LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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_U\METHOD\SOMUTR072120WMA.M
Title : TRACE VOA SOM01.0

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 1.366 | 3 | 8 | 27 | rVB | 445037 | 443489 | 93.89% | 8.155% |
| 2 | 1.498 | 41 | 49 | 63 | rBV | 45296 | 47617 | 10.08% | 0.876% |
| 3 | 1.597 | 70 | 80 | 97 | rBV3 | 243471 | 409789 | 86.75% | 7.536% |
| 4 | 1.668 | 97 | 102 | 106 | rVV | 8735 | 10445 | 2.21% | 0.192% |
| 5 | 1.694 | 106 | 110 | 118 | rVV | 14701 | 18765 | 3.97% | 0.345% |
| 6 | 1.742 | 118 | 125 | 134 | rVB | 32461 | 38627 | 8.18% | 0.710% |
| 7 | 1.922 | 171 | 181 | 195 | rVB2 | 66057 | 99172 | 21.00% | 1.824% |
| 8 | 2.006 | 195 | 207 | 218 | rVB | 53132 | 74348 | 15.74% | 1.367% |
| 9 | 2.166 | 248 | 257 | 265 | rBV | 75935 | 103894 | 21.99% | 1.911% |
| 10 | 2.221 | 265 | 274 | 292 | rVB3 | 63671 | 130669 | 27.66% | 2.403% |
| 11 | 2.330 | 301 | 308 | 315 | rVB2 | 7679 | 9625 | 2.04% | 0.177% |
| 12 | 2.427 | 324 | 338 | 347 | rBV2 | 24829 | 39136 | 8.29% | 0.720% |
| 13 | 2.481 | 347 | 355 | 363 | rVB2 | 5494 | 7789 | 1.65% | 0.143% |
| 14 | 2.581 | 370 | 386 | 398 | rBV | 80166 | 136084 | 28.81% | 2.502% |
| 15 | 2.649 | 399 | 407 | 420 | rVB2 | 9199 | 17320 | 3.67% | 0.318% |
| 16 | 2.877 | 467 | 478 | 488 | rBV4 | 7247 | 12476 | 2.64% | 0.229% |
| 17 | 2.996 | 499 | 515 | 523 | rBV7 | 9001 | 22849 | 4.84% | 0.420% |
| 18 | 3.099 | 539 | 547 | 557 | rVV3 | 6473 | 12170 | 2.58% | 0.224% |
| 19 | 3.153 | 557 | 564 | 577 | rVB7 | 4517 | 9181 | 1.94% | 0.169% |
| 20 | 3.382 | 623 | 635 | 651 | rVB2 | 8183 | 18790 | 3.98% | 0.346% |
| 21 | 3.604 | 687 | 704 | 721 | rVB5 | 25955 | 66144 | 14.00% | 1.216% |
| 22 | 3.719 | 728 | 740 | 755 | rVB3 | 11051 | 23711 | 5.02% | 0.436% |
| 23 | 3.851 | 768 | 781 | 793 | rBV6 | 5007 | 10650 | 2.25% | 0.196% |
| 24 | 3.919 | 793 | 802 | 807 | rBV6 | 3277 | 5890 | 1.25% | 0.108% |
| 25 | 4.215 | 881 | 894 | 908 | rVB5 | 6272 | 15213 | 3.22% | 0.280% |
| 26 | 4.472 | 957 | 974 | 985 | rBV4 | 4729 | 11607 | 2.46% | 0.213% |
| 27 | 4.649 | 1013 | 1029 | 1048 | rBV | 69903 | 182329 | 38.60% | 3.353% |
| 28 | 5.086 | 1150 | 1165 | 1187 | rVB2 | 71958 | 175793 | 37.22% | 3.233% |
| 29 | 5.465 | 1270 | 1283 | 1303 | rVB5 | 4604 | 13658 | 2.89% | 0.251% |
| 30 | 5.610 | 1314 | 1328 | 1342 | rVB5 | 5418 | 11916 | 2.52% | 0.219% |
| 31 | 5.742 | 1348 | 1369 | 1397 | rBV2 | 148169 | 406485 | 86.05% | 7.475% |
| 32 | 5.980 | 1434 | 1443 | 1445 | rBV4 | 5375 | 6152 | 1.30% | 0.113% |
| 33 | 6.034 | 1453 | 1460 | 1478 | rVB7 | 9202 | 22015 | 4.66% | 0.405% |
| 34 | 6.153 | 1487 | 1497 | 1505 | rBV5 | 6113 | 12458 | 2.64% | 0.229% |

