LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Title : VOC Analysis

Signal : TIC: VV020046.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 1.110 | 3 | 6 | 18 | rVB2 | 36061 | 32438 | 0.40% | 0.122% |
| 2 | 1.216 | 34 | 39 | 45 | rBV4 | 9739 | 10908 | 0.14% | 0.041% |
| 3 | 1.309 | 60 | 68 | 81 | rVB | 224690 | 233128 | 2.89% | 0.875% |
| 4 | 1.383 | 86 | 91 | 97 | rBV | 11270 | 11932 | 0.15% | 0.045% |
| 5 | 1.566 | 141 | 148 | 162 | rVV | 194012 | 219992 | 2.73% | 0.826% |
| 6 | 1.634 | 164 | 169 | 181 | rVB2 | 9294 | 12376 | 0.15% | 0.046% |
| 7 | 1.804 | 217 | 222 | 231 | rVB4 | 9592 | 12716 | 0.16% | 0.048% |
| 8 | 2.110 | 303 | 317 | 331 | rBV | 403371 | 612681 | 7.60% | 2.299% |
| 9 | 2.290 | 369 | 373 | 382 | rVV3 | 3141 | 4954 | 0.06% | 0.019% |
| 10 | 2.441 | 414 | 420 | 422 | rBV4 | 1498 | 1276 | 0.02% | 0.005% |
| 11 | 2.531 | 428 | 448 | 471 | rVV | 481872 | 919027 | 11.41% | 3.449% |
| 12 | 2.762 | 503 | 520 | 546 | rBV | 853286 | 1683765 | 20.90% | 6.319% |
| 13 | 2.952 | 574 | 579 | 582 | rBV6 | 1764 | 1928 | 0.02% | 0.007% |
| 14 | 3.042 | 589 | 607 | 635 | rBV | 4131982 | 8057249 | 100.00% | 30.236% |
| 15 | 3.659 | 788 | 799 | 800 | rBV6 | 4947 | 5329 | 0.07% | 0.020% |
| 16 | 3.766 | 810 | 832 | 857 | rBV | 1384371 | 3443813 | 42.74% | 12.923% |
| 17 | 3.881 | 858 | 868 | 900 | rBV2 | 82610 | 228550 | 2.84% | 0.858% |
| 18 | 4.348 | 998 | 1013 | 1039 | rBV | 216668 | 535269 | 6.64% | 2.009% |
| 19 | 4.675 | 1096 | 1115 | 1131 | rBV6 | 31909 | 84553 | 1.05% | 0.317% |
| 20 | 4.766 | 1131 | 1143 | 1161 | rVB3 | 19341 | 46274 | 0.57% | 0.174% |
| 21 | 4.910 | 1171 | 1188 | 1207 | rBV4 | 27066 | 66221 | 0.82% | 0.249% |
| 22 | 5.045 | 1211 | 1230 | 1260 | rBV2 | 588021 | 1524888 | 18.93% | 5.722% |
| 23 | 5.235 | 1281 | 1289 | 1308 | rVB2 | 40171 | 95193 | 1.18% | 0.357% |
| 24 | 5.325 | 1308 | 1317 | 1325 | rBV10 | 6525 | 12697 | 0.16% | 0.048% |
| 25 | 5.492 | 1357 | 1369 | 1388 | rBV4 | 20775 | 50620 | 0.63% | 0.190% |
| 26 | 5.618 | 1396 | 1408 | 1442 | rBV | 304887 | 621596 | 7.71% | 2.333% |
| 27 | 5.907 | 1485 | 1498 | 1512 | rBV3 | 23168 | 49639 | 0.62% | 0.186% |
| 28 | 5.984 | 1518 | 1522 | 1524 | rBV4 | 1768 | 1460 | 0.02% | 0.005% |
| 29 | 6.071 | 1535 | 1549 | 1578 | rBV | 334785 | 727613 | 9.03% | 2.730% |
| 30 | 6.203 | 1581 | 1590 | 1598 | rBV4 | 14427 | 26784 | 0.33% | 0.101% |
| 31 | 6.460 | 1658 | 1670 | 1685 | rVB9 | 11755 | 27447 | 0.34% | 0.103% |
| 32 | 6.656 | 1710 | 1731 | 1749 | rVB2 | 44834 | 100983 | 1.25% | 0.379% |
| 33 | 6.778 | 1753 | 1769 | 1780 | rBV2 | 60632 | 126265 | 1.57% | 0.474% |
| 34 | 6.839 | 1781 | 1788 | 1799 | rVB2 | 16269 | 26219 | 0.33% | 0.098% |
| 35 | 6.920 | 1805 | 1813 | 1820 | rBV3 | 26979 | 44032 | 0.55% | 0.165% |
| 36 | 6.987 | 1825 | 1834 | 1855 | rVB | 189270 | 338835 | 4.21% | 1.272% |

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Title : VOC Analysis

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 7.084 | 1856 | 1864 | 1872 | rBV2 | 33779 | 53437 | 0.66% | 0.201% |
| 38 | 7.183 | 1886 | 1895 | 1905 | rVB4 | 6059 | 11591 | 0.14% | 0.043% |
| 39 | 7.267 | 1907 | 1921 | 1927 | rBV4 | 45245 | 88125 | 1.09% | 0.331% |
| 40 | 7.318 | 1927 | 1937 | 1965 | rVB | 750251 | 1283269 | 15.93% | 4.816% |
| 41 | 7.441 | 1968 | 1975 | 1982 | rBV5 | 11236 | 17351 | 0.22% | 0.065% |
| 42 | 7.569 | 2008 | 2015 | 2023 | rBV2 | 24368 | 34065 | 0.42% | 0.128% |
| 43 | 7.624 | 2023 | 2032 | 2055 | rVB | 112881 | 210837 | 2.62% | 0.791% |
| 44 | 7.788 | 2074 | 2083 | 2098 | rVB3 | 20644 | 37512 | 0.47% | 0.141% |
| 45 | 7.878 | 2108 | 2111 | 2114 | rBV4 | 2864 | 2715 | 0.03% | 0.010% |
| 46 | 7.981 | 2138 | 2143 | 2148 | rBV7 | 3536 | 3772 | 0.05% | 0.014% |
| 47 | 8.013 | 2148 | 2153 | 2165 | rVB2 | 7677 | 11882 | 0.15% | 0.045% |
| 48 | 8.090 | 2167 | 2177 | 2206 | rBV2 | 326467 | 663138 | 8.23% | 2.489% |
| 49 | 8.357 | 2253 | 2260 | 2268 | rBV3 | 9557 | 13752 | 0.17% | 0.052% |
| 50 | 8.515 | 2300 | 2309 | 2315 | rVB8 | 3381 | 5532 | 0.07% | 0.021% |
| 51 | 8.576 | 2315 | 2328 | 2331 | rBV3 | 8607 | 14532 | 0.18% | 0.055% |
| 52 | 8.669 | 2350 | 2357 | 2369 | rVB3 | 33284 | 59015 | 0.73% | 0.221% |
| 53 | 8.794 | 2383 | 2396 | 2404 | rBV | 28452 | 50285 | 0.62% | 0.189% |
| 54 | 8.855 | 2405 | 2415 | 2435 | rVV | 450008 | 737395 | 9.15% | 2.767% |
| 55 | 9.209 | 2517 | 2525 | 2548 | rVB5 | 27085 | 60654 | 0.75% | 0.228% |
| 56 | 9.322 | 2552 | 2560 | 2572 | rVB9 | 4500 | 7573 | 0.09% | 0.028% |
| 57 | 9.479 | 2600 | 2609 | 2620 | rBV9 | 12187 | 23495 | 0.29% | 0.088% |
| 58 | 9.585 | 2638 | 2642 | 2646 | rVV6 | 3888 | 4093 | 0.05% | 0.015% |
| 59 | 9.682 | 2666 | 2672 | 2673 | rBV3 | 4248 | 3553 | 0.04% | 0.013% |
| 60 | 9.807 | 2703 | 2711 | 2724 | rBV9 | 12612 | 26047 | 0.32% | 0.098% |
| 61 | 10.222 | 2827 | 2840 | 2858 | rVB | 416390 | 656072 | 8.14% | 2.462% |
| 62 | 10.521 | 2925 | 2933 | 2942 | rBV10 | 4699 | 8525 | 0.11% | 0.032% |
| 63 | 10.585 | 2947 | 2953 | 2968 | rVB8 | 10118 | 20295 | 0.25% | 0.076% |
| 64 | 10.736 | 2992 | 3000 | 3006 | rBV7 | 5813 | 9406 | 0.12% | 0.035% |
| 65 | 11.061 | 3088 | 3101 | 3106 | rBV5 | 10136 | 18581 | 0.23% | 0.070% |
| 66 | 11.154 | 3124 | 3130 | 3138 | rBV10 | 5300 | 9538 | 0.12% | 0.036% |
| 67 | 11.196 | 3140 | 3143 | 3150 | rVV6 | 4869 | 6183 | 0.08% | 0.023% |
| 68 | 11.254 | 3150 | 3161 | 3183 | rVV | 550880 | 843878 | 10.47% | 3.167% |
| 69 | 11.556 | 3244 | 3255 | 3257 | rBV2 | 19150 | 25244 | 0.31% | 0.095% |
| 70 | 11.630 | 3268 | 3278 | 3300 | rVB | 808011 | 1321597 | 16.40% | 4.960% |
| 71 | 11.752 | 3309 | 3316 | 3324 | rBV6 | 5597 | 7732 | 0.10% | 0.029% |
| 72 | 11.830 | 3332 | 3340 | 3351 | rVB | 19759 | 30580 | 0.38% | 0.115% |
| 73 | 11.923 | 3360 | 3369 | 3373 | rBV2 | 7492 | 12032 | 0.15% | 0.045% |
| 74 | 11.997 | 3386 | 3392 | 3394 | rBV5 | 1518 | 1644 | 0.02% | 0.006% |
| 75 | 12.029 | 3395 | 3402 | 3407 | rBV3 | 6244 | 9132 | 0.11% | 0.034% |
| 76 | 12.132 | 3422 | 3434 | 3439 | rBV5 | 10200 | 18514 | 0.23% | 0.069% |

