LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.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 | 3.520 | 27 | 31 | 33 | rBV2 | 15022 | 20512 | 1.09% | 0.078% |
| 2 | 3.556 | 33 | 37 | 44 | rVB | 40803 | 62378 | 3.31% | 0.236% |
| 3 | 4.713 | 226 | 234 | 241 | rBV | 50501 | 91632 | 4.86% | 0.347% |
| 4 | 5.406 | 346 | 352 | 367 | rBV | 161385 | 267631 | 14.20% | 1.014% |
| 5 | 6.581 | 546 | 552 | 558 | rBV3 | 14903 | 25939 | 1.38% | 0.098% |
| 6 | 7.245 | 658 | 665 | 677 | rBV2 | 65693 | 141748 | 7.52% | 0.537% |
| 7 | 7.416 | 687 | 694 | 704 | rBV | 159188 | 272845 | 14.47% | 1.033% |
| 8 | 7.627 | 722 | 730 | 735 | rVV2 | 190094 | 341404 | 18.11% | 1.293% |
| 9 | 7.674 | 735 | 738 | 743 | rVV2 | 28578 | 52197 | 2.77% | 0.198% |
| 10 | 7.745 | 747 | 750 | 756 | rVB6 | 8410 | 14891 | 0.79% | 0.056% |
| 11 | 7.992 | 787 | 792 | 797 | rVB7 | 9276 | 19175 | 1.02% | 0.073% |
| 12 | 8.097 | 804 | 810 | 818 | rBV | 250960 | 436127 | 23.13% | 1.652% |
| 13 | 8.473 | 865 | 874 | 881 | rVB | 100748 | 177980 | 9.44% | 0.674% |
| 14 | 8.744 | 913 | 920 | 922 | rBV5 | 10959 | 19987 | 1.06% | 0.076% |
| 15 | 8.791 | 922 | 928 | 937 | rVB2 | 174382 | 302927 | 16.07% | 1.147% |
| 16 | 8.990 | 957 | 962 | 966 | rVB2 | 11947 | 16314 | 0.87% | 0.062% |
| 17 | 9.079 | 969 | 977 | 984 | rBV | 220478 | 384381 | 20.39% | 1.456% |
| 18 | 9.202 | 992 | 998 | 1002 | rBV3 | 28008 | 50292 | 2.67% | 0.190% |
| 19 | 9.255 | 1002 | 1007 | 1020 | rVV | 218118 | 398518 | 21.14% | 1.509% |
| 20 | 9.455 | 1036 | 1041 | 1049 | rVB10 | 11675 | 25194 | 1.34% | 0.095% |
| 21 | 9.613 | 1062 | 1068 | 1076 | rBV8 | 28969 | 55325 | 2.93% | 0.210% |
| 22 | 9.678 | 1076 | 1079 | 1083 | rVV5 | 11311 | 17312 | 0.92% | 0.066% |
| 23 | 9.731 | 1083 | 1088 | 1091 | rVB | 15083 | 23121 | 1.23% | 0.088% |
| 24 | 9.848 | 1103 | 1108 | 1112 | rBV6 | 10578 | 17148 | 0.91% | 0.065% |
| 25 | 9.977 | 1124 | 1130 | 1138 | rBV | 196131 | 354217 | 18.79% | 1.342% |
| 26 | 10.107 | 1144 | 1152 | 1158 | rBV10 | 15240 | 41665 | 2.21% | 0.158% |
| 27 | 10.312 | 1181 | 1187 | 1193 | rVV7 | 20524 | 46386 | 2.46% | 0.176% |
| 28 | 10.377 | 1195 | 1198 | 1203 | rVV7 | 10671 | 20153 | 1.07% | 0.076% |
| 29 | 10.442 | 1203 | 1209 | 1216 | rVB6 | 27336 | 67173 | 3.56% | 0.254% |
| 30 | 10.518 | 1216 | 1222 | 1241 | rVB | 330142 | 649200 | 34.44% | 2.459% |
| 31 | 10.800 | 1267 | 1270 | 1277 | rVB4 | 22122 | 40279 | 2.14% | 0.153% |
| 32 | 10.906 | 1281 | 1288 | 1293 | rVV | 319017 | 572861 | 30.39% | 2.170% |
| 33 | 10.959 | 1293 | 1297 | 1303 | rVV2 | 78866 | 137885 | 7.31% | 0.522% |
| 34 | 11.035 | 1303 | 1310 | 1321 | rVB | 241912 | 484632 | 25.71% | 1.836% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 11.158 | 1325 | 1331 | 1333 | rBV7 | 14381 | 22353 | 1.19% | 0.085% |
| 36 | 11.687 | 1413 | 1421 | 1426 | rBV8 | 15524 | 47957 | 2.54% | 0.182% |
| 37 | 11.728 | 1426 | 1428 | 1435 | rVV7 | 13381 | 26909 | 1.43% | 0.102% |
| 38 | 11.852 | 1446 | 1449 | 1453 | rVB6 | 10458 | 16873 | 0.90% | 0.064% |
| 39 | 12.099 | 1486 | 1491 | 1497 | rVB9 | 10646 | 20278 | 1.08% | 0.077% |
| 40 | 12.228 | 1505 | 1513 | 1519 | rBV2 | 84284 | 152448 | 8.09% | 0.577% |
| 41 | 12.392 | 1535 | 1541 | 1546 | rVV3 | 79670 | 147713 | 7.84% | 0.559% |
| 42 | 12.451 | 1547 | 1551 | 1553 | rVV4 | 15774 | 25618 | 1.36% | 0.097% |
| 43 | 12.492 | 1553 | 1558 | 1565 | rVV7 | 30522 | 72525 | 3.85% | 0.275% |
| 44 | 12.557 | 1565 | 1569 | 1572 | rVB6 | 10361 | 16744 | 0.89% | 0.063% |
| 45 | 12.610 | 1572 | 1578 | 1580 | rBV6 | 10945 | 17817 | 0.95% | 0.067% |
| 46 | 12.762 | 1597 | 1604 | 1609 | rVB10 | 26980 | 51942 | 2.76% | 0.197% |
| 47 | 12.874 | 1618 | 1623 | 1628 | rBV6 | 20266 | 32362 | 1.72% | 0.123% |
| 48 | 12.974 | 1636 | 1640 | 1645 | rVB6 | 15923 | 26446 | 1.40% | 0.100% |
| 49 | 13.062 | 1650 | 1655 | 1658 | rBV7 | 11813 | 19191 | 1.02% | 0.073% |
| 50 | 13.520 | 1727 | 1733 | 1736 | rBV8 | 15811 | 34739 | 1.84% | 0.132% |
| 51 | 13.609 | 1743 | 1748 | 1756 | rVB | 334014 | 543815 | 28.85% | 2.060% |
| 52 | 13.773 | 1764 | 1776 | 1778 | rBV9 | 18038 | 49525 | 2.63% | 0.188% |
| 53 | 13.808 | 1778 | 1782 | 1786 | rVB7 | 19502 | 35186 | 1.87% | 0.133% |
| 54 | 13.891 | 1793 | 1796 | 1798 | rBV4 | 14764 | 18508 | 0.98% | 0.070% |
| 55 | 14.049 | 1816 | 1823 | 1827 | rBV7 | 34210 | 85829 | 4.55% | 0.325% |
| 56 | 14.120 | 1829 | 1835 | 1840 | rVV | 633597 | 904265 | 47.97% | 3.425% |
| 57 | 14.167 | 1840 | 1843 | 1847 | rVB | 113255 | 147878 | 7.84% | 0.560% |
| 58 | 14.267 | 1857 | 1860 | 1865 | rBV7 | 24101 | 42104 | 2.23% | 0.159% |
| 59 | 14.366 | 1874 | 1877 | 1878 | rBV3 | 16891 | 14793 | 0.78% | 0.056% |
| 60 | 14.413 | 1879 | 1885 | 1895 | rVB | 629246 | 1023658 | 54.30% | 3.877% |
| 61 | 14.519 | 1899 | 1903 | 1909 | rBV5 | 21657 | 38163 | 2.02% | 0.145% |
| 62 | 14.643 | 1920 | 1924 | 1928 | rBV7 | 38886 | 58142 | 3.08% | 0.220% |
| 63 | 14.719 | 1931 | 1937 | 1944 | rVV | 506000 | 835405 | 44.31% | 3.164% |
| 64 | 14.784 | 1944 | 1948 | 1954 | rVB2 | 198596 | 291024 | 15.44% | 1.102% |
| 65 | 14.848 | 1955 | 1959 | 1964 | rVV | 29266 | 49334 | 2.62% | 0.187% |
| 66 | 14.907 | 1965 | 1969 | 1977 | rVB3 | 54685 | 88829 | 4.71% | 0.336% |
| 67 | 15.119 | 2000 | 2005 | 2011 | rVB | 75559 | 124316 | 6.59% | 0.471% |
| 68 | 15.324 | 2035 | 2040 | 2041 | rBV5 | 17213 | 30526 | 1.62% | 0.116% |
| 69 | 15.448 | 2057 | 2061 | 2070 | rBV | 122926 | 192485 | 10.21% | 0.729% |
| 70 | 15.712 | 2099 | 2106 | 2111 | rBV | 832679 | 1283714 | 68.09% | 4.862% |
| 71 | 15.765 | 2111 | 2115 | 2119 | rVV2 | 93405 | 139008 | 7.37% | 0.527% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 15.818 | 2119 | 2124 | 2132 | rVV2 | 191084 | 317066 | 16.82% | 1.201% |
| 73 | 15.918 | 2138 | 2141 | 2144 | rBV | 31248 | 36224 | 1.92% | 0.137% |
| 74 | 15.965 | 2146 | 2149 | 2155 | rVB4 | 47820 | 78775 | 4.18% | 0.298% |
| 75 | 16.217 | 2186 | 2192 | 2197 | rBV | 546415 | 816820 | 43.33% | 3.094% |
| 76 | 16.394 | 2214 | 2222 | 2227 | rBV | 251772 | 431998 | 22.91% | 1.636% |
| 77 | 16.546 | 2244 | 2248 | 2251 | rVB2 | 23355 | 29767 | 1.58% | 0.113% |
| 78 | 16.728 | 2273 | 2279 | 2283 | rBV | 332238 | 491965 | 26.10% | 1.863% |
| 79 | 16.987 | 2320 | 2323 | 2334 | rVB2 | 49086 | 92278 | 4.89% | 0.350% |
| 80 | 17.263 | 2365 | 2370 | 2376 | rVB4 | 67827 | 145253 | 7.70% | 0.550% |
| 81 | 17.463 | 2399 | 2404 | 2410 | rVB | 685942 | 1039891 | 55.16% | 3.939% |
| 82 | 17.563 | 2415 | 2421 | 2429 | rVB | 963010 | 1473207 | 78.14% | 5.580% |
| 83 | 17.862 | 2469 | 2472 | 2475 | rBV | 31343 | 41624 | 2.21% | 0.158% |
| 84 | 18.350 | 2550 | 2555 | 2562 | rBV | 326206 | 467860 | 24.82% | 1.772% |
| 85 | 18.456 | 2570 | 2573 | 2578 | rVB5 | 33488 | 42193 | 2.24% | 0.160% |
| 86 | 19.102 | 2680 | 2683 | 2685 | rBV4 | 26623 | 34410 | 1.83% | 0.130% |
| 87 | 19.472 | 2743 | 2746 | 2750 | rBV4 | 27085 | 50713 | 2.69% | 0.192% |
| 88 | 19.549 | 2756 | 2759 | 2764 | rVB2 | 48592 | 58746 | 3.12% | 0.223% |
| 89 | 19.613 | 2766 | 2770 | 2779 | rVB3 | 158499 | 232371 | 12.33% | 0.880% |
| 90 | 19.848 | 2804 | 2810 | 2813 | rBV | 1279178 | 1885236 | 100.00% | 7.140% |
| 91 | 19.878 | 2813 | 2815 | 2826 | rVB | 585186 | 842875 | 44.71% | 3.192% |
| 92 | 20.242 | 2875 | 2877 | 2883 | rVB2 | 40284 | 46476 | 2.47% | 0.176% |
| 93 | 20.307 | 2884 | 2888 | 2893 | rVV | 86245 | 109848 | 5.83% | 0.416% |
| 94 | 21.376 | 3066 | 3070 | 3080 | rVB2 | 218859 | 374740 | 19.88% | 1.419% |
| 95 | 21.740 | 3125 | 3132 | 3139 | rVB | 816158 | 1340044 | 71.08% | 5.075% |
| 96 | 24.073 | 3524 | 3529 | 3535 | rVB6 | 34333 | 60958 | 3.23% | 0.231% |
| 97 | 24.778 | 3639 | 3649 | 3661 | rVB | 687872 | 1853229 | 98.30% | 7.019% |
| 98 | 25.007 | 3676 | 3688 | 3700 | rVB2 | 482732 | 1450760 | 76.95% | 5.495% |
| 99 | 26.194 | 3883 | 3890 | 3900 | rVB6 | 31641 | 90479 | 4.80% | 0.343% |
| 100 | 27.575 | 4117 | 4125 | 4135 | rVB8 | 27546 | 88565 | 4.70% | 0.335% |