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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_U\METHOD\SOMUTR072120WMA.M
Title : TRACE VOA SOM01.0

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 6.266 | 1515 | 1532 | 1553 | rBV | 155961 | 307497 | 65.10% | 5.655% |
| 36 | 6.568 | 1613 | 1626 | 1635 | rVB8 | 4333 | 7886 | 1.67% | 0.145% |
| 37 | 6.706 | 1653 | 1669 | 1682 | rBV2 | 95199 | 185981 | 39.37% | 3.420% |
| 38 | 6.777 | 1682 | 1691 | 1704 | rVB8 | 5636 | 12522 | 2.65% | 0.230% |
| 39 | 6.890 | 1716 | 1726 | 1737 | rBV5 | 5382 | 11555 | 2.45% | 0.212% |
| 40 | 7.427 | 1883 | 1893 | 1904 | rBV9 | 2397 | 5933 | 1.26% | 0.109% |
| 41 | 7.578 | 1928 | 1940 | 1961 | rBV | 48233 | 88833 | 18.81% | 1.634% |
| 42 | 7.906 | 2029 | 2042 | 2055 | rBV | 183630 | 313036 | 66.27% | 5.756% |
| 43 | 7.976 | 2056 | 2064 | 2074 | rVB | 22589 | 35540 | 7.52% | 0.654% |
| 44 | 8.185 | 2119 | 2129 | 2140 | rBV | 29610 | 50763 | 10.75% | 0.933% |
| 45 | 8.558 | 2236 | 2245 | 2255 | rVB5 | 9924 | 15841 | 3.35% | 0.291% |
| 46 | 8.642 | 2255 | 2271 | 2298 | rBV | 257597 | 472360 | 100.00% | 8.686% |
| 47 | 8.793 | 2308 | 2318 | 2331 | rVB3 | 8954 | 16053 | 3.40% | 0.295% |
| 48 | 9.420 | 2501 | 2513 | 2540 | rBV | 223852 | 380968 | 80.65% | 7.006% |
| 49 | 9.574 | 2551 | 2561 | 2570 | rBV2 | 16775 | 26807 | 5.68% | 0.493% |
| 50 | 9.693 | 2584 | 2598 | 2608 | rBV2 | 8599 | 15485 | 3.28% | 0.285% |
| 51 | 10.102 | 2716 | 2725 | 2742 | rBV3 | 9936 | 18115 | 3.83% | 0.333% |
| 52 | 10.346 | 2794 | 2801 | 2810 | rVB5 | 3586 | 5482 | 1.16% | 0.101% |
| 53 | 10.558 | 2857 | 2867 | 2877 | rBV3 | 8192 | 12331 | 2.61% | 0.227% |
| 54 | 10.758 | 2917 | 2929 | 2941 | rBV | 76619 | 124286 | 26.31% | 2.285% |
| 55 | 11.465 | 3141 | 3149 | 3160 | rVB3 | 6082 | 9346 | 1.98% | 0.172% |
| 56 | 11.680 | 3208 | 3216 | 3226 | rVB7 | 4983 | 8398 | 1.78% | 0.154% |
| 57 | 11.809 | 3243 | 3256 | 3273 | rVB | 224606 | 372186 | 78.79% | 6.844% |
| 58 | 12.188 | 3362 | 3374 | 3387 | rVB | 190206 | 307654 | 65.13% | 5.657% |
| 59 | 12.880 | 3580 | 3589 | 3600 | rVB6 | 4670 | 6924 | 1.47% | 0.127% |