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Title : VOC Analysis

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 12.254 | 3462 | 3472 | 3483 | rBV5 | 8689 | 22011 | 0.27% | 0.083% |
| 78 | 12.363 | 3499 | 3506 | 3507 | rBV7 | 2776 | 3163 | 0.04% | 0.012% |
| 79 | 12.421 | 3517 | 3524 | 3535 | rVB2 | 29357 | 46152 | 0.57% | 0.173% |
| 80 | 12.476 | 3535 | 3541 | 3552 | rVB2 | 3547 | 5862 | 0.07% | 0.022% |
| 81 | 12.559 | 3558 | 3567 | 3576 | rBV2 | 16523 | 24658 | 0.31% | 0.093% |
| 82 | 12.694 | 3592 | 3609 | 3617 | rBV2 | 9642 | 25021 | 0.31% | 0.094% |
| 83 | 12.736 | 3617 | 3622 | 3636 | rVB8 | 7492 | 12362 | 0.15% | 0.046% |
| 84 | 12.807 | 3637 | 3644 | 3649 | rBV7 | 7125 | 9426 | 0.12% | 0.035% |
| 85 | 12.839 | 3651 | 3654 | 3656 | rBV4 | 1930 | 1494 | 0.02% | 0.006% |
| 86 | 12.958 | 3686 | 3691 | 3701 | rVB7 | 7080 | 10718 | 0.13% | 0.040% |
| 87 | 13.013 | 3701 | 3708 | 3715 | rBV4 | 4635 | 7049 | 0.09% | 0.026% |
| 88 | 13.174 | 3750 | 3758 | 3766 | rBV6 | 5278 | 8385 | 0.10% | 0.031% |
| 89 | 13.276 | 3783 | 3790 | 3797 | rBV2 | 16989 | 23633 | 0.29% | 0.089% |
| 90 | 13.379 | 3817 | 3822 | 3826 | rVB8 | 1955 | 2061 | 0.03% | 0.008% |
| 91 | 13.624 | 3891 | 3898 | 3903 | rVB9 | 1748 | 2451 | 0.03% | 0.009% |
| 92 | 13.720 | 3923 | 3928 | 3932 | rVV5 | 1693 | 1489 | 0.02% | 0.006% |
| 93 | 14.103 | 4041 | 4047 | 4049 | rBV5 | 1548 | 1622 | 0.02% | 0.006% |
| 94 | 14.241 | 4083 | 4090 | 4100 | rBV9 | 4253 | 6067 | 0.08% | 0.023% |
| 95 | 14.315 | 4107 | 4113 | 4117 | rBV9 | 1771 | 1622 | 0.02% | 0.006% |
| 96 | 14.382 | 4131 | 4134 | 4141 | rVB5 | 2319 | 2303 | 0.03% | 0.009% |
| 97 | 14.646 | 4210 | 4216 | 4221 | rBV6 | 1396 | 1643 | 0.02% | 0.006% |
| 98 | 14.717 | 4235 | 4238 | 4241 | rBV6 | 1762 | 1441 | 0.02% | 0.005% |
| 99 | 15.678 | 4531 | 4537 | 4544 | rVV10 | 2178 | 3321 | 0.04% | 0.012% |
| 100 | 15.817 | 4573 | 4580 | 4583 | rBV6 | 2350 | 2523 | 0.03% | 0.009% |

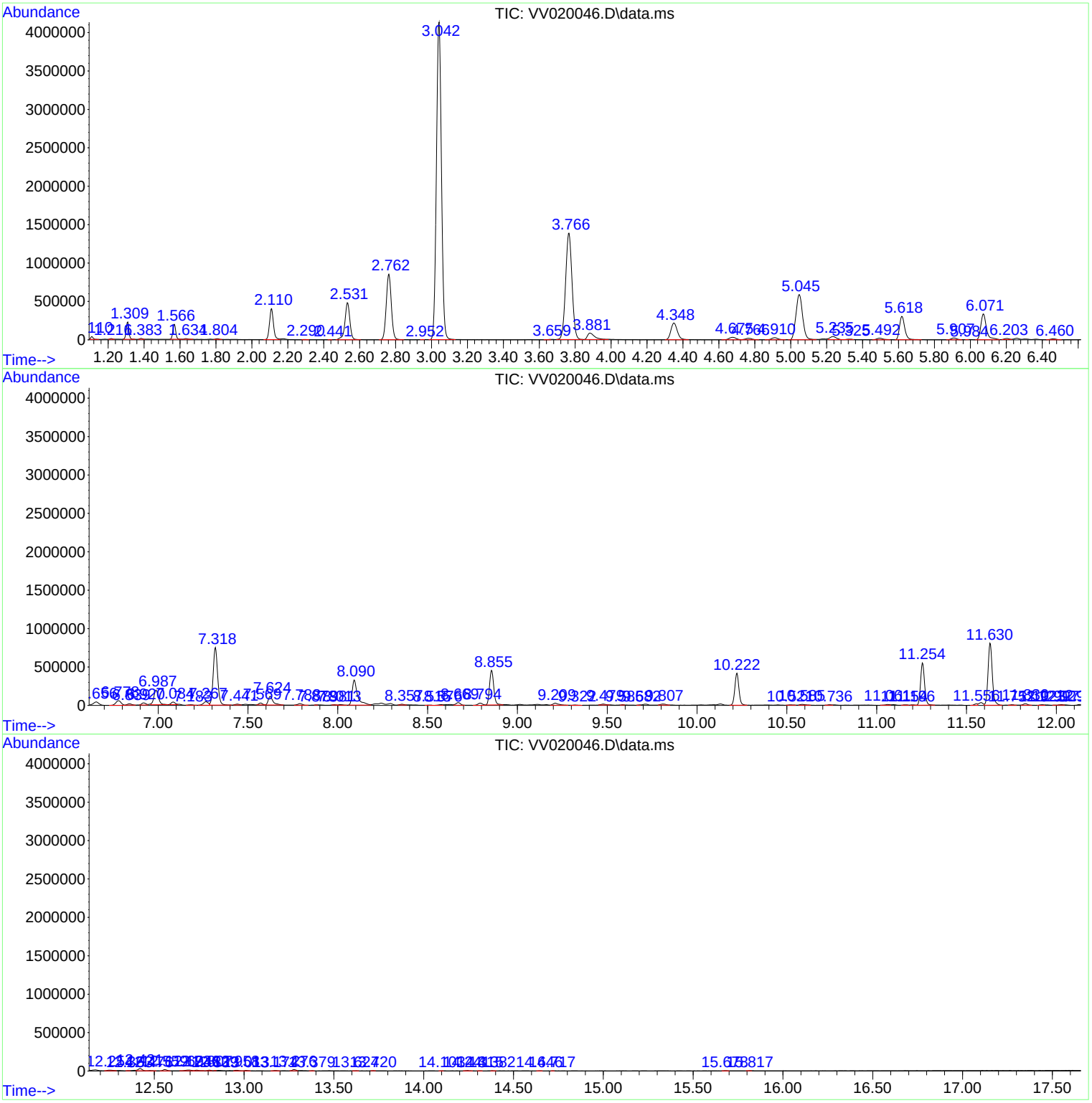
Sum of corrected areas: 26647700

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :**
MSVOA_V
**ClientSampleId :**
BG4Y7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

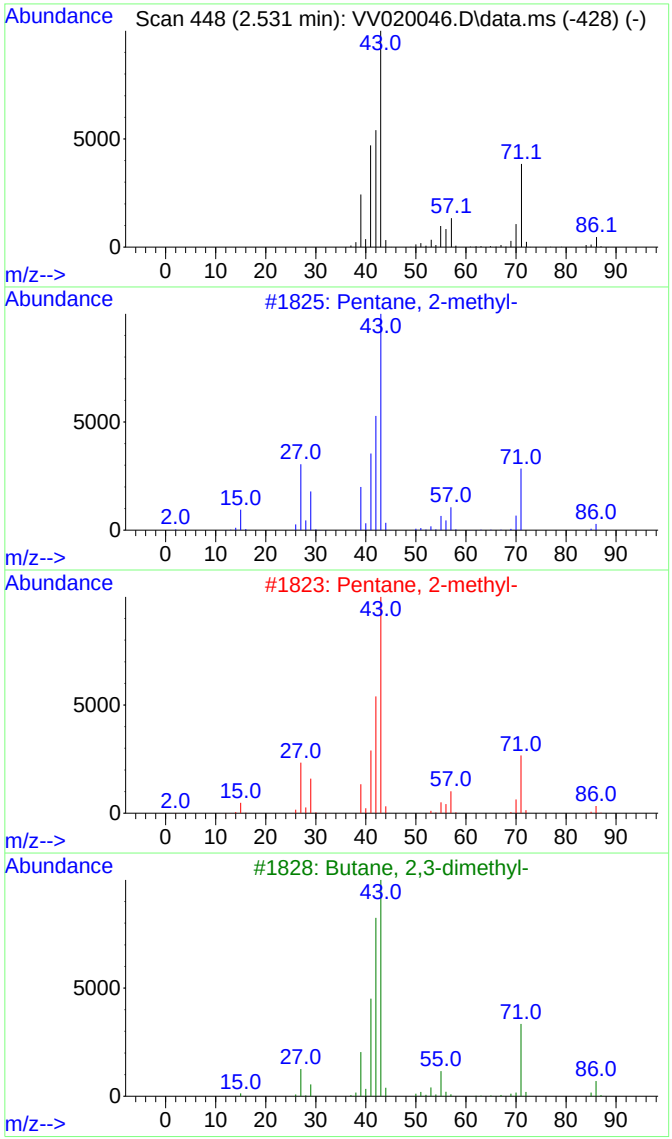
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

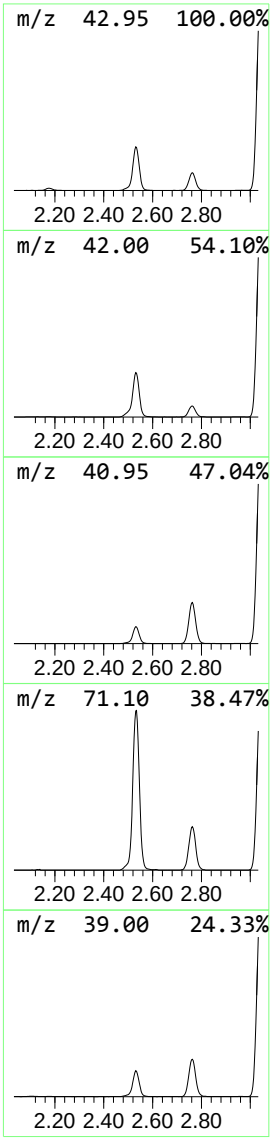
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.531 | 73.92 ug/L | 919027 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentane, 2-methyl- | 86 | C6H14 | 000107-83-5 | 91 |
| 2 | Pentane, 2-methyl- | 86 | C6H14 | 000107-83-5 | 90 |
| 3 | Butane, 2,3-dimethyl- | 86 | C6H14 | 000079-29-8 | 64 |
| 4 | Butane, 2,3-dimethyl- | 86 | C6H14 | 000079-29-8 | 47 |
| 5 | 1-Butanol, 2,3-dimethyl- | 102 | C6H14O | 019550-30-2 | 45 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

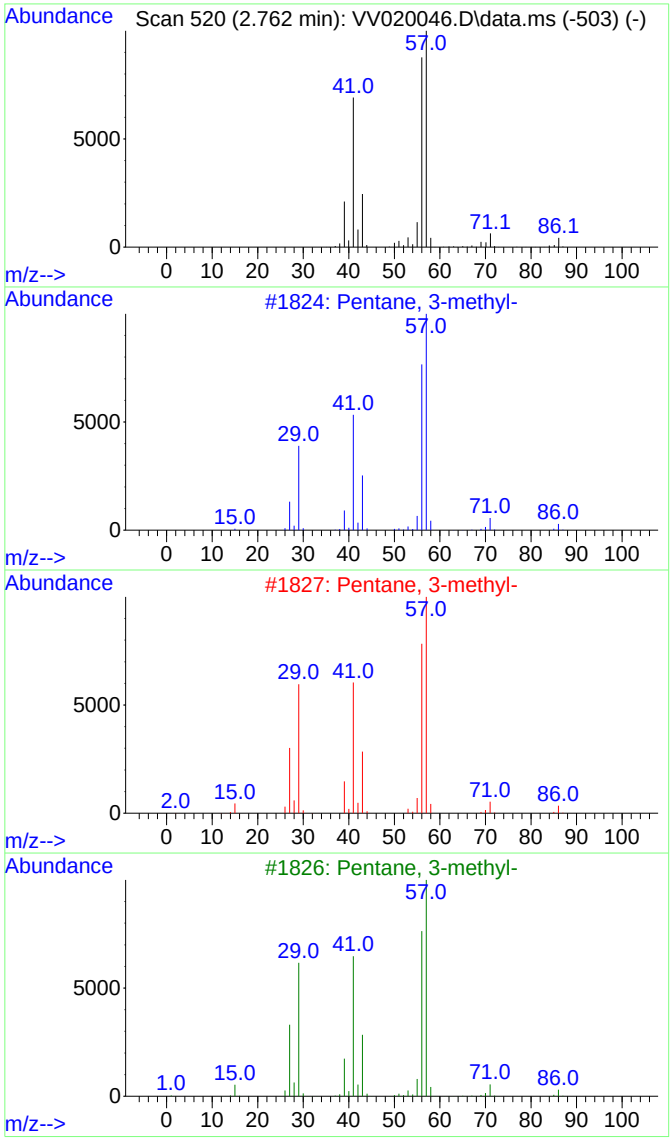
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