Sum of corrected areas: 26402222

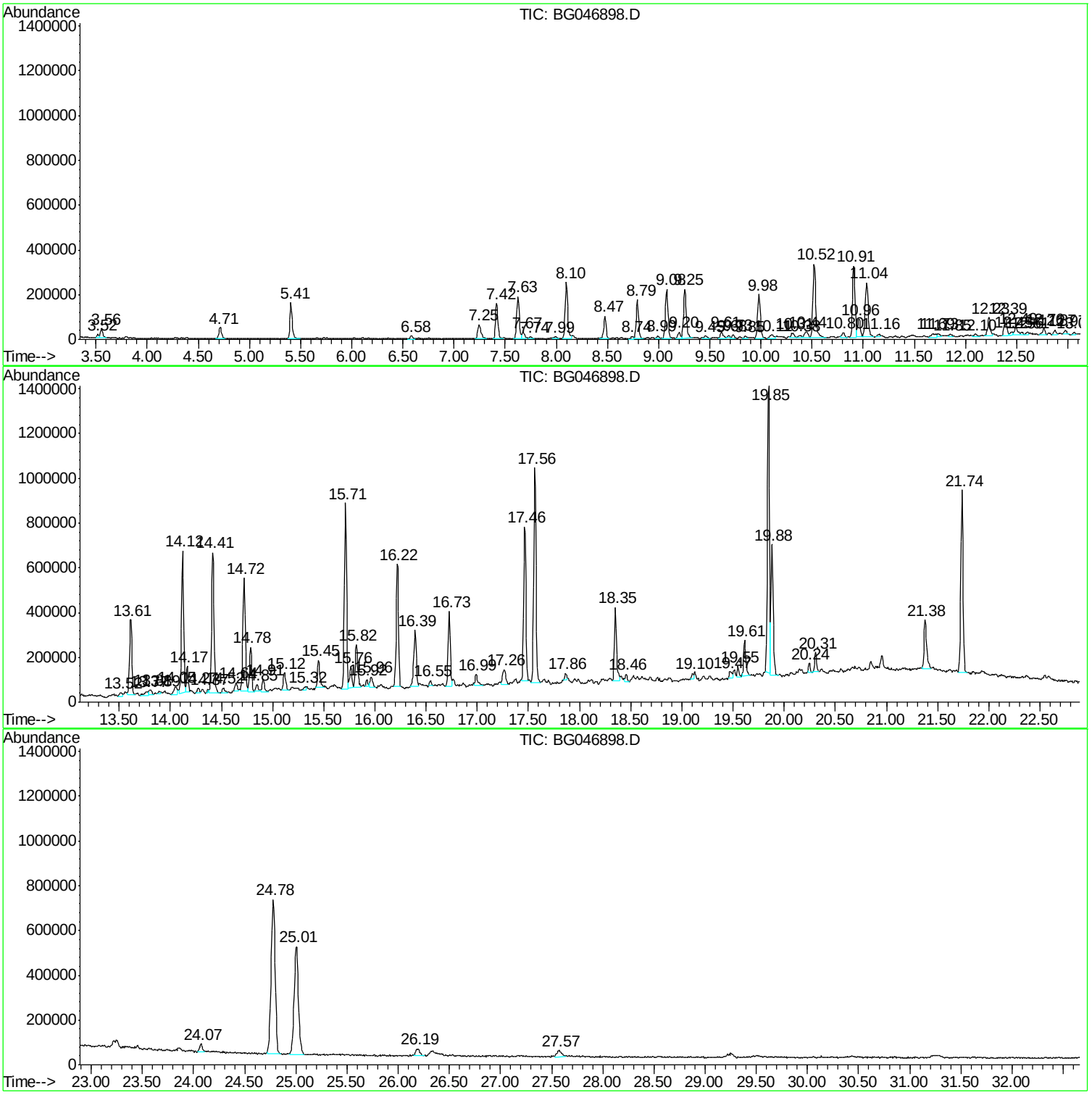
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

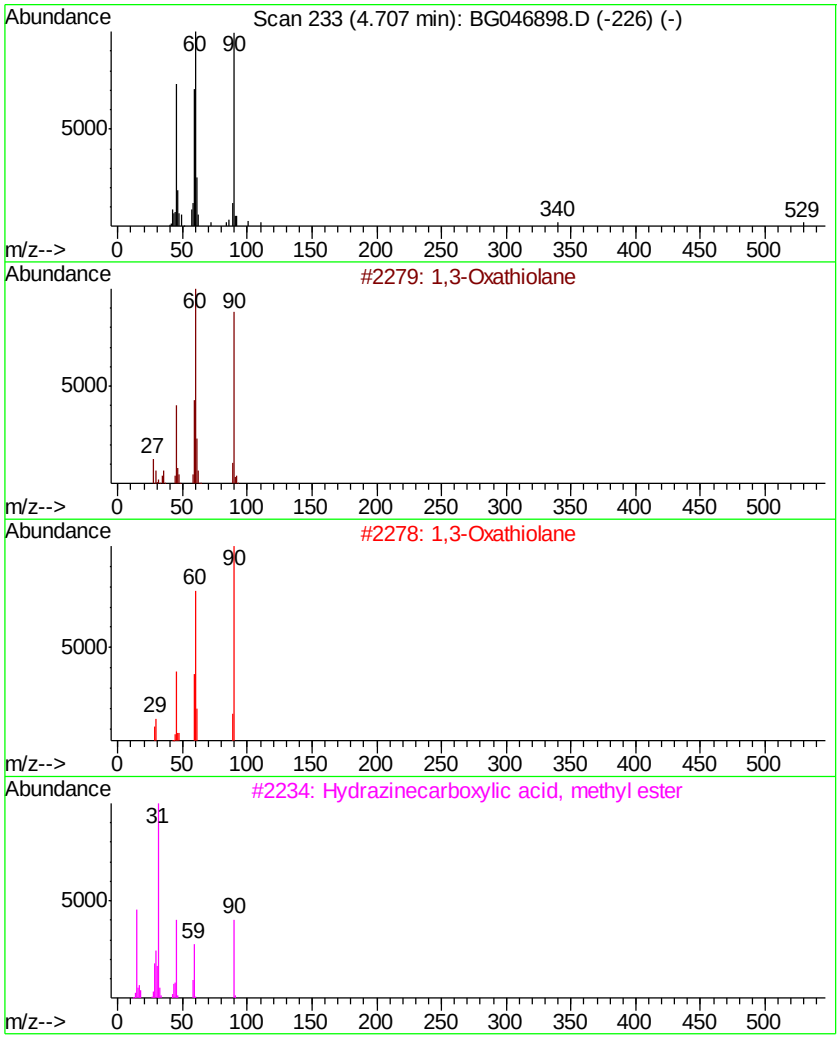
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

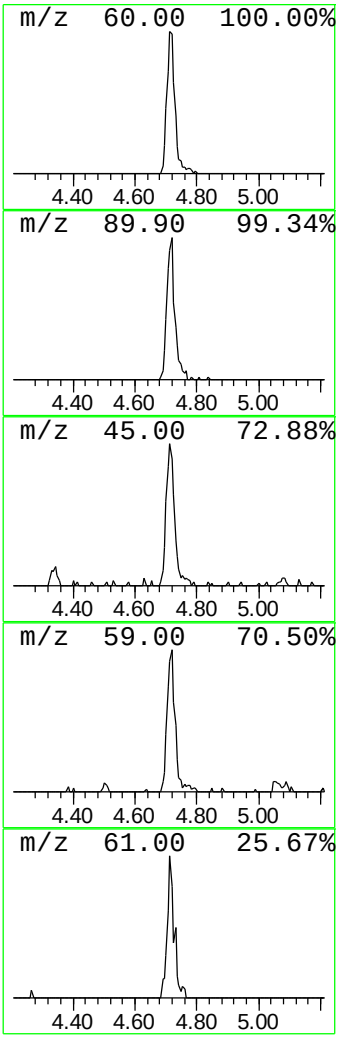
**********

Peak Number 1 1,3-Oxathiolane Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.71 | 4.20 ng/ul | 91632 | 1,4-Dichlorobenzene-d4 | 8.10 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,3-Oxathiolane | 90 | C3H6OS | 002094-97-5 | 86 |
| 2 | 1,3-Oxathiolane | 90 | C3H6OS | 002094-97-5 | 74 |
| 3 | Hydrazinecarboxylic acid, methyl... | 90 | C2H6N2O2 | 006294-89-9 | 59 |
| 4 | Thiourea, methyl- | 90 | C2H6N2S | 000598-52-7 | 50 |
| 5 | 3-Thietanol | 90 | C3H6OS | 010304-16-2 | 42 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

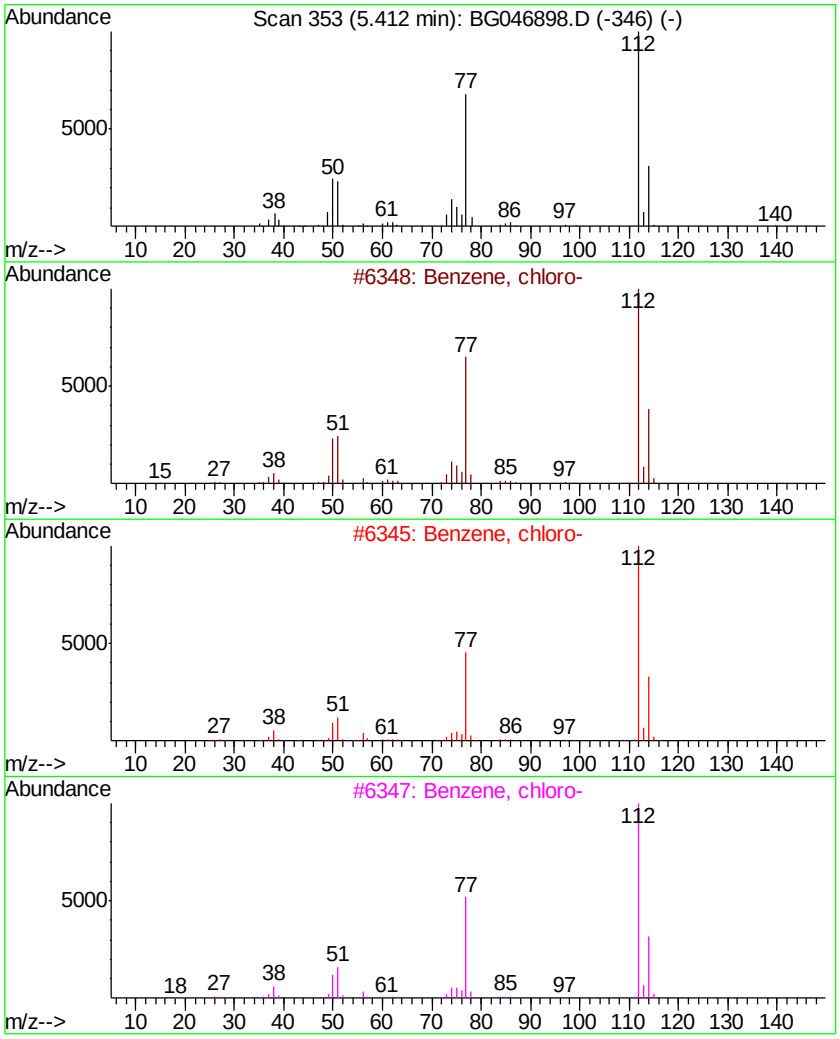
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

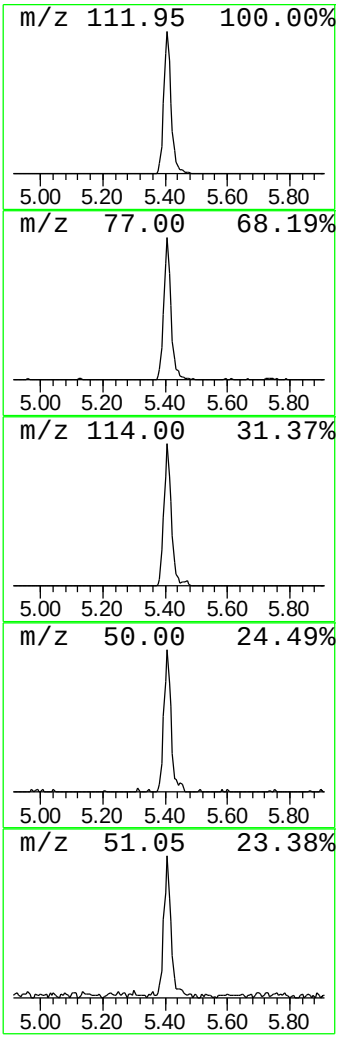
**********

Peak Number 2 Benzene, chloro- Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.41 | 12.27 ng/ul | 267631 | 1,4-Dichlorobenzene-d4 | 8.10 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, chloro- | 112 | C6H5Cl | 000108-90-7 | 96 |
| 2 | Benzene, chloro- | 112 | C6H5Cl | 000108-90-7 | 91 |
| 3 | Benzene, chloro- | 112 | C6H5Cl | 000108-90-7 | 91 |
| 4 | Benzene, chloro- | 112 | C6H5Cl | 000108-90-7 | 90 |
| 5 | 2-Chloro-3-nitropyridine | 158 | C5H3ClN2O2 | 005470-18-8 | 45 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

