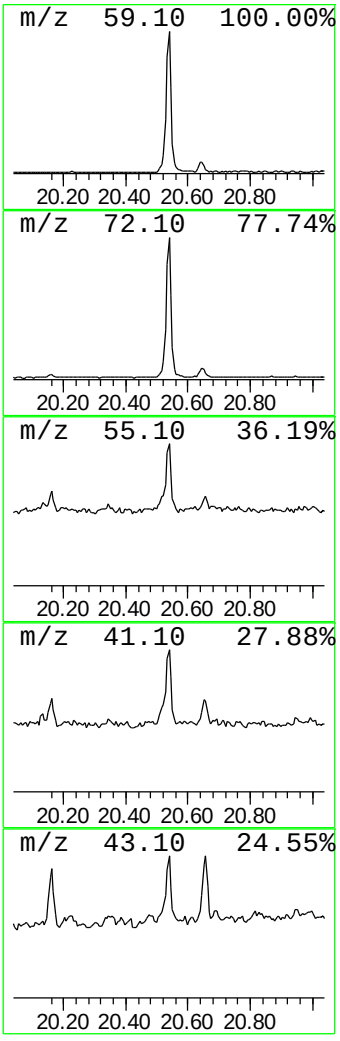
**********

Peak Number 19 9-Octadecenamide, (Z)- Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.54 | 4.71 ng/ul | 1850690 | Chrysene-d12 | 21.40 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9-Octadecenamide, (Z)- | 281 | C18H35NO | 000301-02-0 | 96 |
| 2 | 9-Octadecenamide, (Z)- | 281 | C18H35NO | 000301-02-0 | 91 |
| 3 | 9-Octadecenamide, (Z)- | 281 | C18H35NO | 000301-02-0 | 91 |
| 4 | 9-Octadecenamide, (Z)- | 281 | C18H35NO | 000301-02-0 | 72 |
| 5 | Benzeneethanamine, 2-fluoro-.bet... | 229 | C11H16FNO3 | 061338-98-5 | 64 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM103017\
Data File : BM012597.D
Acq On : 31 Oct 2017 05:22
Operator : SJ/JU
Sample : I5786-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** B-8(5-6.25)_101017

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA 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 |
| 2-Pentanone | 3.16 | 13.1 | ng/ul | 1973140 | 1 | 7.86 | 3020300 | 20.0 |
| Propanoic acid, 2... | 3.80 | 44.2 | ng/ul | 6676800 | 1 | 7.86 | 3020300 | 20.0 |
| Butanoic acid | 4.17 | 62.5 | ng/ul | 9431570 | 1 | 7.86 | 3020300 | 20.0 |
| 2-Heptanone | 5.72 | 19.0 | ng/ul | 2865400 | 1 | 7.86 | 3020300 | 20.0 |
| 2-Decanol | 5.87 | 2.1 | ng/ul | 324151 | 1 | 7.86 | 3020300 | 20.0 |
| (DEL) Alkane: Str... | 6.53 | 2.6 | ng/ul | 396978 | 1 | 7.86 | 3020300 | 20.0 |
| unknown-01 | 7.79 | 3.1 | ng/ul | 476373 | 1 | 7.86 | 3020300 | 20.0 |
| (DEL) Alkane: Str... | 8.05 | 2.3 | ng/ul | 346800 | 1 | 7.86 | 3020300 | 20.0 |
| (DEL) Alkane: Cyc... | 8.42 | 2.9 | ng/ul | 441485 | 1 | 7.86 | 3020300 | 20.0 |
| (DEL) Alkane: Cyc... | 8.96 | 3.7 | ng/ul | 558171 | 1 | 7.86 | 3020300 | 20.0 |
| Octanoic acid | 10.02 | 17.1 | ng/ul | 3010690 | 2 | 10.65 | 3527600 | 20.0 |
| 2-Hydroxy-iso-but... | 11.96 | 2.6 | ng/ul | 461688 | 2 | 10.65 | 3527600 | 20.0 |
| Benzophenone | 15.84 | 6.6 | ng/ul | 1521470 | 3 | 14.49 | 4585920 | 20.0 |
| (DEL) Alkane: Str... | 16.20 | 2.9 | ng/ul | 768016 | 4 | 17.23 | 5276350 | 20.0 |
| (DEL) Alkane: Str... | 16.99 | 2.2 | ng/ul | 570215 | 4 | 17.23 | 5276350 | 20.0 |
| (DEL) Alkane: Str... | 17.04 | 2.7 | ng/ul | 718468 | 4 | 17.23 | 5276350 | 20.0 |
| n-Hexadecanoic acid | 18.13 | 17.9 | ng/ul | 4721820 | 4 | 17.23 | 5276350 | 20.0 |
| Octadecanoic acid | 19.40 | 2.0 | ng/ul | 787778 | 5 | 21.40 | 7859790 | 20.0 |
| 9-Octadecenamide,... | 20.54 | 4.7 | ng/ul | 1850690 | 5 | 21.40 | 7859790 | 20.0 |