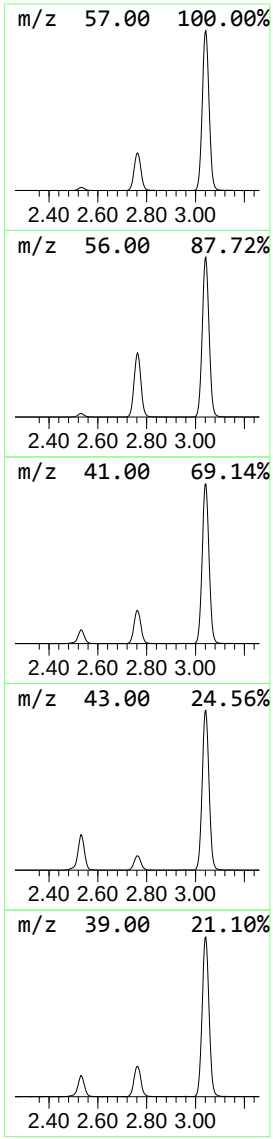
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.762 | 135.44 ug/L | 1683770 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentane, 3-methyl- | 86 | C6H14 | 000096-14-0 | 90 |
| 2 | Pentane, 3-methyl- | 86 | C6H14 | 000096-14-0 | 90 |
| 3 | Pentane, 3-methyl- | 86 | C6H14 | 000096-14-0 | 90 |
| 4 | Hexane, 2,2,3-trimethyl- | 128 | C9H20 | 016747-25-4 | 64 |
| 5 | Hexane, 2,2,3-trimethyl- | 128 | C9H20 | 016747-25-4 | 50 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

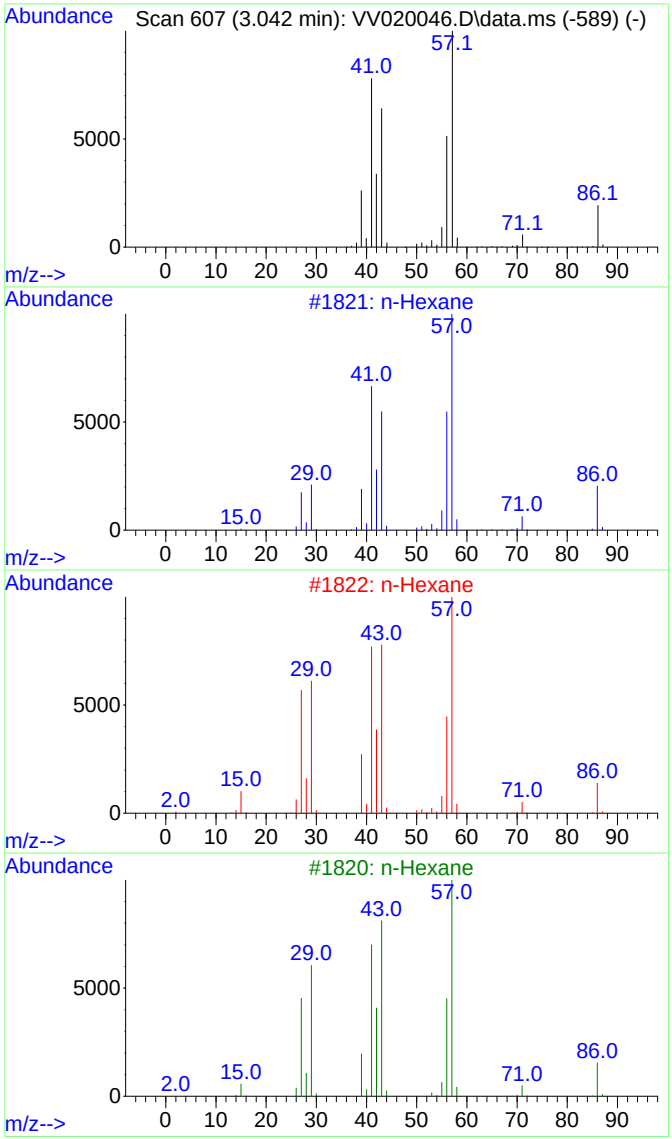
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

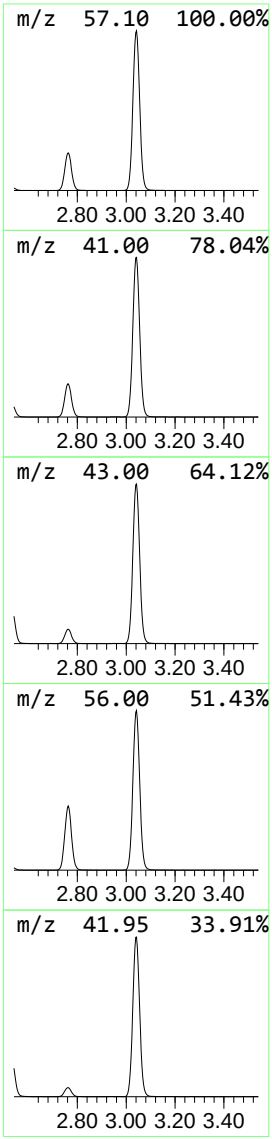
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.042 | 648.11 ug/L | 8057250 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexane | 86 | C6H14 | 000110-54-3 | 94 |
| 2 | n-Hexane | 86 | C6H14 | 000110-54-3 | 64 |
| 3 | n-Hexane | 86 | C6H14 | 000110-54-3 | 59 |
| 4 | Pentane, 2,2,3,4-tetramethyl- | 128 | C9H20 | 001186-53-4 | 38 |
| 5 | Propanal, 2,2-dimethyl- | 86 | C5H10O | 000630-19-3 | 38 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

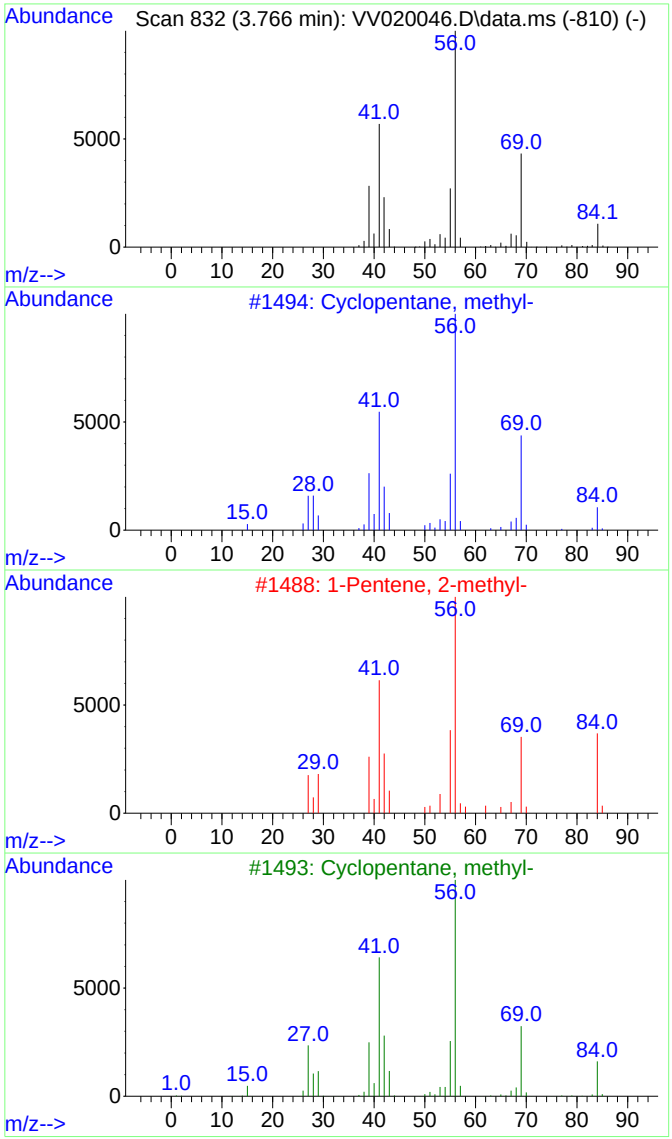
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

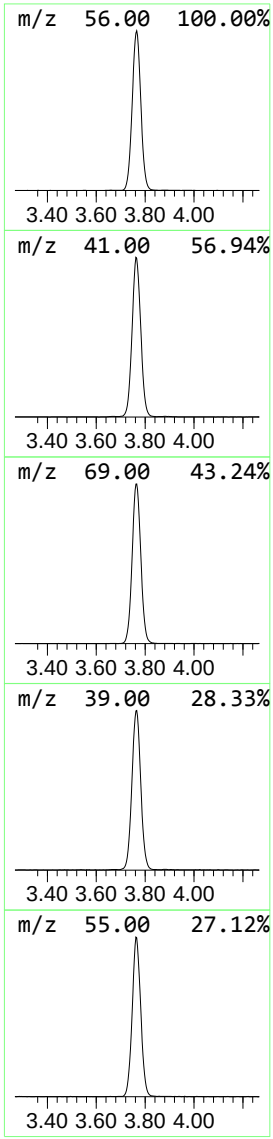
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 5 (DEL) Alkane: Cyclic3.766 Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.766 | 277.01 ug/L | 3443810 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclopentane, methyl- | 84 | C6H12 | 000096-37-7 | 94 |
| 2 | 1-Pentene, 2-methyl- | 84 | C6H12 | 000763-29-1 | 87 |
| 3 | Cyclopentane, methyl- | 84 | C6H12 | 000096-37-7 | 86 |
| 4 | Cyclopentane, methyl- | 84 | C6H12 | 000096-37-7 | 78 |
| 5 | 1H-Tetrazole, 5-methyl- | 84 | C2H4N4 | 004076-36-2 | 72 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