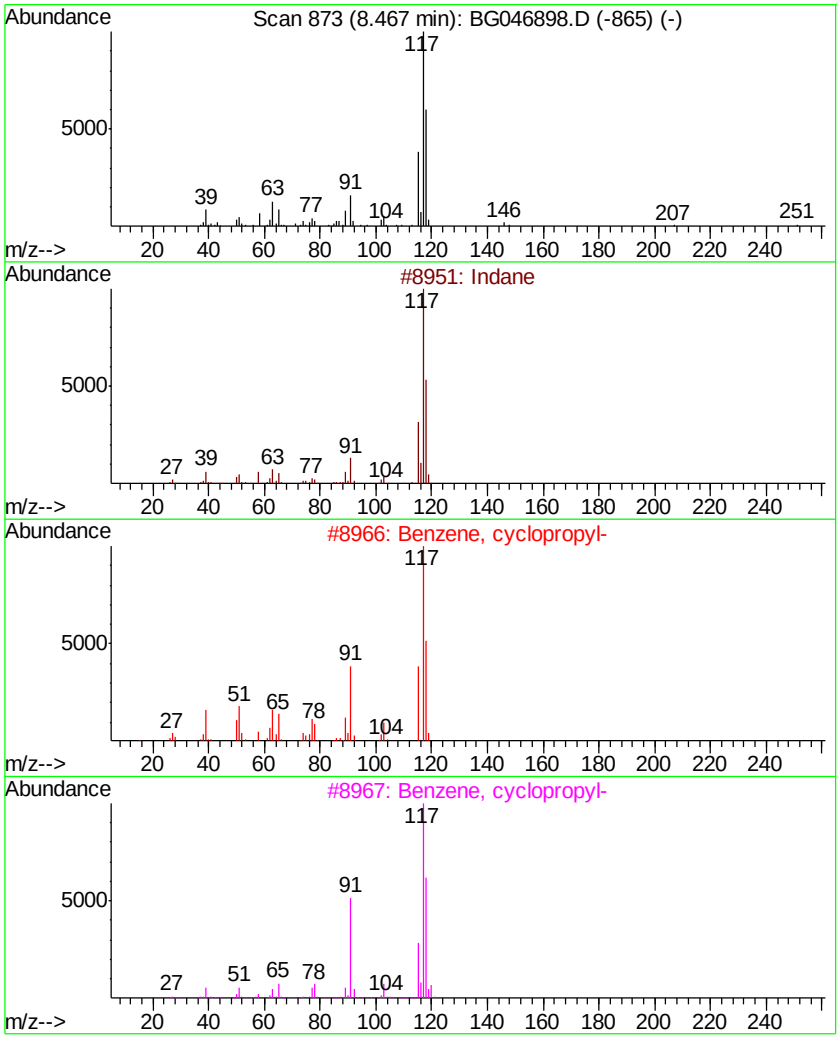
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

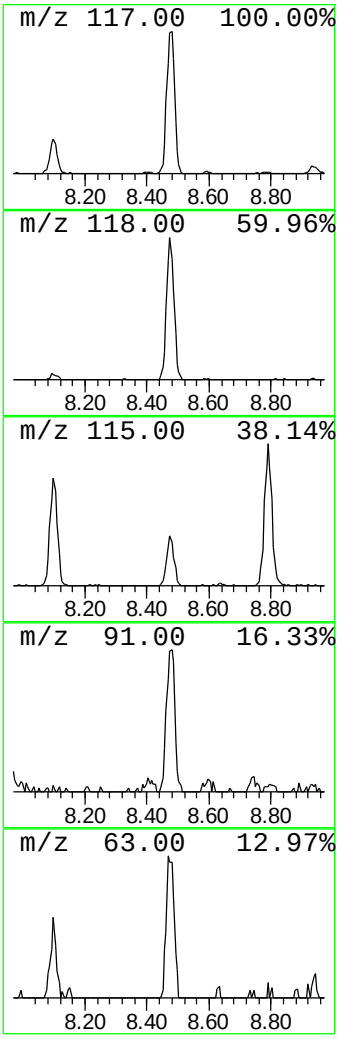
**********

Peak Number 3 Indane Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.47 | 8.16 ng/ul | 177980 | 1,4-Dichlorobenzene-d4 | 8.10 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Indane | 118 | C9H10 | 000496-11-7 | 87 |
| 2 | Benzene, cyclopropyl- | 118 | C9H10 | 000873-49-4 | 68 |
| 3 | Benzene, cyclopropyl- | 118 | C9H10 | 000873-49-4 | 64 |
| 4 | Benzene, 1-ethenyl-2-methyl- | 118 | C9H10 | 000611-15-4 | 64 |
| 5 | 1H-Indene-5-carboxylic acid, 2,3... | 162 | C10H10O2 | 1000337-61-3 | 59 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

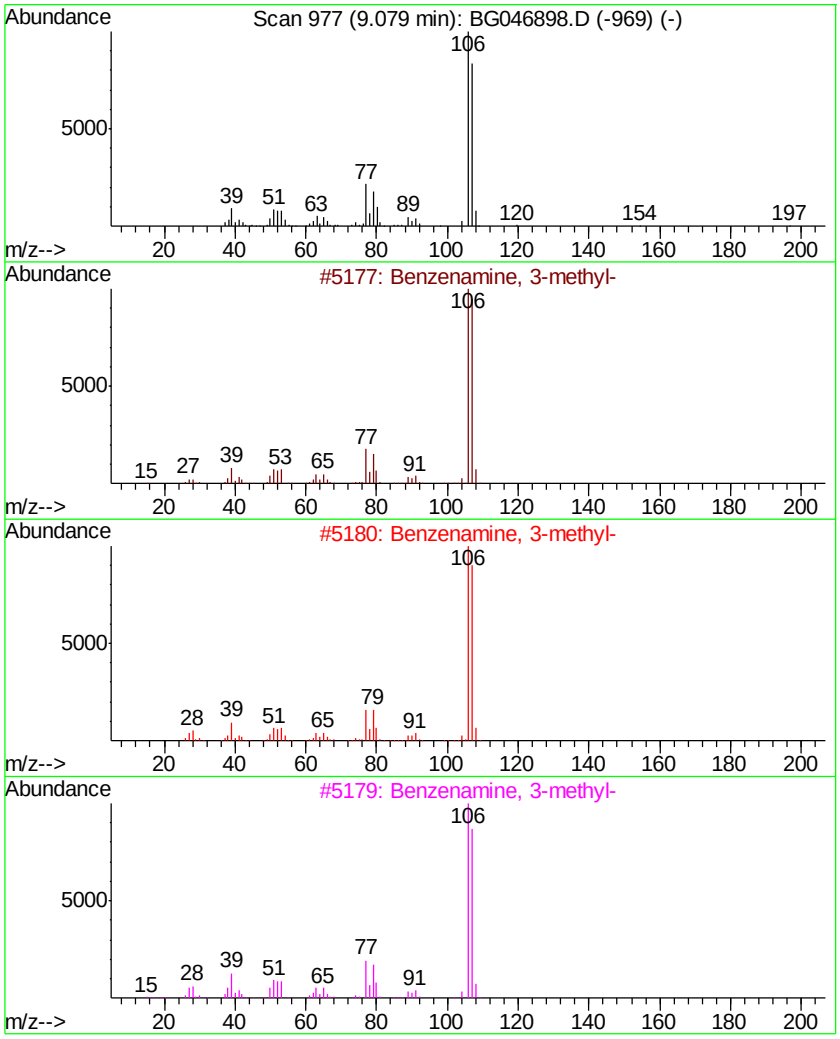
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

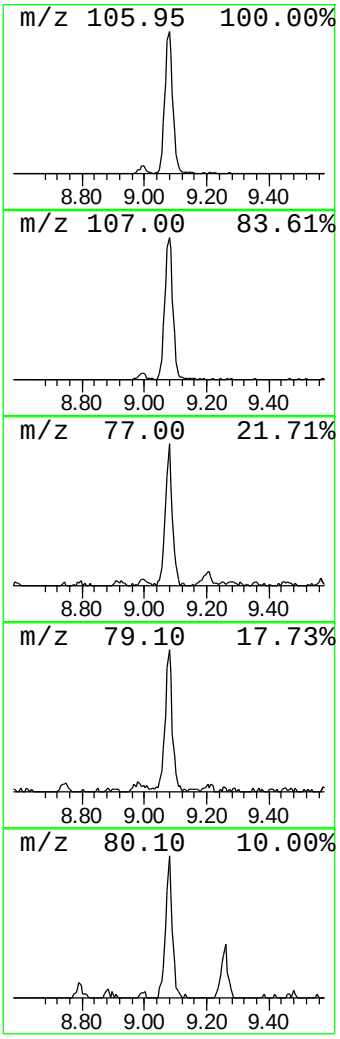
**********

Peak Number 4 Benzenamine, 3-methyl- Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.08 | 17.63 ng/ul | 384381 | 1,4-Dichlorobenzene-d4 | 8.10 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 95 |
| 2 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 94 |
| 3 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 94 |
| 4 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 94 |
| 5 | o-Toluidine | 107 | C7H9N | 000095-53-4 | 91 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

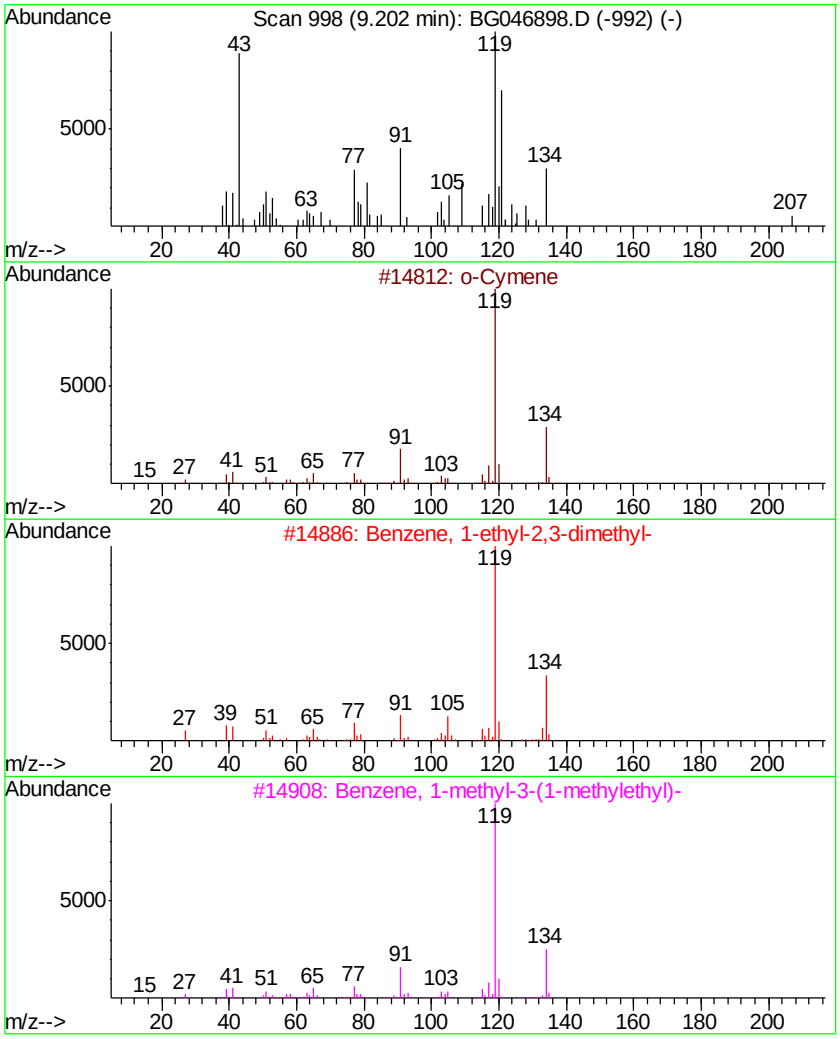
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

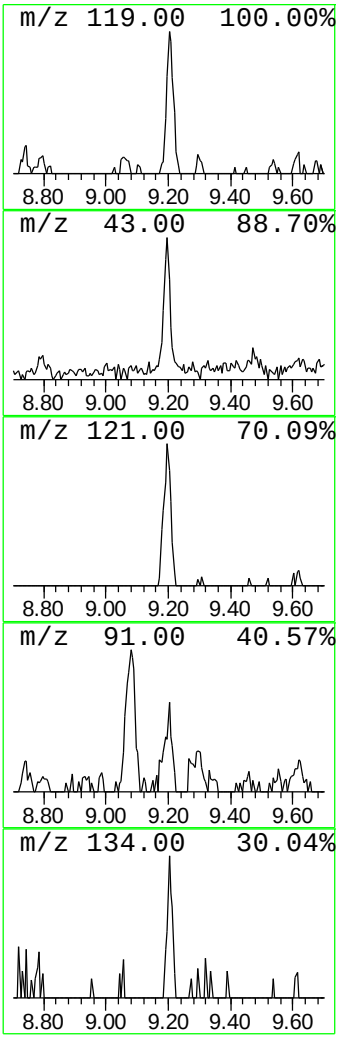
**********

Peak Number 5 unknown-01 Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.20 | 2.31 ng/ul | 50292 | 1,4-Dichlorobenzene-d4 | 8.10 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | o-Cymene | 134 | C10H14 | 000527-84-4 | 49 |
| 2 | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14 | 000933-98-2 | 47 |
| 3 | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14 | 000535-77-3 | 47 |
| 4 | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14 | 000095-93-2 | 47 |
| 5 | o-Cymene | 134 | C10H14 | 000527-84-4 | 47 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