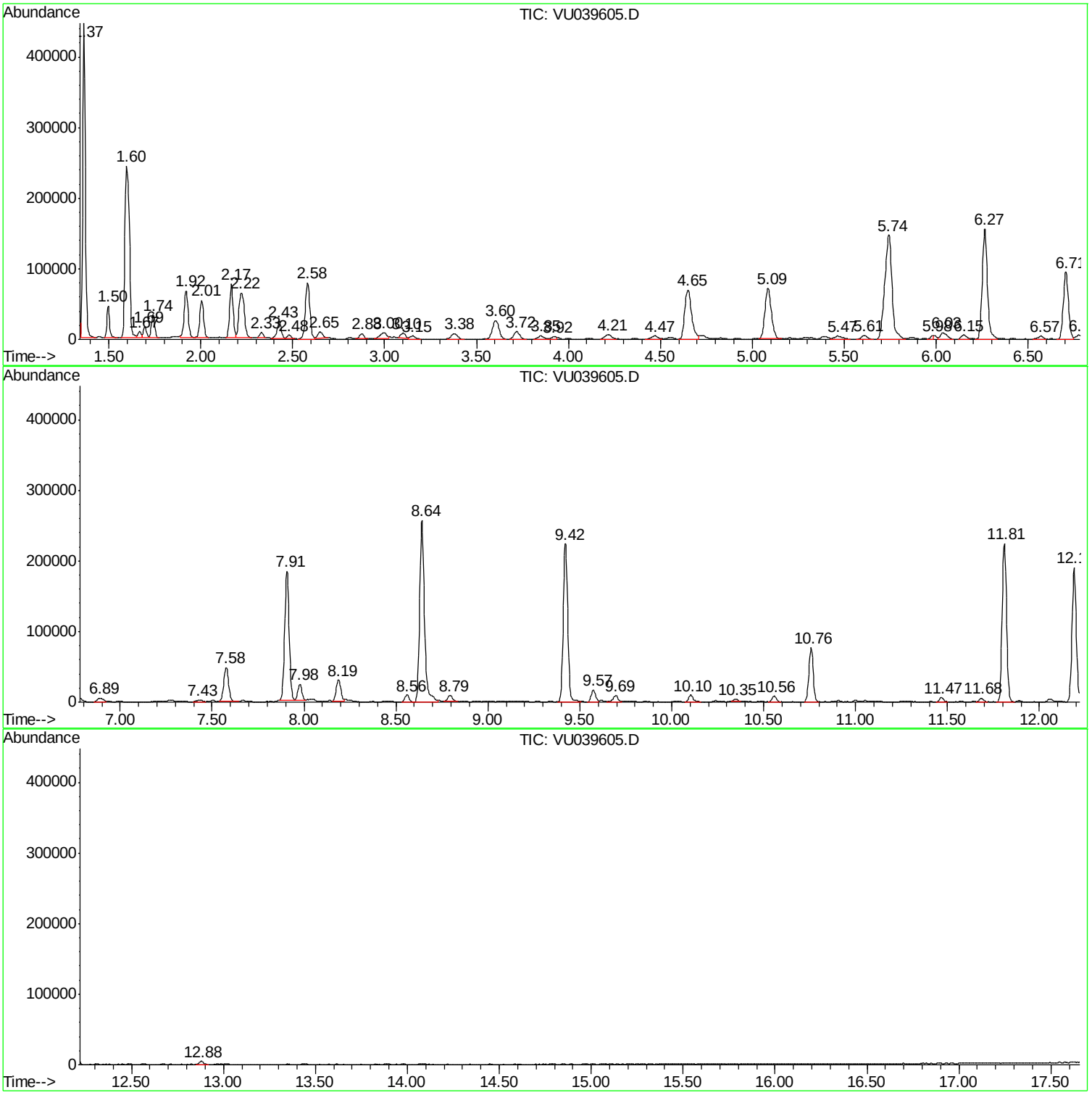
Sum of corrected areas: 5438038

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR072120WMA.M
Quant Title : TRACE VOA SOM01.0