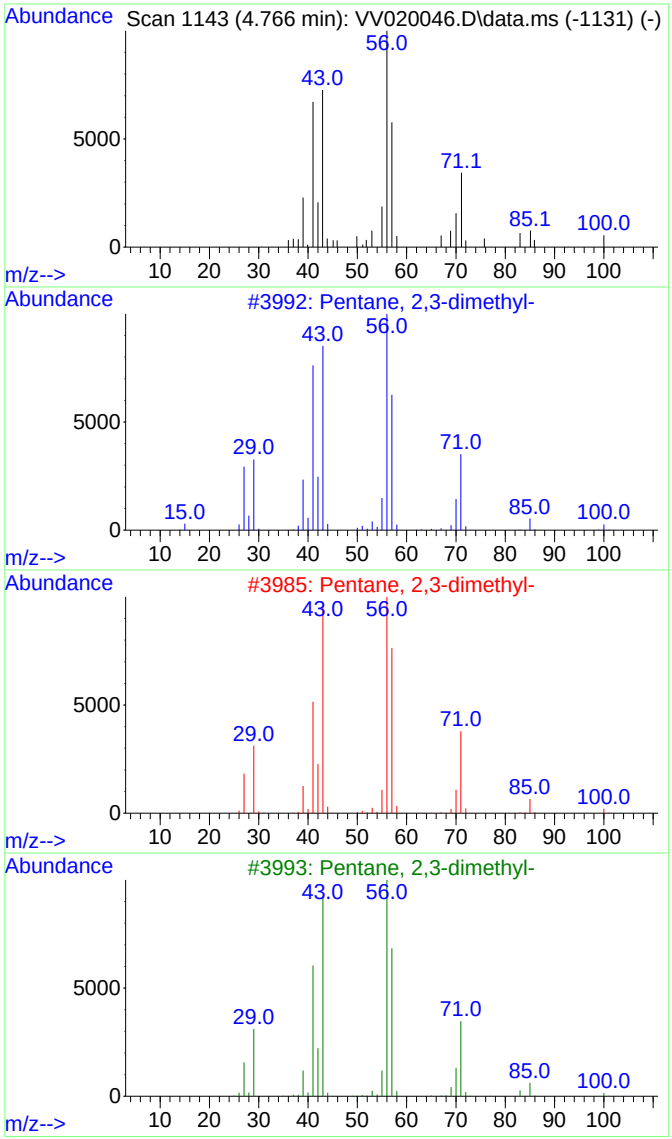
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

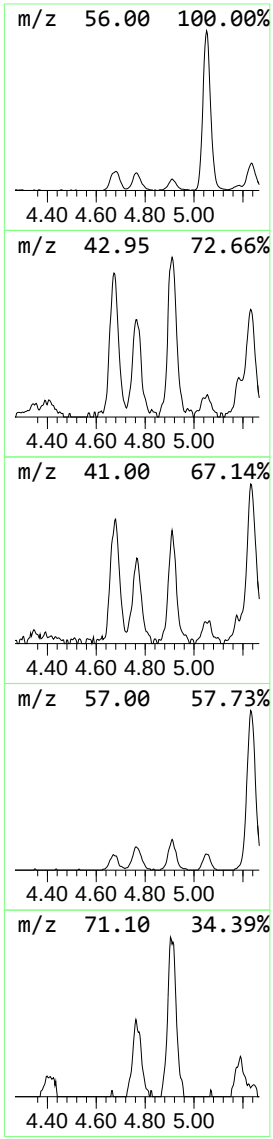
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.766 | 3.72 ug/L | 46274 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentane, 2,3-dimethyl- | 100 | C7H16 | 000565-59-3 | 78 |
| 2 | Pentane, 2,3-dimethyl- | 100 | C7H16 | 000565-59-3 | 78 |
| 3 | Pentane, 2,3-dimethyl- | 100 | C7H16 | 000565-59-3 | 78 |
| 4 | Pentane, 2,3-dimethyl- | 100 | C7H16 | 000565-59-3 | 56 |
| 5 | Aziridine, 2-methyl- | 57 | C3H7N | 000075-55-8 | 53 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :**
MSVOA_V
**ClientSampleId :**
BG4Y7

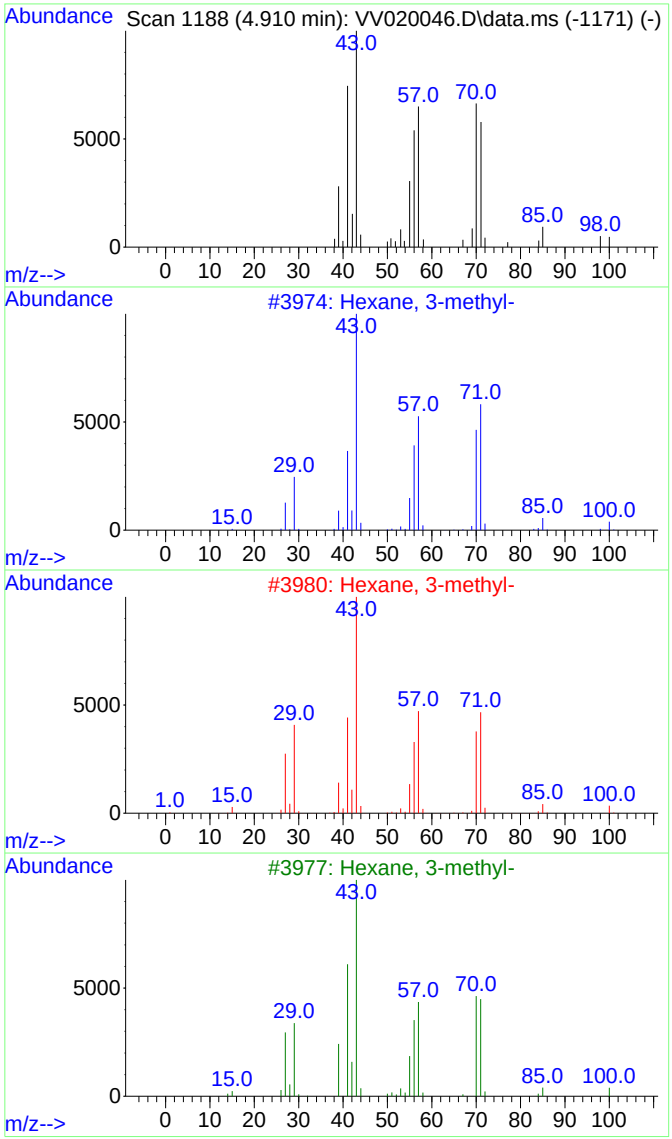
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

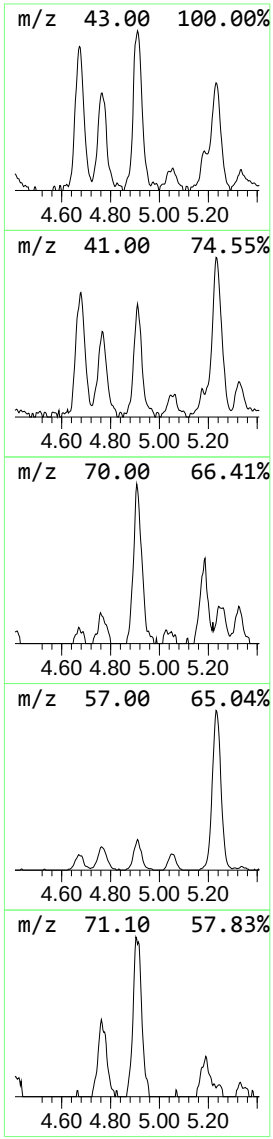
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.910 | 5.33 ug/L | 66221 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexane, 3-methyl- | 100 | C7H16 | 000589-34-4 | 91 |
| 2 | Hexane, 3-methyl- | 100 | C7H16 | 000589-34-4 | 87 |
| 3 | Hexane, 3-methyl- | 100 | C7H16 | 000589-34-4 | 81 |
| 4 | Heptane, 3,4-dimethyl- | 128 | C9H20 | 000922-28-1 | 72 |
| 5 | Hexane, 3-ethyl-4-methyl- | 128 | C9H20 | 003074-77-9 | 59 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :**
MSVOA_V
**ClientSampleId :**
BG4Y7

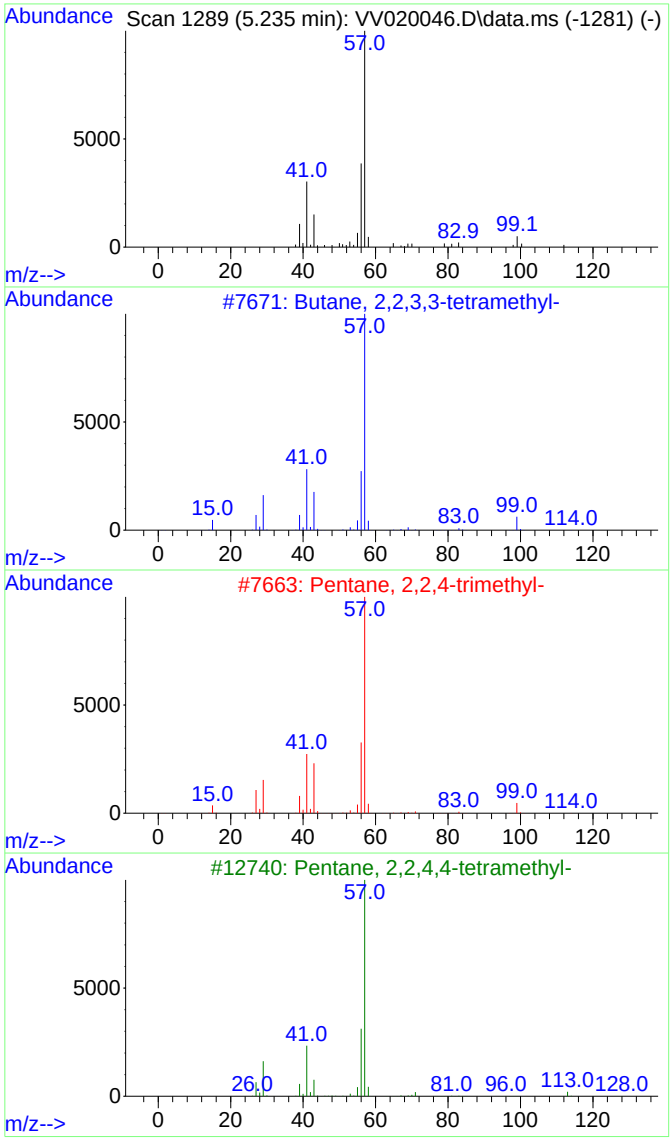
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

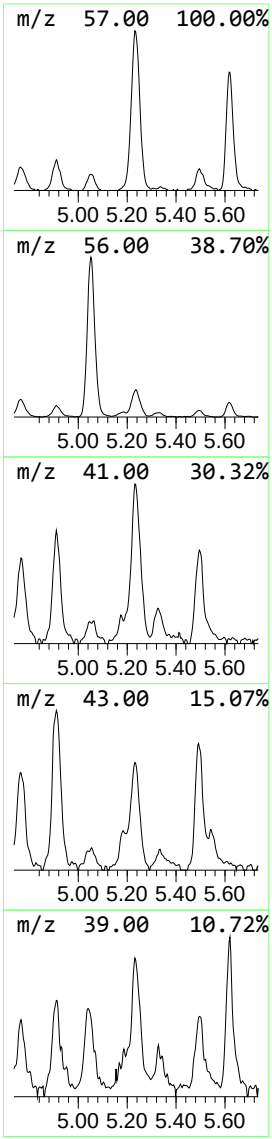
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.235 | 7.66 ug/L | 95193 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2,2,3,3-tetramethyl- | 114 | C8H18 | 000594-82-1 | 72 |
| 2 | Pentane, 2,2,4-trimethyl- | 114 | C8H18 | 000540-84-1 | 72 |
| 3 | Pentane, 2,2,4,4-tetramethyl- | 128 | C9H20 | 001070-87-7 | 64 |
| 4 | Pentane, 2,2,3-trimethyl- | 114 | C8H18 | 000564-02-3 | 59 |
| 5 | Pentane, 2,2,4-trimethyl- | 114 | C8H18 | 000540-84-1 | 56 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