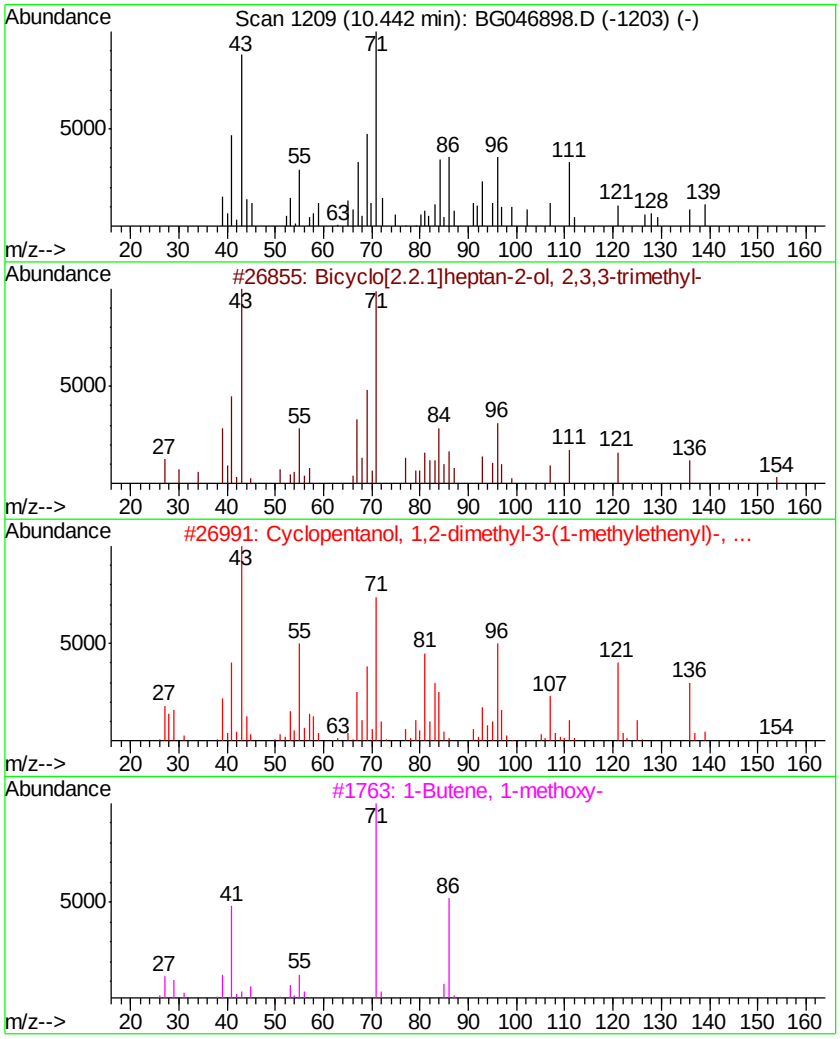
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

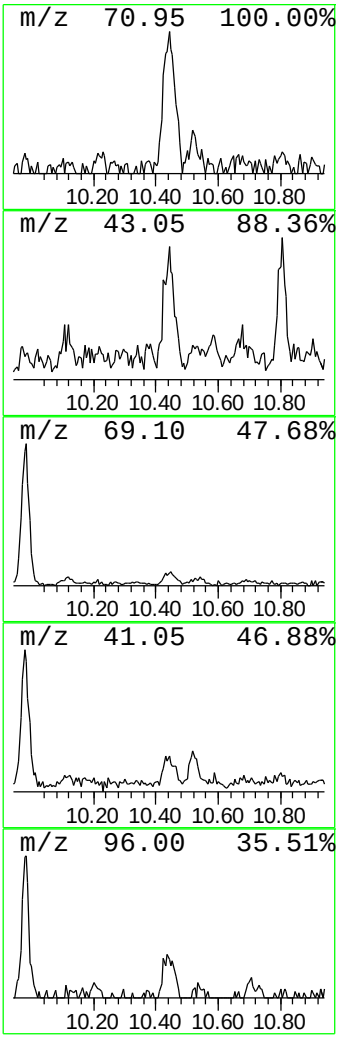
**********

Peak Number 6 unknown-02 Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.44 | 2.35 ng/ul | 67173 | Naphthalene-d8 | 10.91 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Bicyclo[2.2.1]heptan-2-ol, 2,3,3... | 154 | C10H18O | 000465-31-6 | 45 |
| 2 | Cyclopentanol, 1,2-dimethyl-3-(1... | 154 | C10H18O | 004028-59-5 | 38 |
| 3 | 1-Butene, 1-methoxy- | 86 | C5H10O | 029512-02-5 | 38 |
| 4 | 1-Propene, 3-methoxy-2-methyl- | 86 | C5H10O | 022418-49-1 | 35 |
| 5 | 2-Methoxycyclohexanone | 128 | C7H12O2 | 007429-44-9 | 35 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

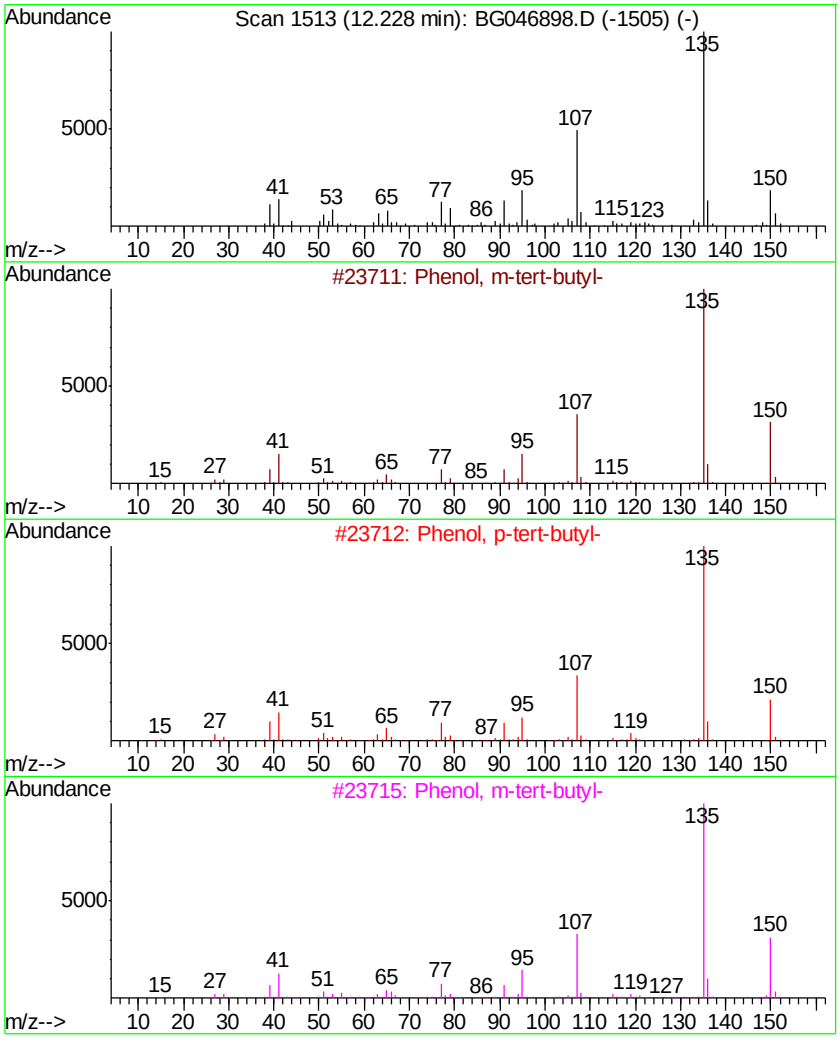
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

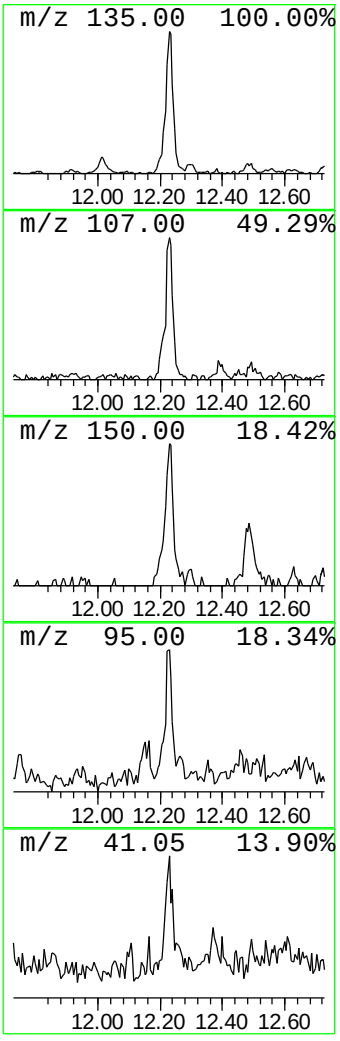
**********

Peak Number 7 Phenol, m-tert-butyl- Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.23 | 5.32 ng/ul | 152448 | Naphthalene-d8 | 10.91 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 90 |
| 2 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 87 |
| 3 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 83 |
| 4 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 78 |
| 5 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 74 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

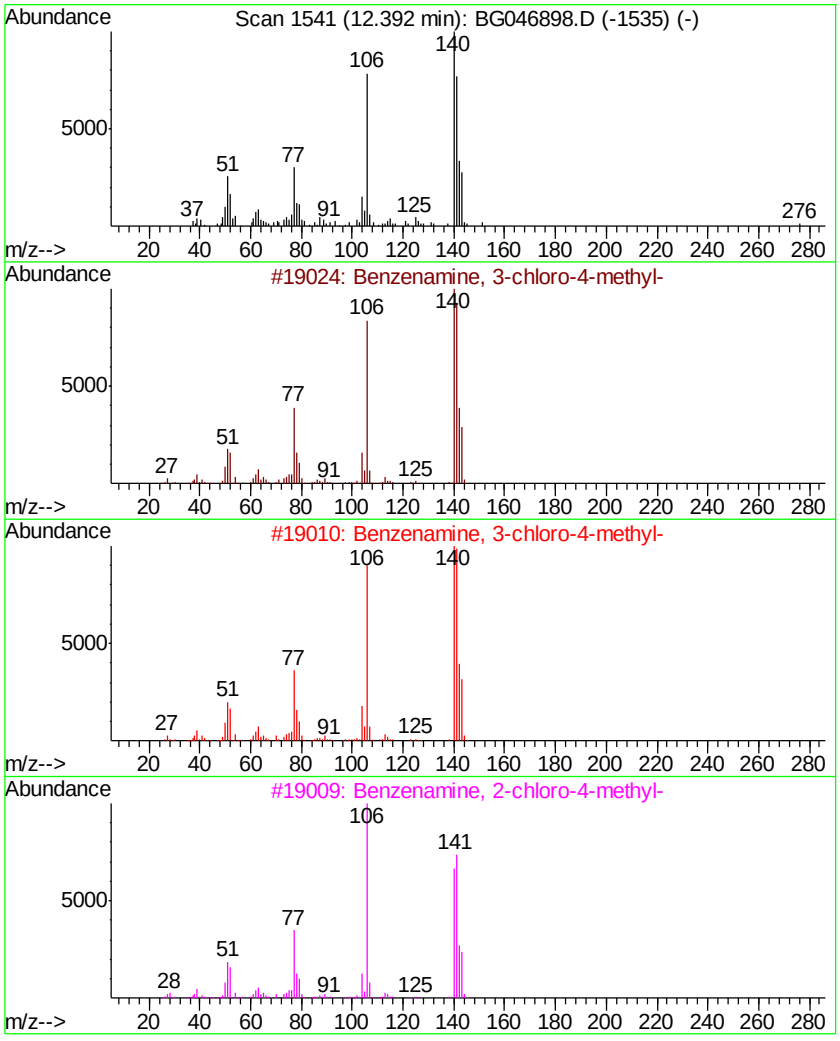
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

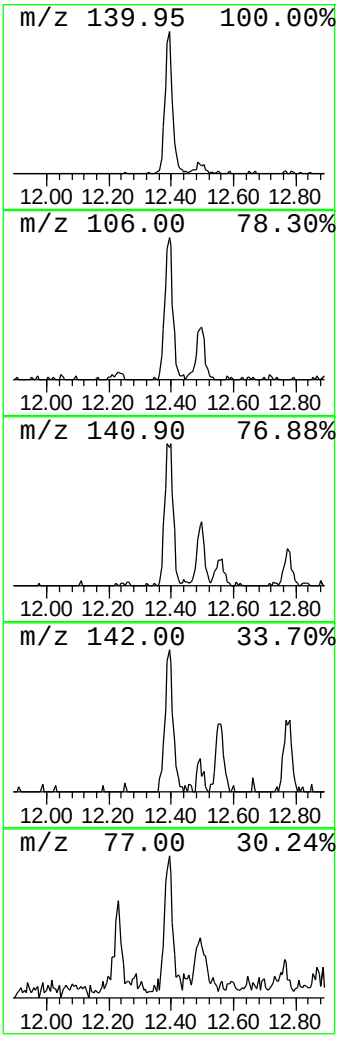
**********

Peak Number 8 Benzenamine, 3-chloro-4-met... Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.39 | 5.16 ng/ul | 147713 | Naphthalene-d8 | 10.91 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenamine, 3-chloro-4-methyl- | 141 | C7H8ClN | 000095-74-9 | 98 |
| 2 | Benzenamine, 3-chloro-4-methyl- | 141 | C7H8ClN | 000095-74-9 | 97 |
| 3 | Benzenamine, 2-chloro-4-methyl- | 141 | C7H8ClN | 000615-65-6 | 96 |
| 4 | Benzenamine, 2-chloro-4-methyl- | 141 | C7H8ClN | 000615-65-6 | 90 |
| 5 | Benzenamine, 5-chloro-2-methyl- | 141 | C7H8ClN | 000095-79-4 | 87 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