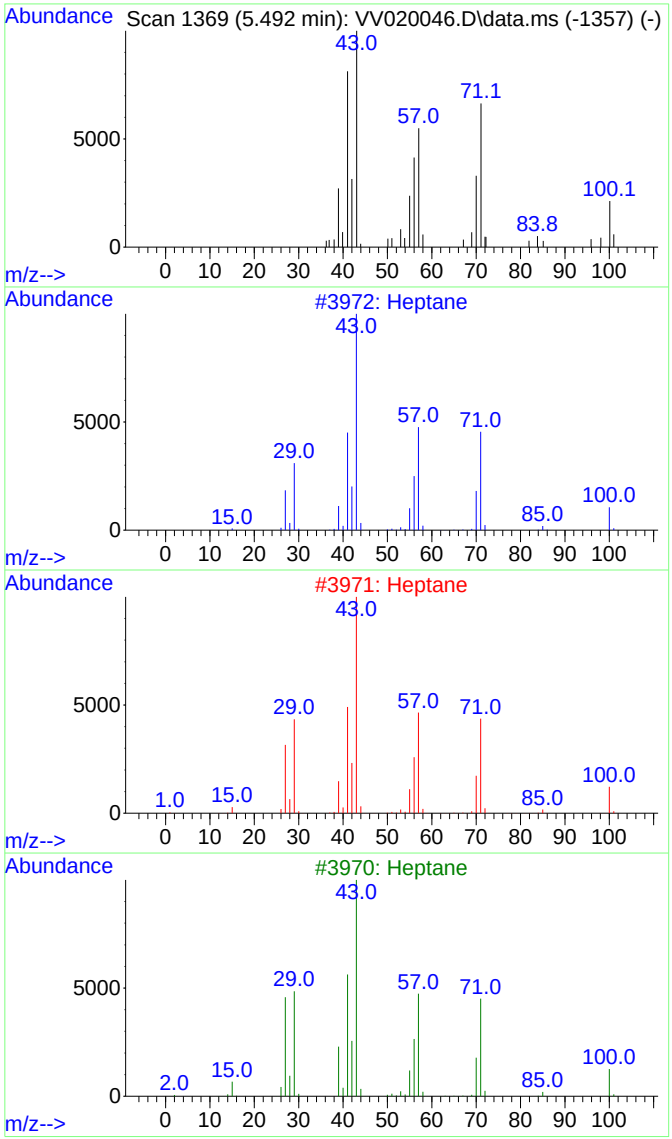
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

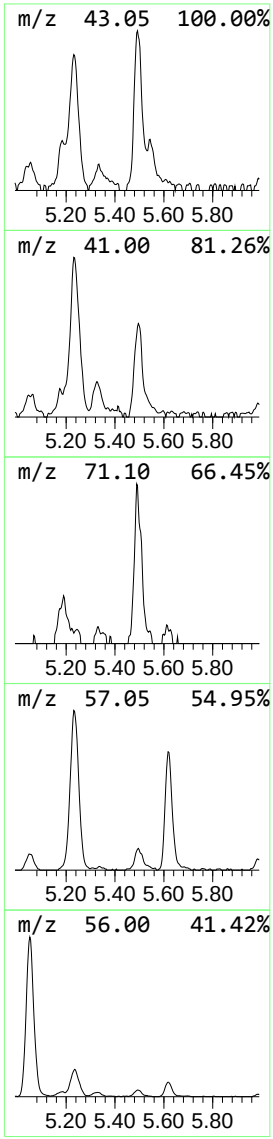
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.492 | 4.07 ug/L | 50620 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptane | 100 | C7H16 | 000142-82-5 | 80 |
| 2 | Heptane | 100 | C7H16 | 000142-82-5 | 72 |
| 3 | Heptane | 100 | C7H16 | 000142-82-5 | 72 |
| 4 | Heptane | 100 | C7H16 | 000142-82-5 | 59 |
| 5 | 2,2-Dimethylperhydro-1,3-oxazine | 115 | C6H13NO | 073155-29-0 | 50 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :**
MSVOA_V
**ClientSampleId :**
BG4Y7

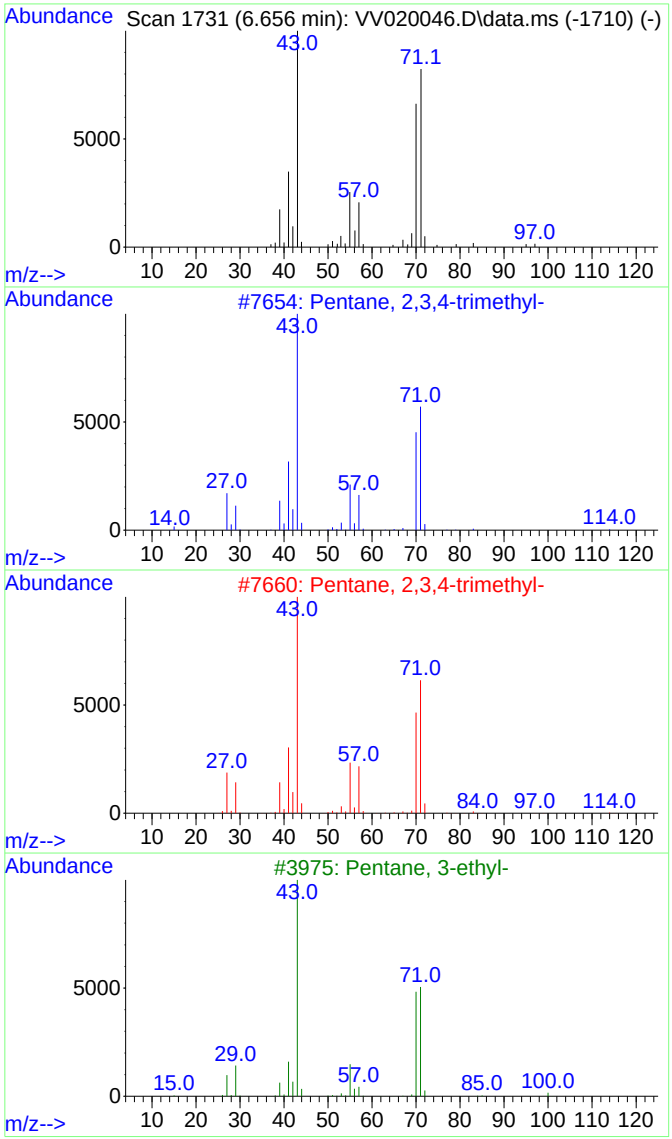
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

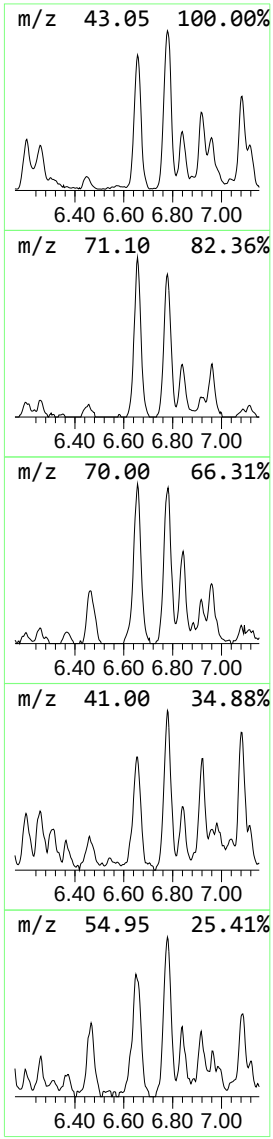
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.656 | 8.12 ug/L | 100983 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentane, 2,3,4-trimethyl- | 114 | C8H18 | 000565-75-3 | 90 |
| 2 | Pentane, 2,3,4-trimethyl- | 114 | C8H18 | 000565-75-3 | 83 |
| 3 | Pentane, 3-ethyl- | 100 | C7H16 | 000617-78-7 | 83 |
| 4 | Pentane, 2,3,4-trimethyl- | 114 | C8H18 | 000565-75-3 | 83 |
| 5 | Pentane, 2,3,3-trimethyl- | 114 | C8H18 | 000560-21-4 | 78 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

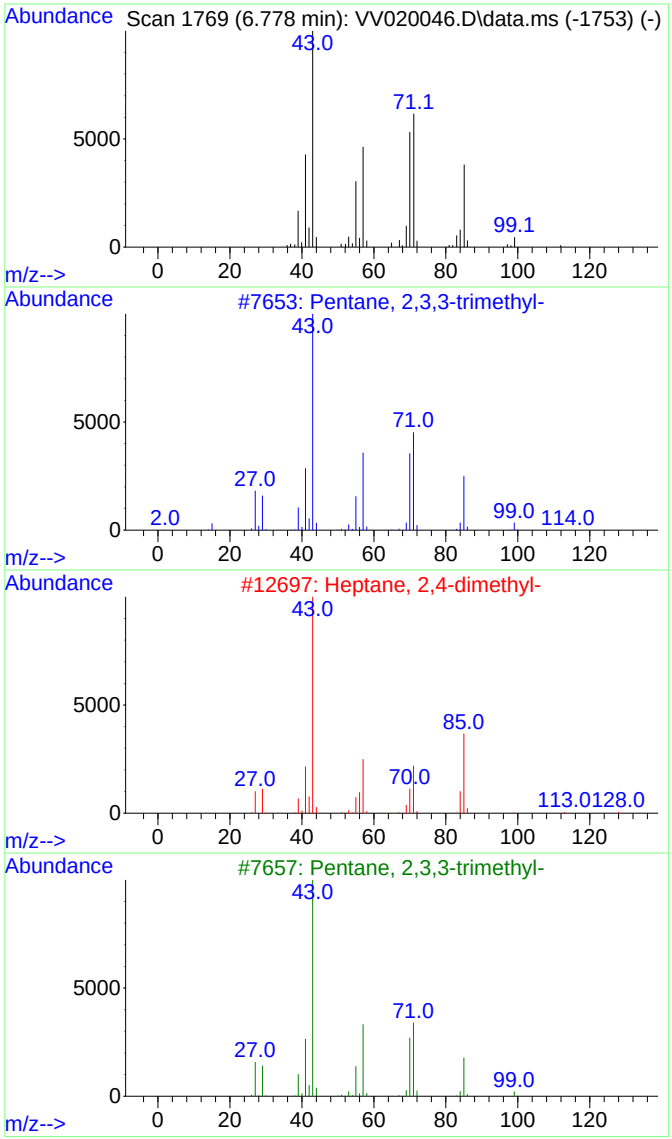
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

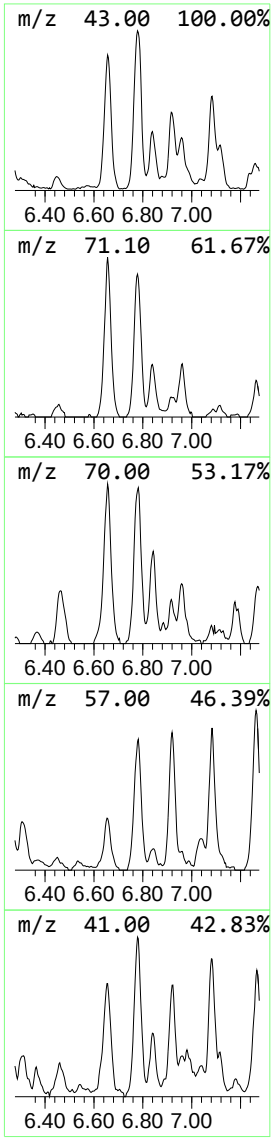
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.778 | 10.16 ug/L | 126265 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentane, 2,3,3-trimethyl- | 114 | C8H18 | 000560-21-4 | 83 |
| 2 | Heptane, 2,4-dimethyl- | 128 | C9H20 | 002213-23-2 | 64 |
| 3 | Pentane, 2,3,3-trimethyl- | 114 | C8H18 | 000560-21-4 | 64 |
| 4 | Hexane, 2,3,4-trimethyl- | 128 | C9H20 | 000921-47-1 | 64 |
| 5 | Oxalic acid, isohexyl neopentyl ... | 244 | C13H24O4 | 1000309-73-0 | 53 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