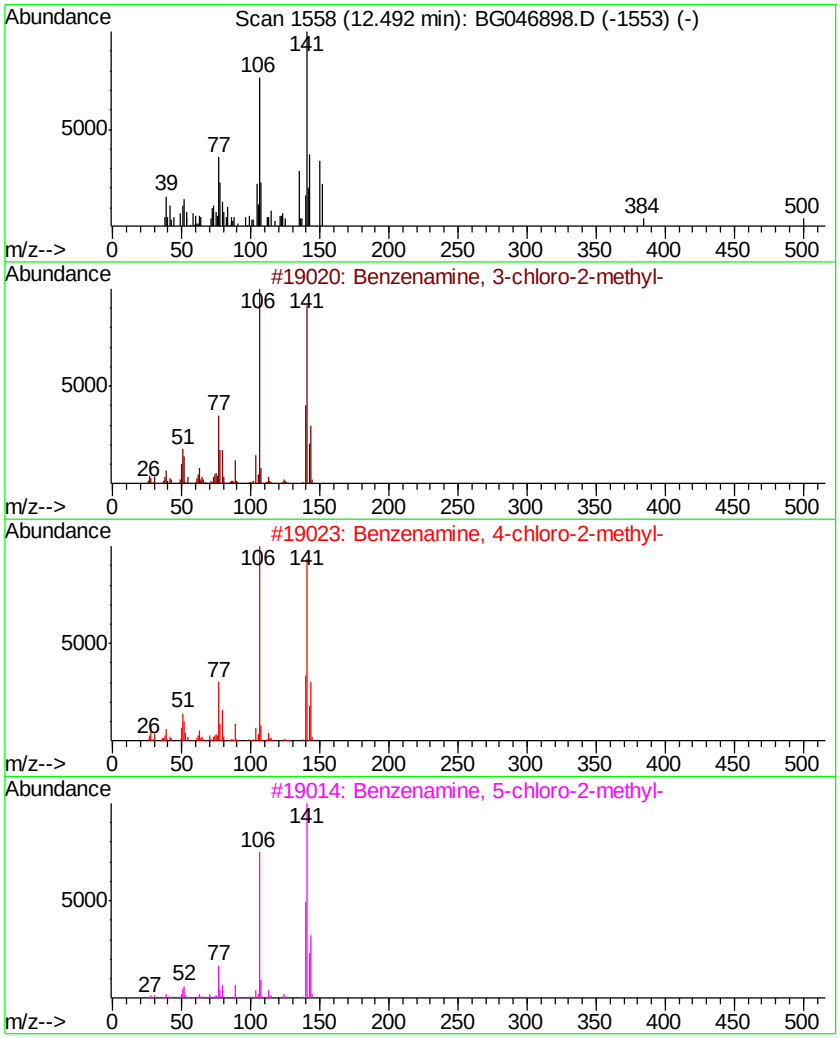
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

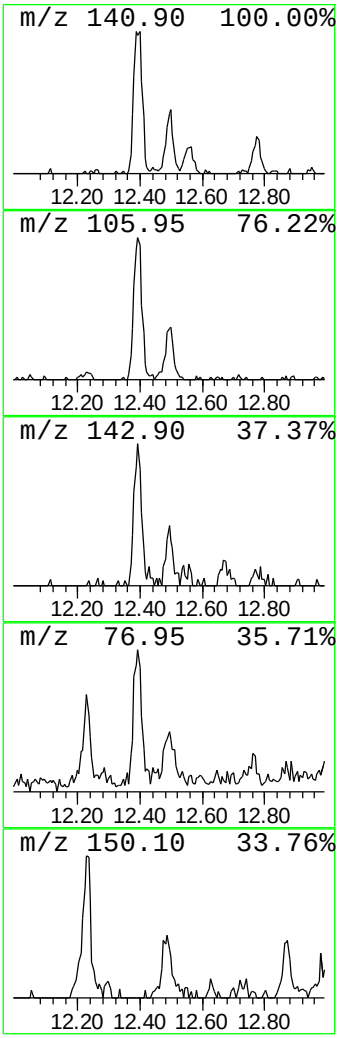
**********

Peak Number 9 Benzenamine, 3-chloro-2-met... Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.49 | 2.53 ng/ul | 72525 | Naphthalene-d8 | 10.91 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenamine, 3-chloro-2-methyl- | 141 | C7H8ClN | 000087-60-5 | 53 |
| 2 | Benzenamine, 4-chloro-2-methyl- | 141 | C7H8ClN | 000095-69-2 | 52 |
| 3 | Benzenamine, 5-chloro-2-methyl- | 141 | C7H8ClN | 000095-79-4 | 52 |
| 4 | Benzenamine, 5-chloro-2-methyl- | 141 | C7H8ClN | 000095-79-4 | 50 |
| 5 | Benzenamine, 4-chloro-2-methyl- | 141 | C7H8ClN | 000095-69-2 | 50 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

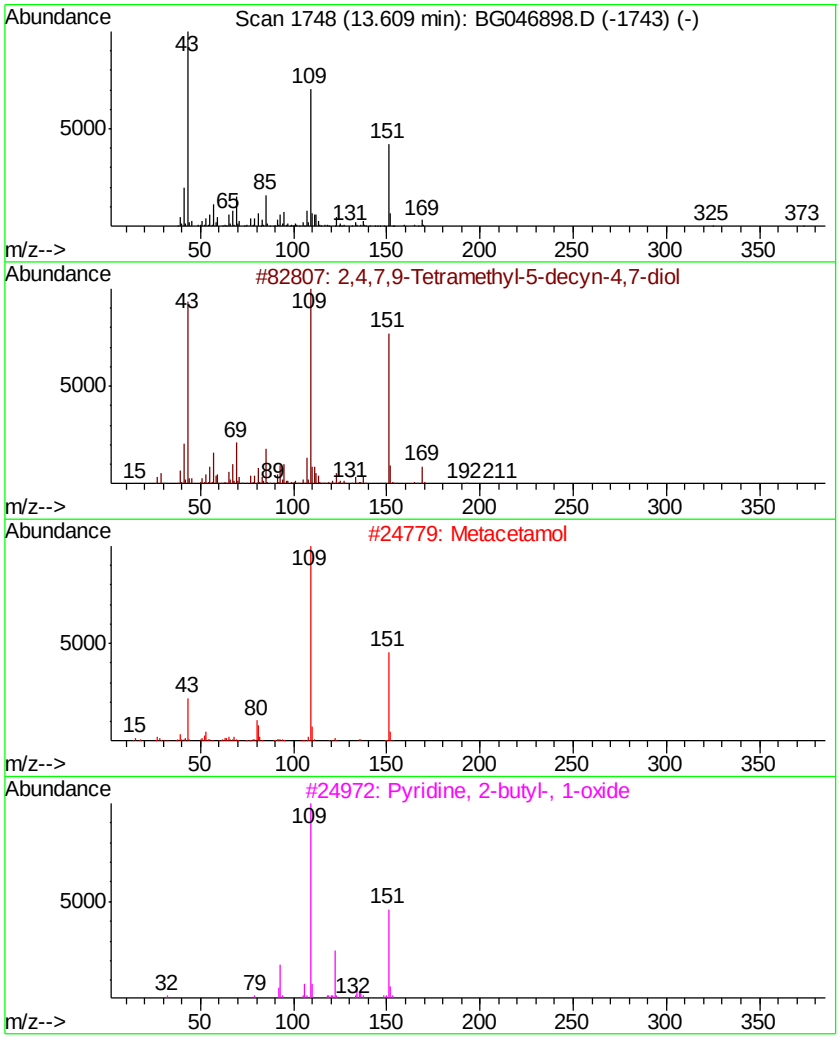
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

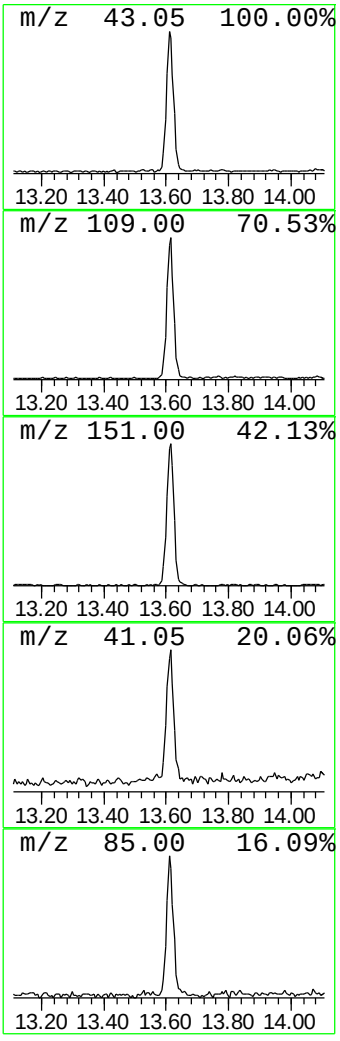
**********

Peak Number 10 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.61 | 13.02 ng/ul | 543815 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,4,7,9-Tetramethyl-5-decyn-4,7-... | 226 | C14H26O2 | 000126-86-3 | 90 |
| 2 | Metacetamol | 151 | C8H9NO2 | 000621-42-1 | 58 |
| 3 | Pyridine, 2-butyl-, 1-oxide | 151 | C9H13NO | 031396-32-4 | 53 |
| 4 | 2-Propenal, 3-(2,2,6-trimethyl-7... | 194 | C12H18O2 | 055759-91-6 | 50 |
| 5 | 2,4,7,9-Tetramethyl-5-decyn-4,7-... | 226 | C14H26O2 | 000126-86-3 | 45 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

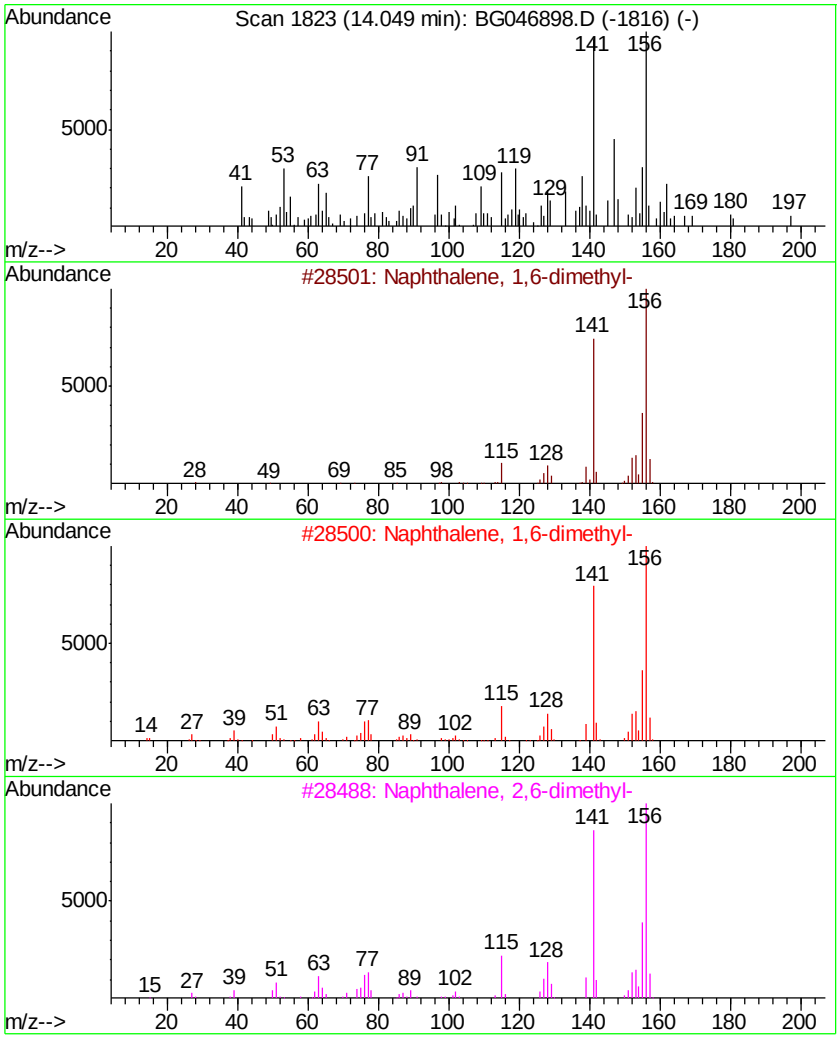
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

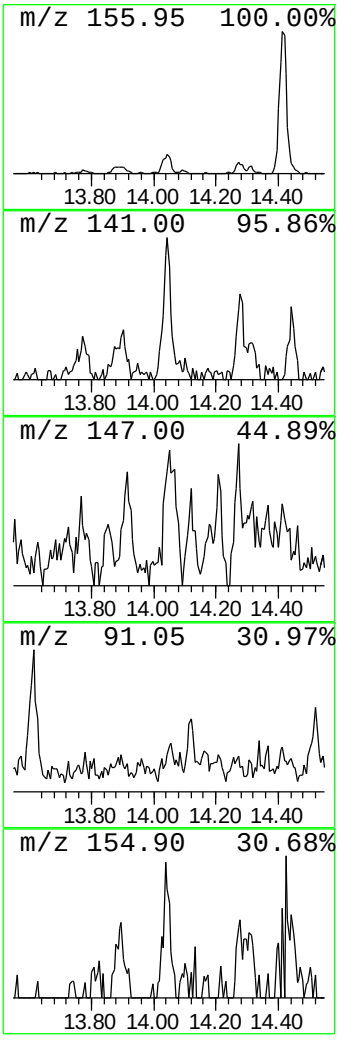
**********

Peak Number 11 Naphthalene, 1,6-dimethyl- Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.05 | 2.05 ng/ul | 85829 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 70 |
| 2 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 56 |
| 3 | Naphthalene, 2,6-dimethyl- | 156 | C12H12 | 000581-42-0 | 50 |
| 4 | Naphthalene, 1,5-dimethyl- | 156 | C12H12 | 000571-61-9 | 50 |
| 5 | Naphthalene, 1,4-dimethyl- | 156 | C12H12 | 000571-58-4 | 49 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