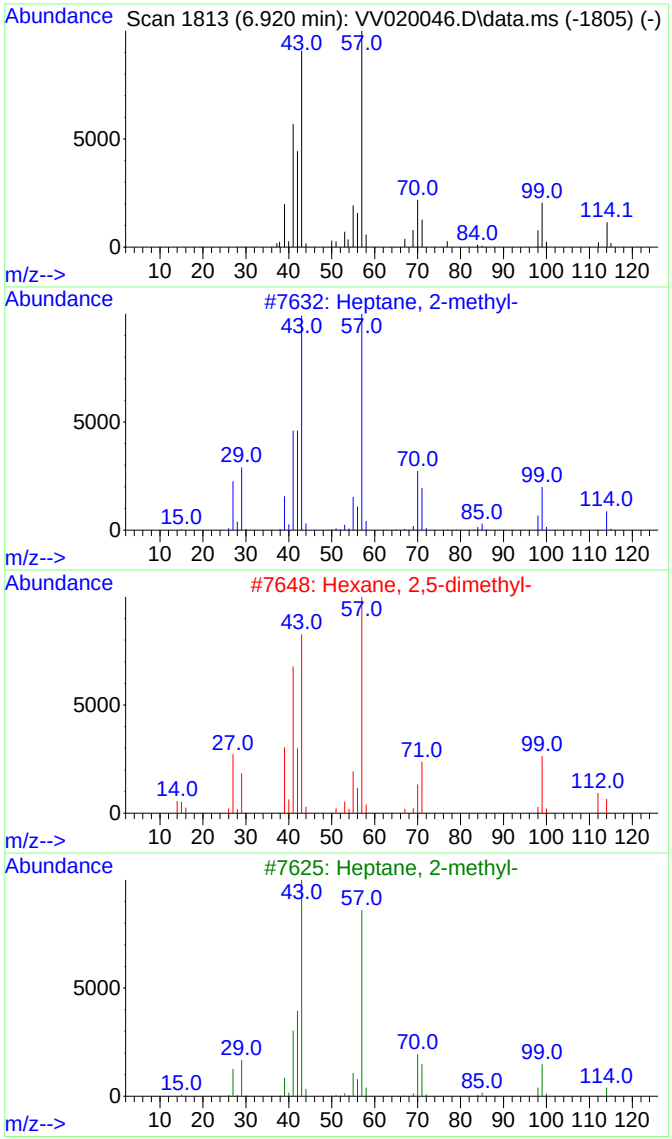
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

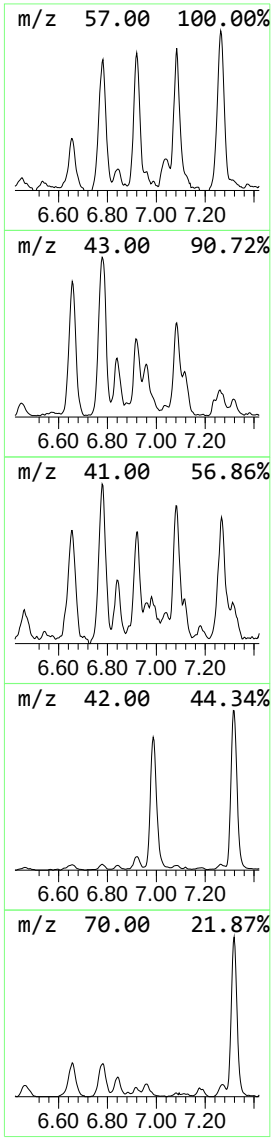
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.920 | 3.54 ug/L | 44032 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptane, 2-methyl- | 114 | C8H18 | 000592-27-8 | 72 |
| 2 | Hexane, 2,5-dimethyl- | 114 | C8H18 | 000592-13-2 | 58 |
| 3 | Heptane, 2-methyl- | 114 | C8H18 | 000592-27-8 | 56 |
| 4 | Hexane, 2,5-dimethyl- | 114 | C8H18 | 000592-13-2 | 52 |
| 5 | Heptane, 1,1'-oxybis- | 214 | C14H30O | 000629-64-1 | 50 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :**
MSVOA_V
**ClientSampleId :**
BG4Y7

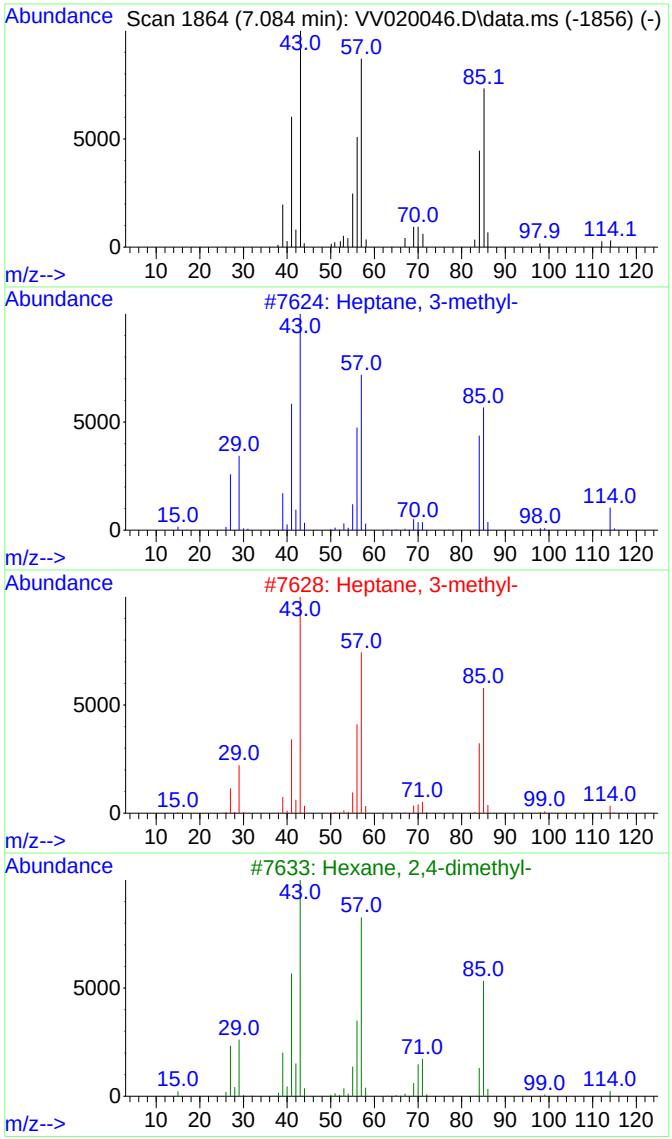
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

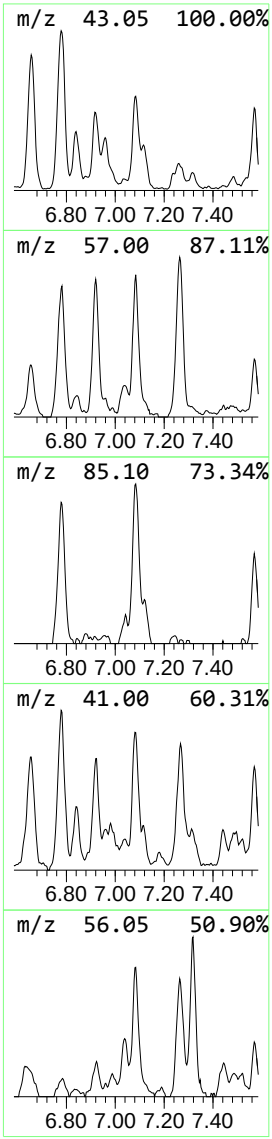
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.084 | 4.30 ug/L | 53437 | 1,4-Difluorobenzene | 5.618 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptane, 3-methyl- | 114 | C8H18 | 000589-81-1 | 91 |
| 2 | Heptane, 3-methyl- | 114 | C8H18 | 000589-81-1 | 91 |
| 3 | Hexane, 2,4-dimethyl- | 114 | C8H18 | 000589-43-5 | 72 |
| 4 | Hexane, 2,4-dimethyl- | 114 | C8H18 | 000589-43-5 | 64 |
| 5 | Hexane, 2-methyl- | 100 | C7H16 | 000591-76-4 | 53 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

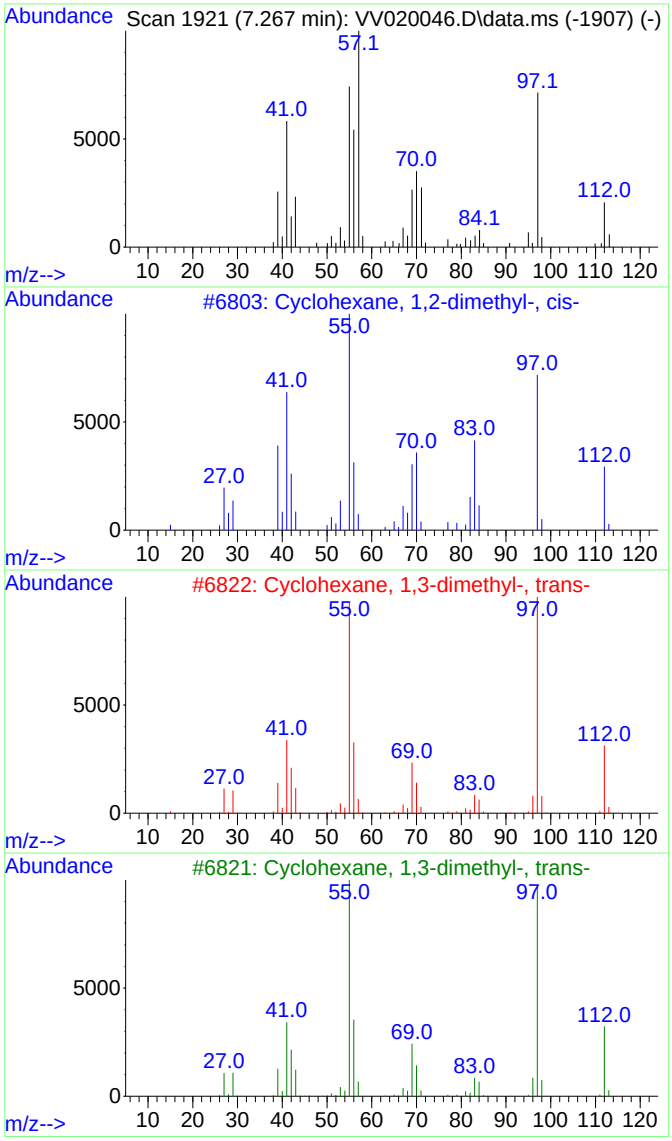
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

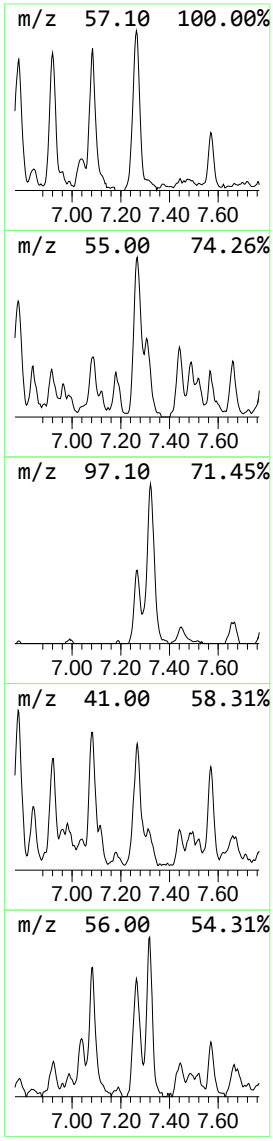
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 16 (DEL) Alkane: Cyclic7.267 Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.267 | 5.98 ug/L | 88125 | Chlorobenzene-d5 | 8.855 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclohexane, 1,2-dimethyl-, cis- | 112 | C8H16 | 002207-01-4 | 58 |
| 2 | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16 | 002207-03-6 | 49 |
| 3 | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16 | 002207-03-6 | 49 |
| 4 | 1H-Pyrazol-4-amine, 3-methyl- | 97 | C4H7N3 | 1000338-28-2 | 49 |
| 5 | 2-Octene, (Z)- | 112 | C8H16 | 007642-04-8 | 49 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

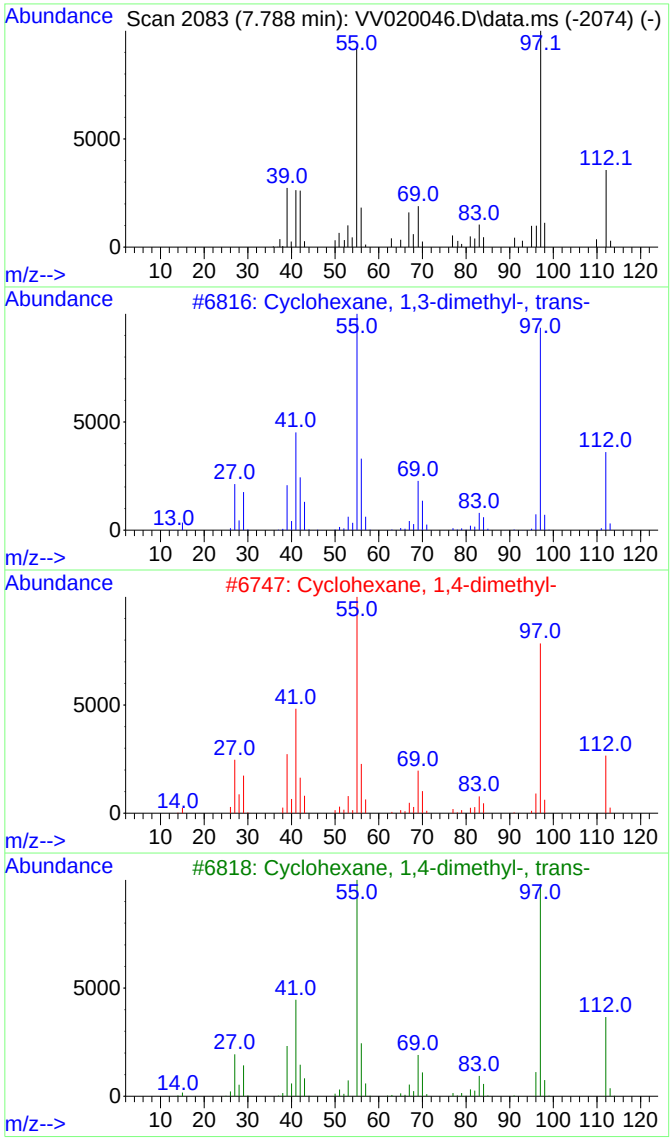
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

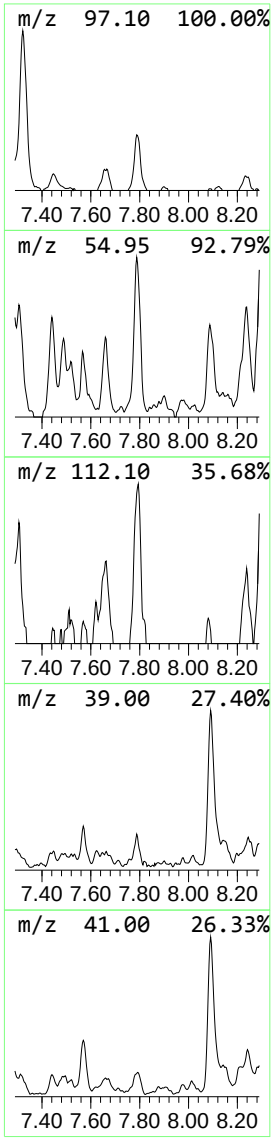
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 17 (DEL) Alkane: Cyclic7.788 Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.788 | 2.54 ug/L | 37512 | Chlorobenzene-d5 | 8.855 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16 | 002207-03-6 | 87 |
| 2 | Cyclohexane, 1,4-dimethyl- | 112 | C8H16 | 000589-90-2 | 87 |
| 3 | Cyclohexane, 1,4-dimethyl-, trans- | 112 | C8H16 | 002207-04-7 | 87 |
| 4 | Cyclohexane, 1,3-dimethyl-, cis- | 112 | C8H16 | 000638-04-0 | 87 |
| 5 | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16 | 002207-03-6 | 83 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :**
MSVOA_V
**ClientSampleId :**
BG4Y7

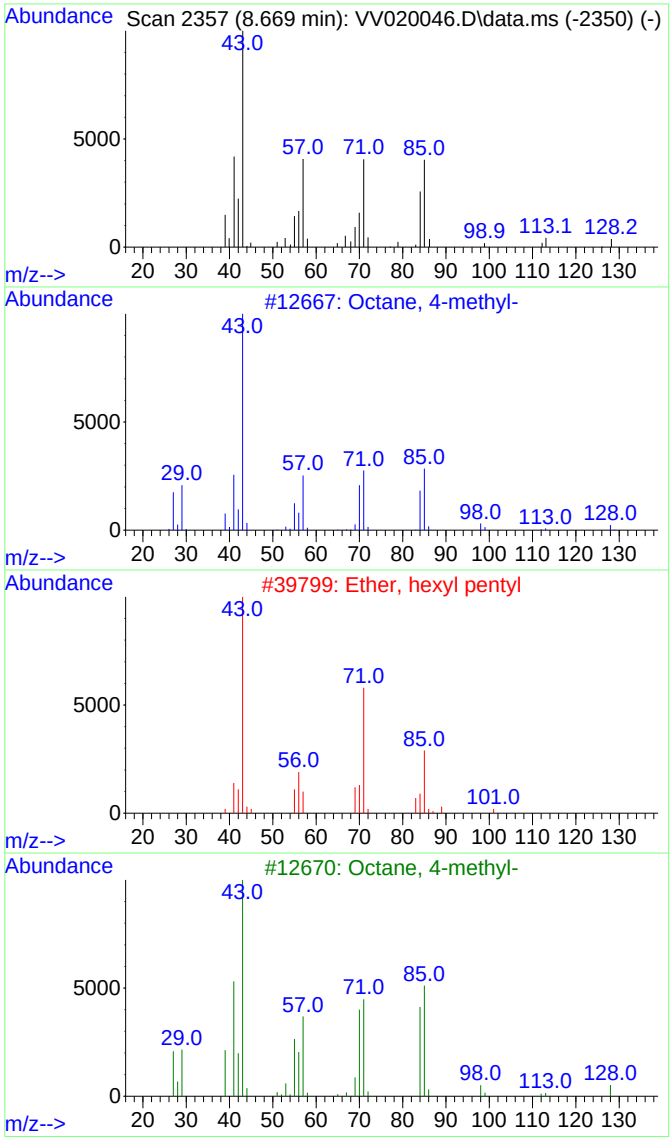
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

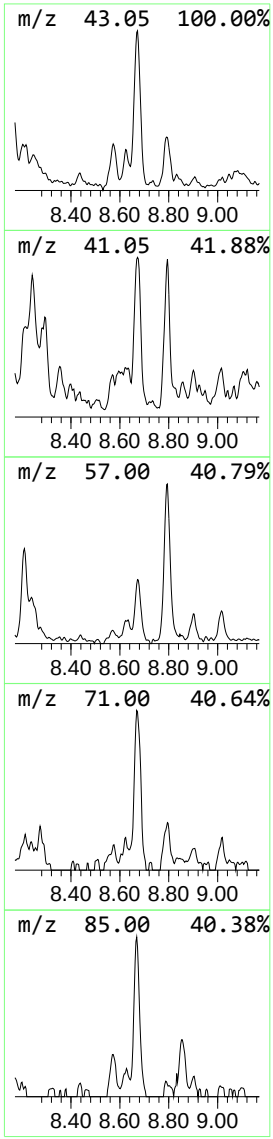
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.669 | 4.00 ug/L | 59015 | Chlorobenzene-d5 | 8.855 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octane, 4-methyl- | 128 | C9H20 | 002216-34-4 | 64 |
| 2 | Ether, hexyl pentyl | 172 | C11H24O | 032357-83-8 | 64 |
| 3 | Octane, 4-methyl- | 128 | C9H20 | 002216-34-4 | 64 |
| 4 | Octane, 4-methyl- | 128 | C9H20 | 002216-34-4 | 64 |
| 5 | Dodecane, 5-methyl- | 184 | C13H28 | 017453-93-9 | 59 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

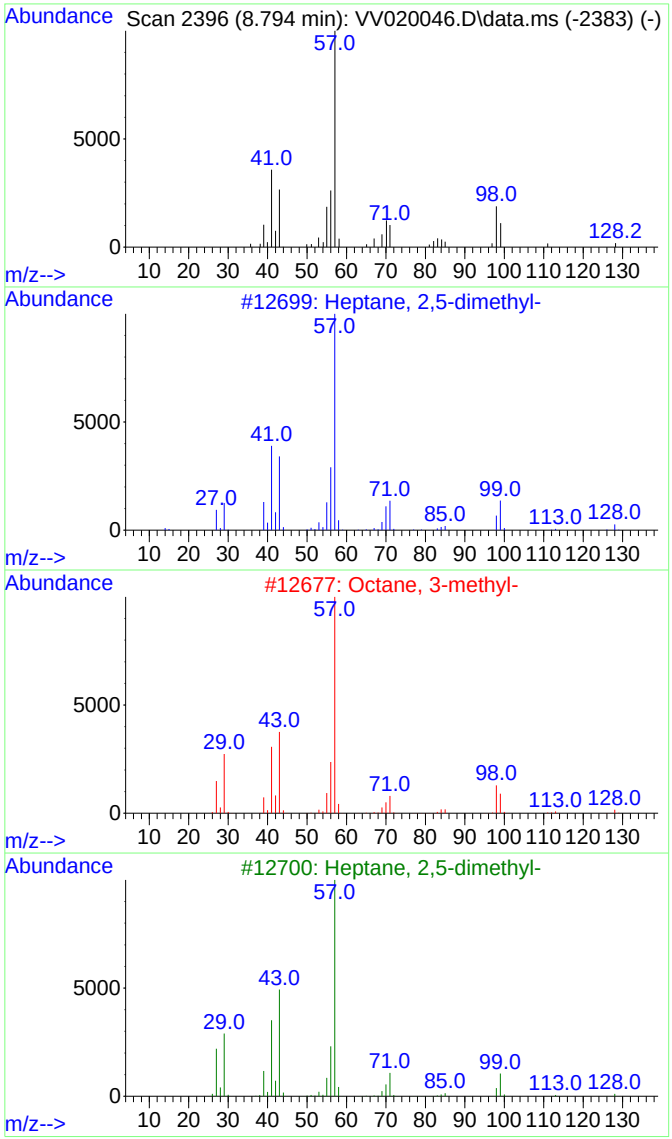
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