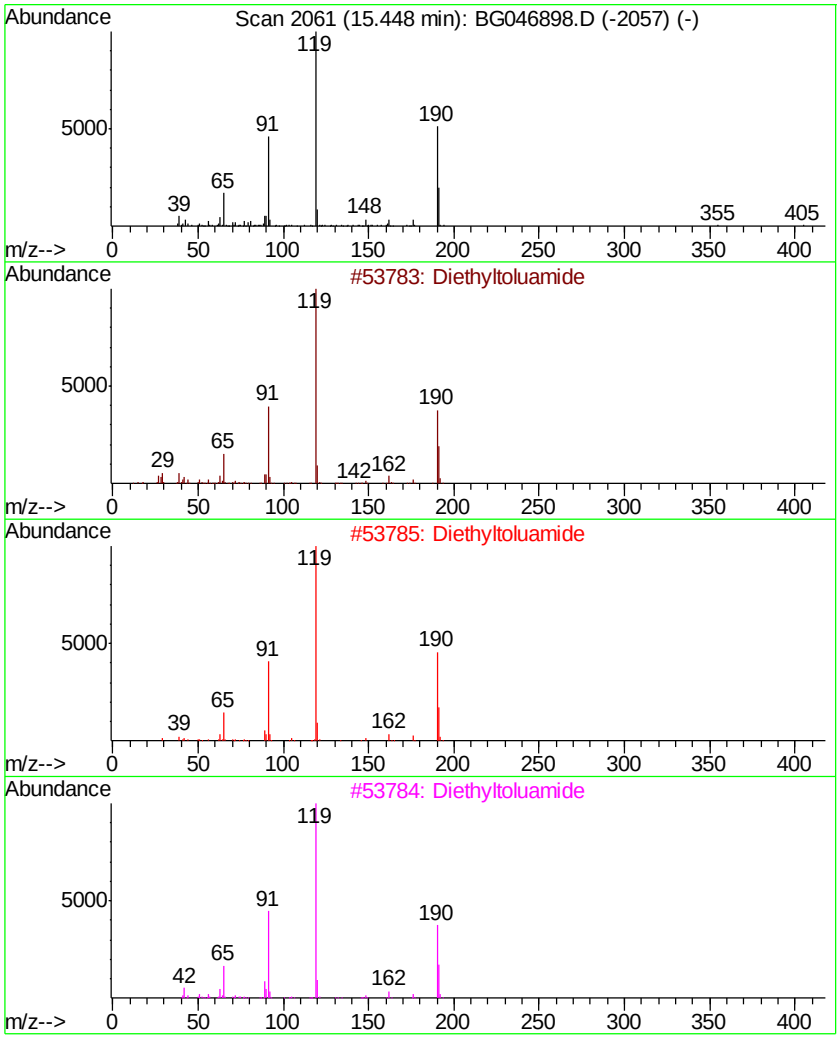
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

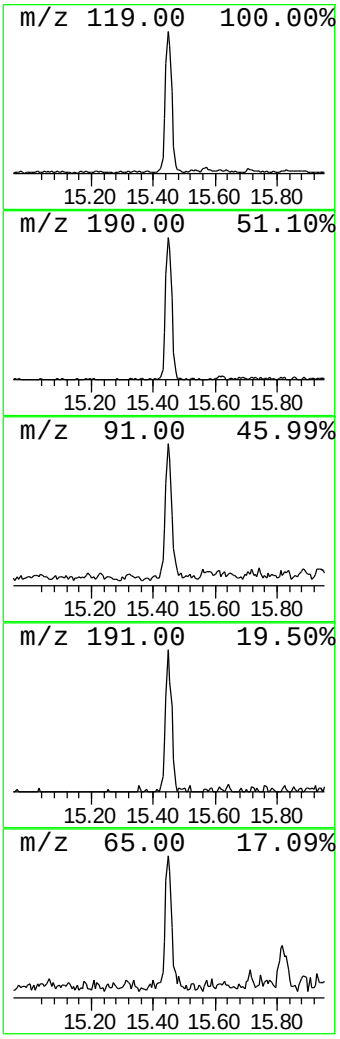
**********

Peak Number 12 Diethyltoluamide Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.45 | 4.61 ng/ul | 192485 | Acenaphthene-d10 | 14.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Diethyltoluamide | 191 | C12H17NO | 000134-62-3 | 94 |
| 2 | Diethyltoluamide | 191 | C12H17NO | 000134-62-3 | 94 |
| 3 | Diethyltoluamide | 191 | C12H17NO | 000134-62-3 | 91 |
| 4 | Benzamide, N,N-diethyl-4-methyl- | 191 | C12H17NO | 002728-05-4 | 87 |
| 5 | Diethyltoluamide | 191 | C12H17NO | 000134-62-3 | 64 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

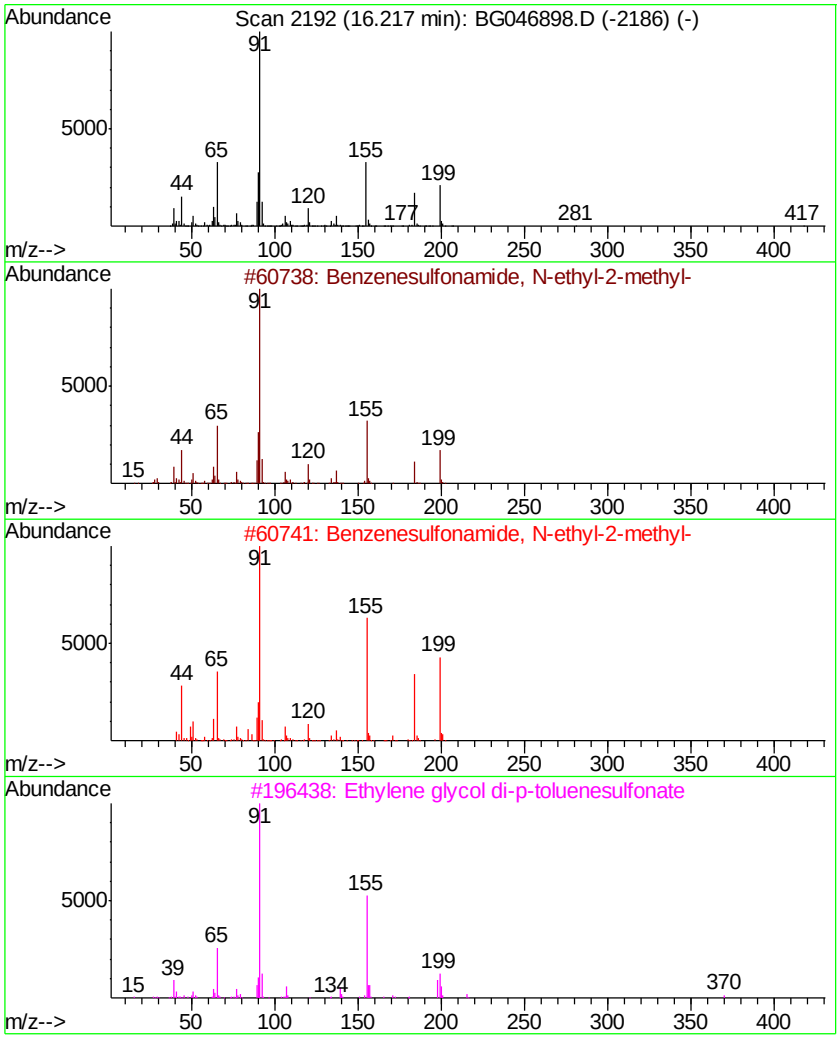
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

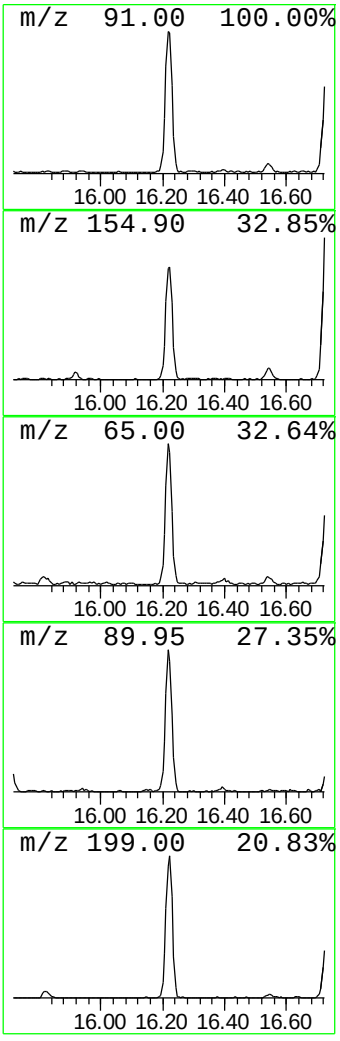
**********

Peak Number 13 Benzenesulfonamide, N-ethyl... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.22 | 15.71 ng/ul | 816820 | Phenanthrene-d10 | 17.46 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenesulfonamide, N-ethyl-2-me... | 199 | C9H13NO2S | 001077-56-1 | 94 |
| 2 | Benzenesulfonamide, N-ethyl-2-me... | 199 | C9H13NO2S | 001077-56-1 | 62 |
| 3 | Ethylene glycol di-p-toluenesulf... | 370 | C16H18O6S2 | 006315-52-2 | 47 |
| 4 | N3, N.alpha.-bis*ptoluenesulfony... | 462 | C20H22N4O5S2 | 1000130-21-9 | 47 |
| 5 | Benzenesulfonamide, N-ethyl-4-me... | 199 | C9H13NO2S | 000080-39-7 | 46 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

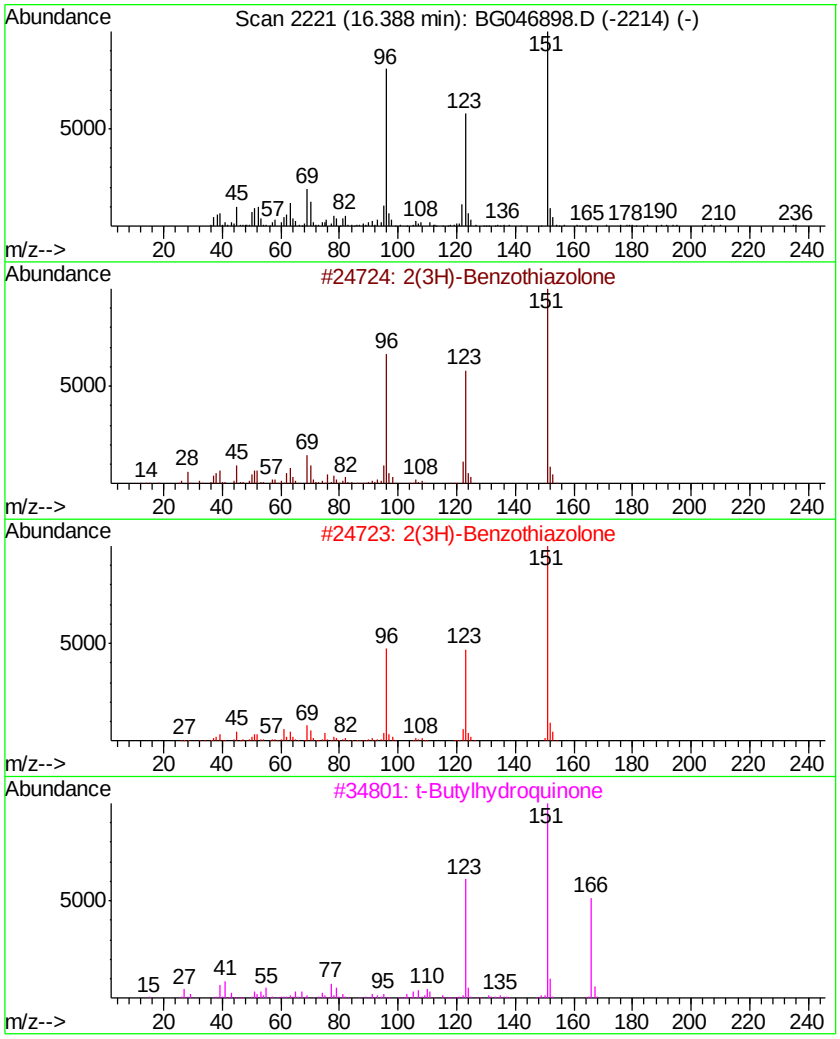
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