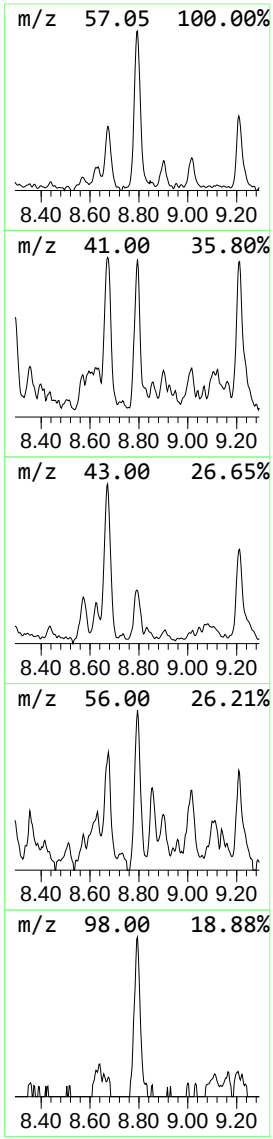
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.794 | 3.41 ug/L | 50285 | Chlorobenzene-d5 | 8.855 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptane, 2,5-dimethyl- | 128 | C9H20 | 002216-30-0 | 87 |
| 2 | Octane, 3-methyl- | 128 | C9H20 | 002216-33-3 | 87 |
| 3 | Heptane, 2,5-dimethyl- | 128 | C9H20 | 002216-30-0 | 64 |
| 4 | Octane, 3-methyl- | 128 | C9H20 | 002216-33-3 | 53 |
| 5 | Heptane, 3,5-dimethyl- | 128 | C9H20 | 000926-82-9 | 53 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

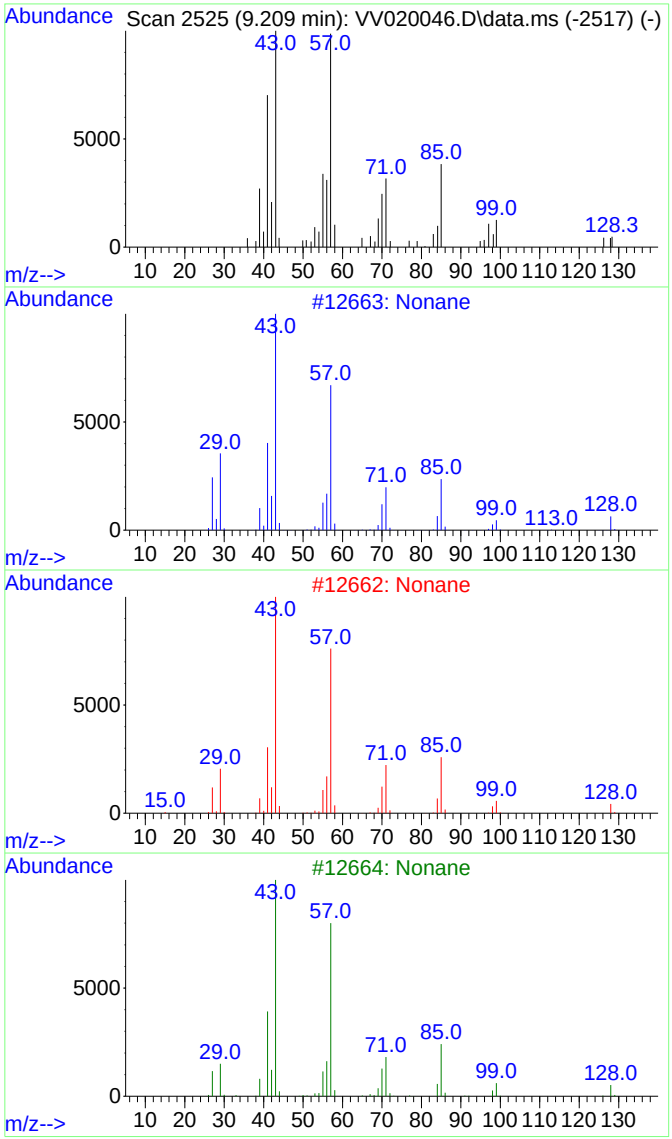
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

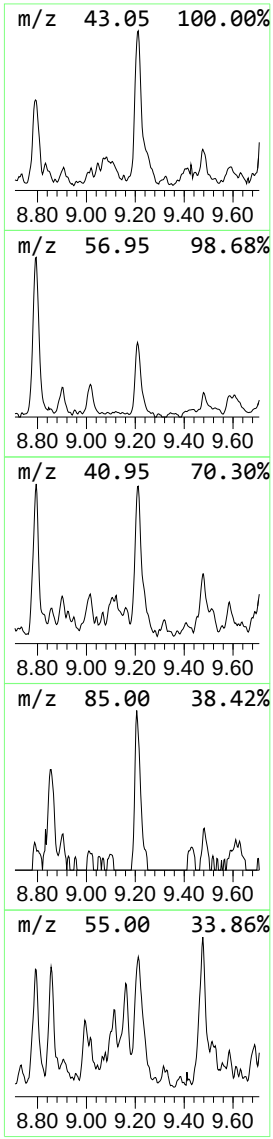
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.209 | 4.11 ug/L | 60654 | Chlorobenzene-d5 | 8.855 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Nonane | 128 | C9H20 | 000111-84-2 | 91 |
| 2 | Nonane | 128 | C9H20 | 000111-84-2 | 87 |
| 3 | Nonane | 128 | C9H20 | 000111-84-2 | 81 |
| 4 | 1-Eicosanol | 298 | C20H42O | 000629-96-9 | 47 |
| 5 | Sulfurous acid, hexyl heptyl ester | 264 | C13H28O3S | 1000309-12-9 | 47 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

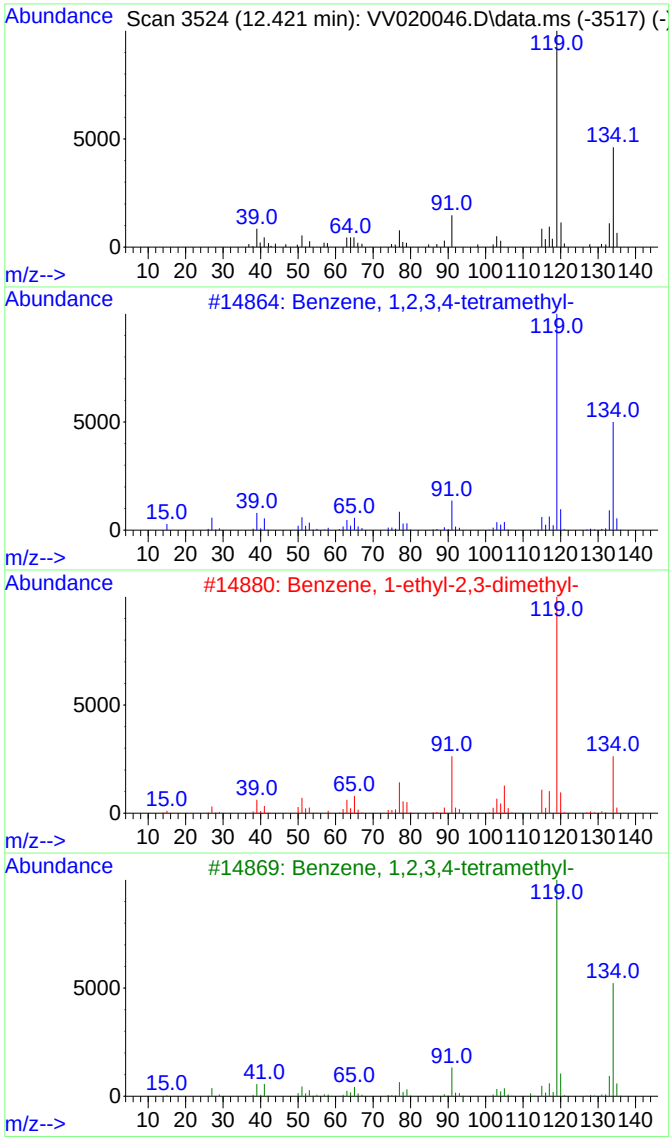
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

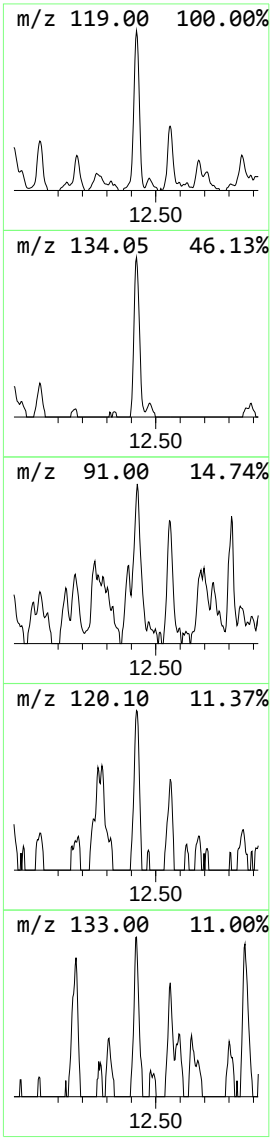
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 21 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.421 | 2.73 ug/L | 46152 | 1,4-Dichlorobenzene-d4 | 11.254 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14 | 000488-23-3 | 94 |
| 2 | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14 | 000933-98-2 | 93 |
| 3 | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14 | 000488-23-3 | 91 |
| 4 | 1,3-Cyclopentadiene, 1,2,3,4-tet... | 134 | C10H14 | 076089-59-3 | 91 |
| 5 | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14 | 000095-93-2 | 91 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011821\
Data File : VV020046.D
Acq On : 18 Jan 2021 16:18
Operator : SY/MD
Sample : M1147-15
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_V
**ClientSampleId :** BG4Y7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM011121WMA.M
Quant Title : VOC Analysis

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

| | | | | | --Internal Standard--- | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| (DEL) Alkane: S... | 2.531 | 73.9 | ug/L | 919027 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: S... | 2.762 | 135.4 | ug/L | 1683770 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: S... | 3.042 | 648.1 | ug/L | 8057250 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: C... | 3.766 | 277.0 | ug/L | 3443810 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: S... | 4.766 | 3.7 | ug/L | 46274 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: S... | 4.910 | 5.3 | ug/L | 66221 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: S... | 5.235 | 7.7 | ug/L | 95193 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: S... | 5.492 | 4.1 | ug/L | 50620 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: S... | 6.656 | 8.1 | ug/L | 100983 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: S... | 6.778 | 10.2 | ug/L | 126265 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: S... | 6.920 | 3.5 | ug/L | 44032 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: S... | 7.084 | 4.3 | ug/L | 53437 | 1 | 5.618 | 621596 | 50.0 |
| (DEL) Alkane: C... | 7.267 | 6.0 | ug/L | 88125 | 2 | 8.855 | 737395 | 50.0 |
| (DEL) Alkane: C... | 7.788 | 2.5 | ug/L | 37512 | 2 | 8.855 | 737395 | 50.0 |
| (DEL) Alkane: S... | 8.669 | 4.0 | ug/L | 59015 | 2 | 8.855 | 737395 | 50.0 |
| (DEL) Alkane: S... | 8.794 | 3.4 | ug/L | 50285 | 2 | 8.855 | 737395 | 50.0 |
| (DEL) Alkane: S... | 9.209 | 4.1 | ug/L | 60654 | 2 | 8.855 | 737395 | 50.0 |
| Benzene, 1,2,3,... | 12.421 | 2.7 | ug/L | 46152 | 3 | 11.254 | 843878 | 50.0 |