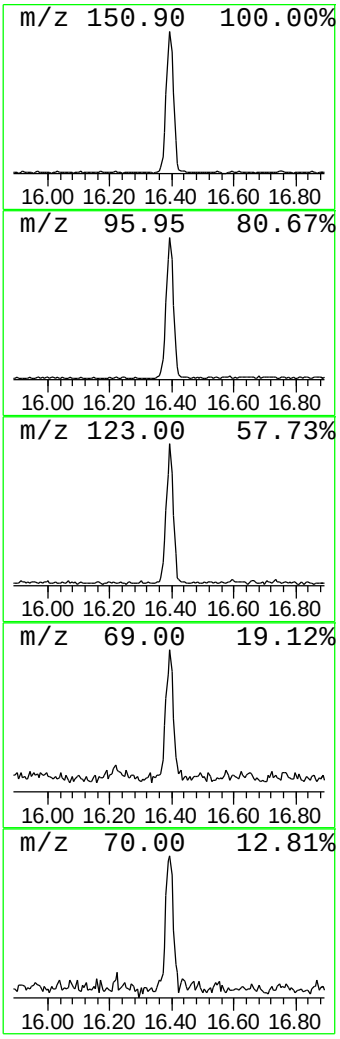
**********

Peak Number 14 2(3H)-Benzothiazolone Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.39 | 8.31 ng/ul | 431998 | Phenanthrene-d10 | 17.46 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2(3H)-Benzothiazolone | 151 | C7H5NOS | 000934-34-9 | 93 |
| 2 | 2(3H)-Benzothiazolone | 151 | C7H5NOS | 000934-34-9 | 91 |
| 3 | t-Butylhydroquinone | 166 | C10H14O2 | 001948-33-0 | 46 |
| 4 | 2,4-Dioxabicyclo[4.4.0]decan-5-o... | 212 | C12H20O3 | 124899-16-7 | 38 |
| 5 | Benzooxazole-2-thiol | 151 | C7H5NOS | 1000318-31-6 | 35 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

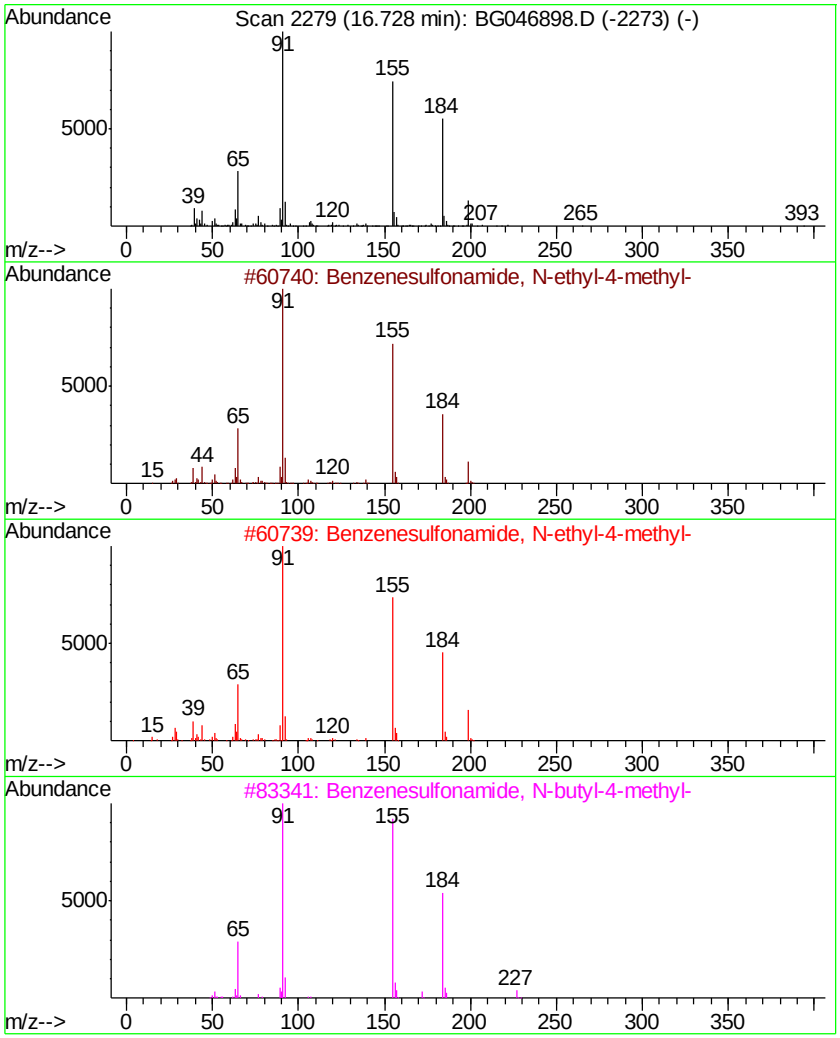
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

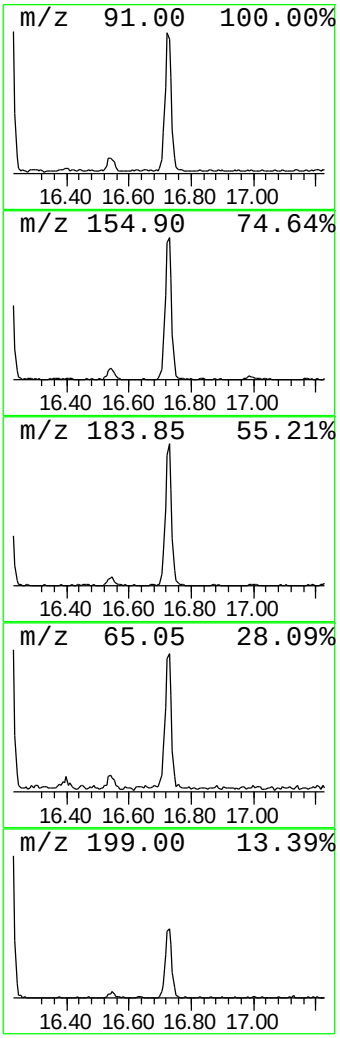
**********

Peak Number 15 Benzenesulfonamide, N-ethyl... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.73 | 9.46 ng/ul | 491965 | Phenanthrene-d10 | 17.46 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenesulfonamide, N-ethyl-4-me... | 199 | C9H13NO2S | 000080-39-7 | 96 |
| 2 | Benzenesulfonamide, N-ethyl-4-me... | 199 | C9H13NO2S | 000080-39-7 | 95 |
| 3 | Benzenesulfonamide, N-butyl-4-me... | 227 | C11H17NO2S | 001907-65-9 | 90 |
| 4 | Acetic acid, 2-(4-methylphenylsu... | 243 | C10H13NO4S | 002645-02-5 | 83 |
| 5 | N-Ethoxycarbonylmethyl-p-toluene... | 257 | C11H15NO4S | 005465-67-8 | 78 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

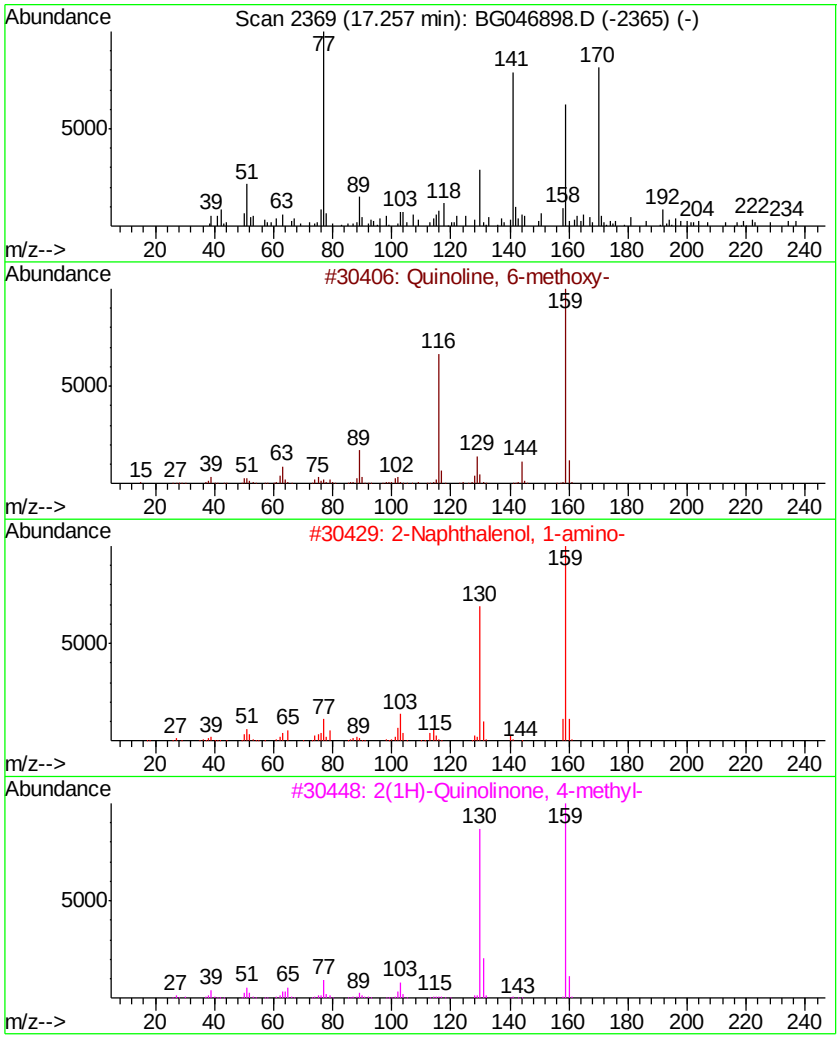
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

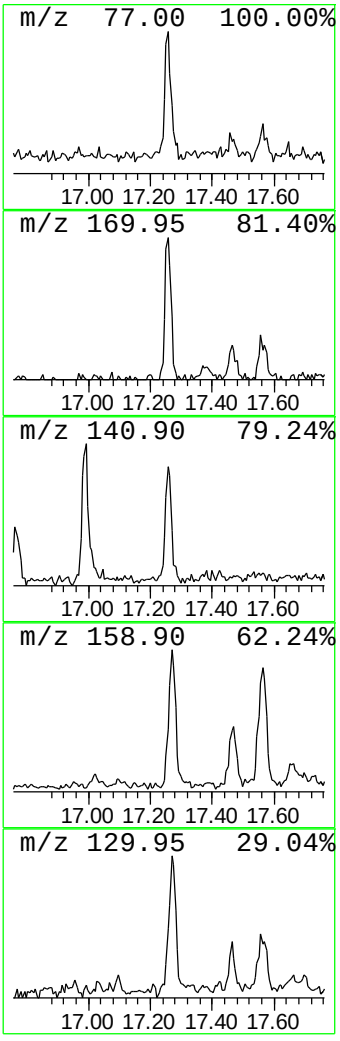
**********

Peak Number 16 unknown-03 Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.26 | 2.79 ng/ul | 145253 | Phenanthrene-d10 | 17.46 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Quinoline, 6-methoxy- | 159 | C10H9NO | 005263-87-6 | 46 |
| 2 | 2-Naphthalenol, 1-amino- | 159 | C10H9NO | 002834-92-6 | 46 |
| 3 | 2(1H)-Quinolinone, 4-methyl- | 159 | C10H9NO | 000607-66-9 | 43 |
| 4 | 1-Methyl-2-formylindole | 159 | C10H9NO | 027421-51-8 | 43 |
| 5 | 2(1H)-Quinolinone, 3-methyl- | 159 | C10H9NO | 002721-59-7 | 43 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

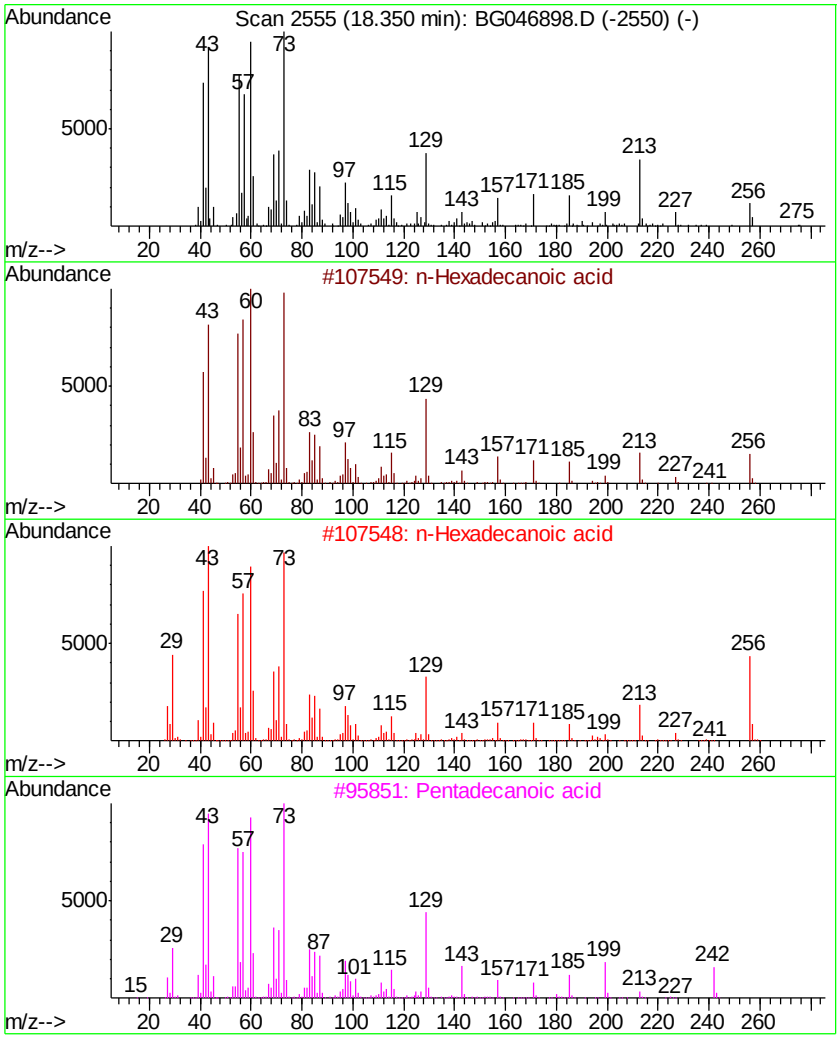
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

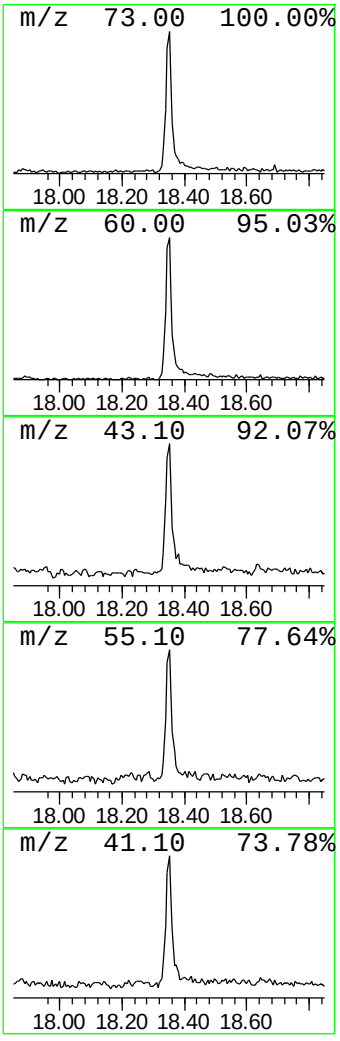
**********

Peak Number 17 n-Hexadecanoic acid Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.35 | 9.00 ng/ul | 467860 | Phenanthrene-d10 | 17.46 |

| 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 | 90 |
| 3 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 87 |
| 4 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 86 |
| 5 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 80 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

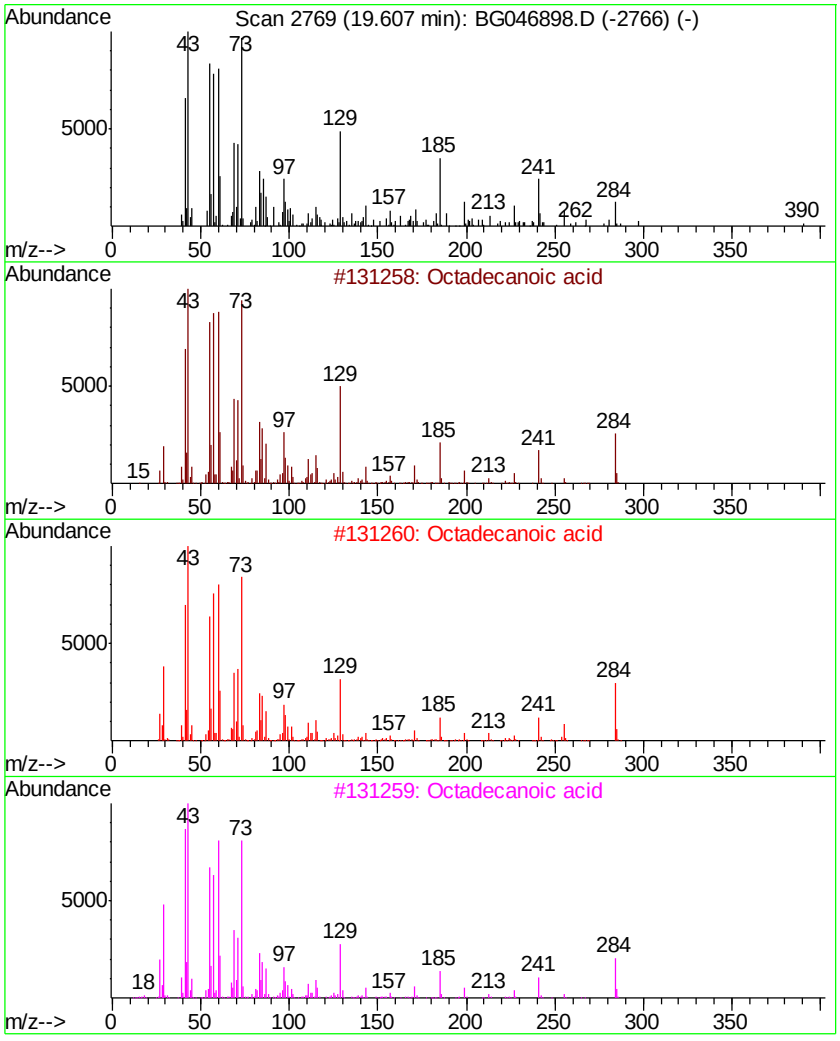
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

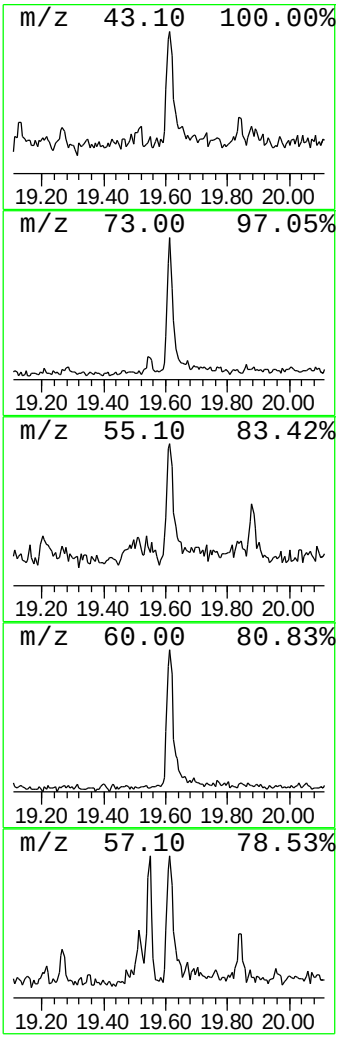
**********

Peak Number 18 Octadecanoic acid Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.61 | 3.47 ng/ul | 232371 | Chrysene-d12 | 21.74 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 98 |
| 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 | 83 |
| 5 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 78 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

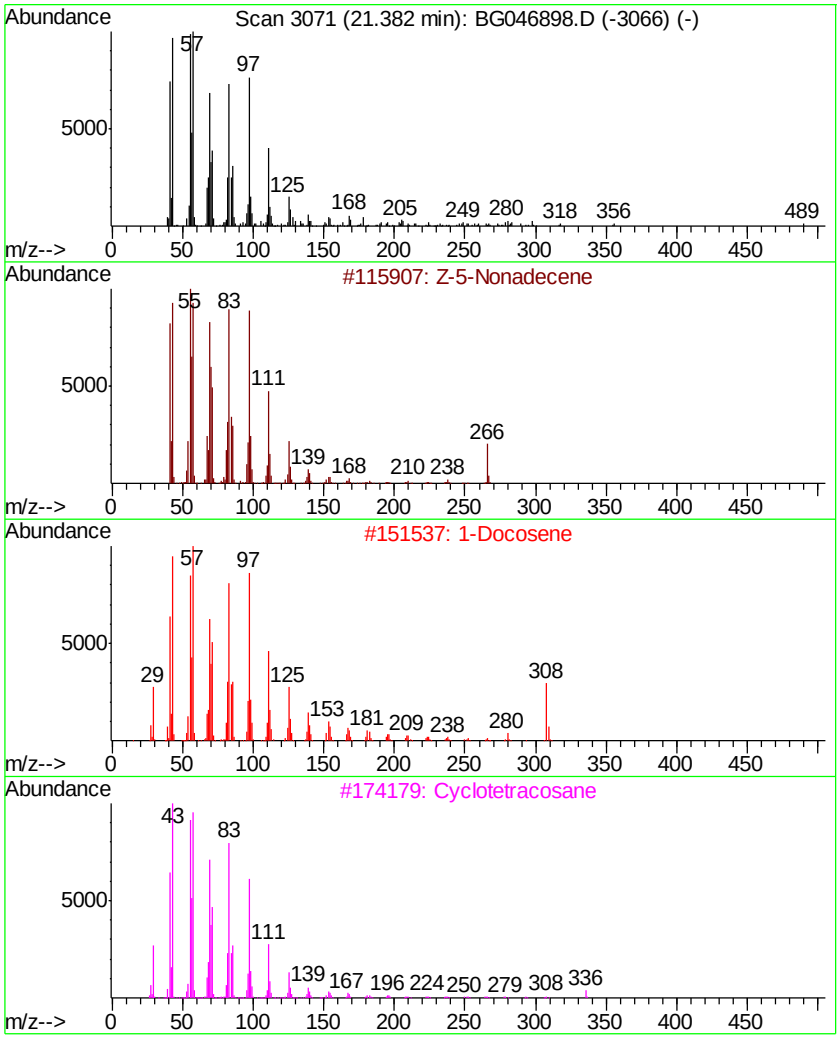
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

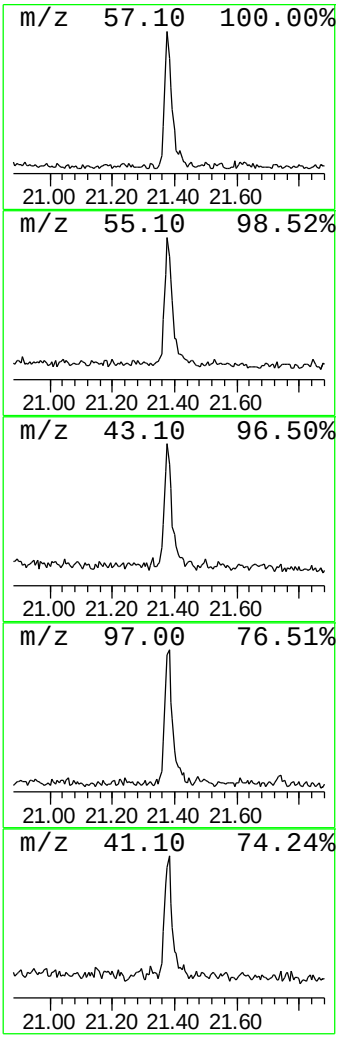
**********

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.38 | 5.59 ng/ul | 374740 | Chrysene-d12 | 21.74 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Z-5-Nonadecene | 266 | C19H38 | 1000131-11-8 | 95 |
| 2 | 1-Docosene | 308 | C22H44 | 001599-67-3 | 93 |
| 3 | Cyclotetracosane | 336 | C24H48 | 000297-03-0 | 91 |
| 4 | 1-Docosene | 308 | C22H44 | 001599-67-3 | 90 |
| 5 | 1-Nonadecene | 266 | C19H38 | 018435-45-5 | 89 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101420\
Data File : BG046898.D
Acq On : 14 Oct 2020 18:26
Operator : CG/JU
Sample : L4360-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** CB7W6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.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 |
| 1,3-Oxathiolane | 4.71 | 4.2 | ng/ul | 91632 | 1 | 8.10 | 436127 | 20.0 |
| Benzene, chloro- | 5.41 | 12.3 | ng/ul | 267631 | 1 | 8.10 | 436127 | 20.0 |
| Indane | 8.47 | 8.2 | ng/ul | 177980 | 1 | 8.10 | 436127 | 20.0 |
| Benzenamine, 3-me... | 9.08 | 17.6 | ng/ul | 384381 | 1 | 8.10 | 436127 | 20.0 |
| unknown-01 | 9.20 | 2.3 | ng/ul | 50292 | 1 | 8.10 | 436127 | 20.0 |
| unknown-02 | 10.44 | 2.4 | ng/ul | 67173 | 2 | 10.91 | 572861 | 20.0 |
| Phenol, m-tert-bu... | 12.23 | 5.3 | ng/ul | 152448 | 2 | 10.91 | 572861 | 20.0 |
| Benzenamine, 3-ch... | 12.39 | 5.2 | ng/ul | 147713 | 2 | 10.91 | 572861 | 20.0 |
| Benzenamine, 3-ch... | 12.49 | 2.5 | ng/ul | 72525 | 2 | 10.91 | 572861 | 20.0 |
| 2,4,7,9-Tetrameth... | 13.61 | 13.0 | ng/ul | 543815 | 3 | 14.72 | 835405 | 20.0 |
| Naphthalene, 1,6-... | 14.05 | 2.0 | ng/ul | 85829 | 3 | 14.72 | 835405 | 20.0 |
| Diethyltoluamide | 15.45 | 4.6 | ng/ul | 192485 | 3 | 14.72 | 835405 | 20.0 |
| Benzenesulfonamid... | 16.22 | 15.7 | ng/ul | 816820 | 4 | 17.46 | 1039890 | 20.0 |
| 2(3H)-Benzothiazo... | 16.39 | 8.3 | ng/ul | 431998 | 4 | 17.46 | 1039890 | 20.0 |
| Benzenesulfonamid... | 16.73 | 9.5 | ng/ul | 491965 | 4 | 17.46 | 1039890 | 20.0 |
| unknown-03 | 17.26 | 2.8 | ng/ul | 145253 | 4 | 17.46 | 1039890 | 20.0 |
| n-Hexadecanoic acid | 18.35 | 9.0 | ng/ul | 467860 | 4 | 17.46 | 1039890 | 20.0 |
| Octadecanoic acid | 19.61 | 3.5 | ng/ul | 232371 | 5 | 21.74 | 1340040 | 20.0 |
| (DEL) Alkane: Str... | 21.38 | 5.6 | ng/ul | 374740 | 5 | 21.74 | 1340040 | 20.0 |