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



# Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

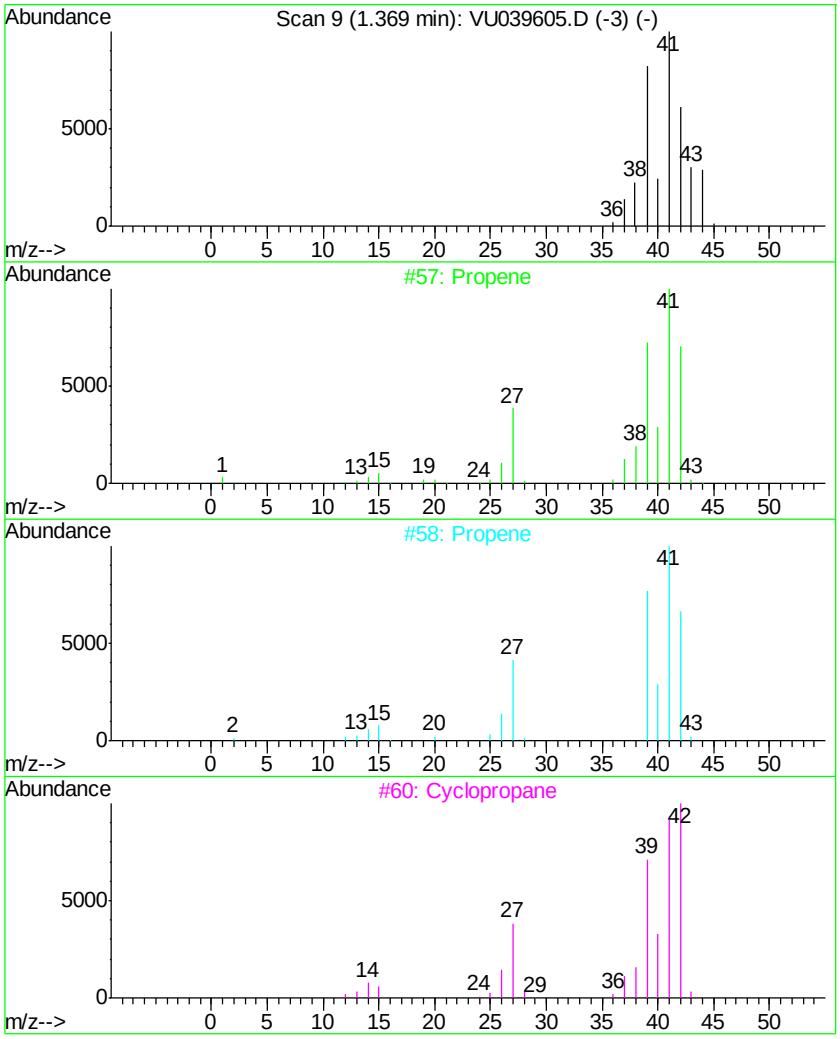
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR072120WMA.M
Quant Title : TRACE VOA SOM01.0

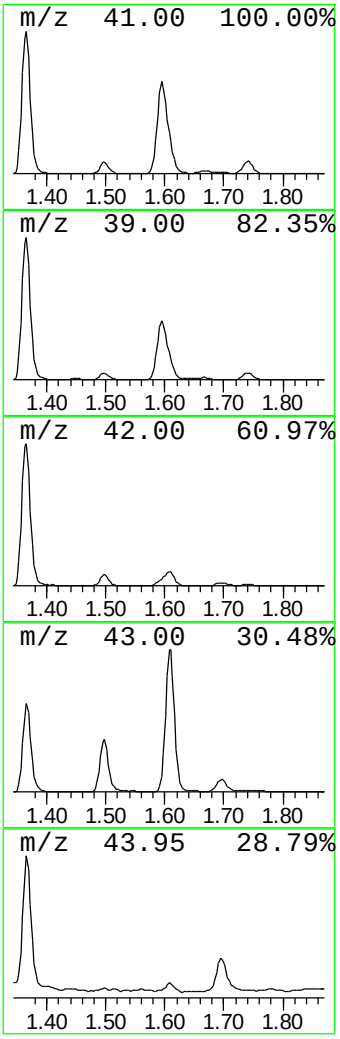
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 1.37 | 7.21 ug/L | 443489 | 1,4-Difluorobenzene | 6.27 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propene | 42 | C3H6 | 000115-07-1 | 90 |
| 2 | Propene | 42 | C3H6 | 000115-07-1 | 42 |
| 3 | Cyclopropane | 42 | C3H6 | 000075-19-4 | 9 |
| 4 | Cyclopropene | 40 | C3H4 | 002781-85-3 | 9 |
| 5 | Cyclopropane | 42 | C3H6 | 000075-19-4 | 7 |



# Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR072120WMA.M
Quant Title : TRACE VOA SOM01.0

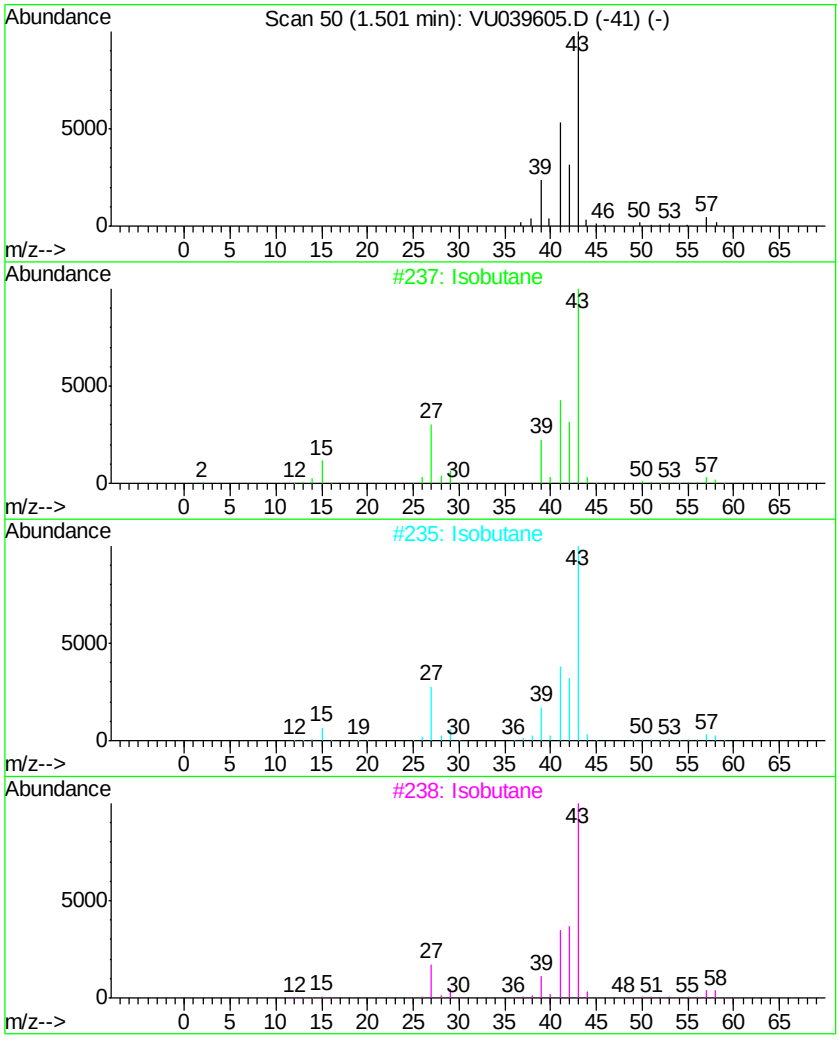
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

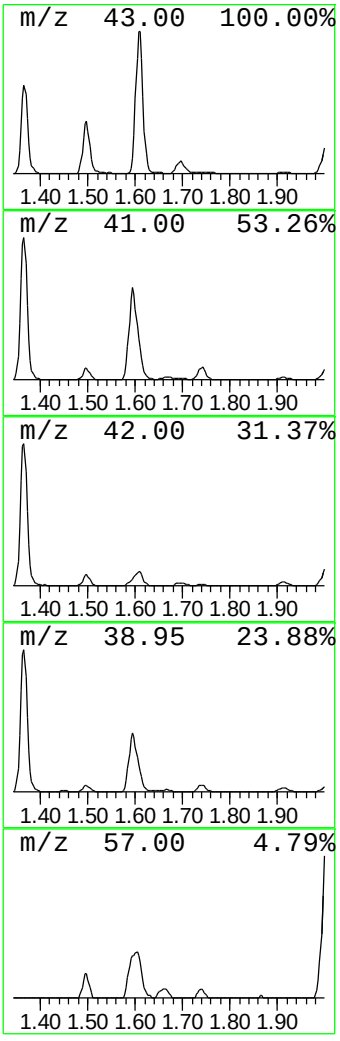
**********

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 1.50 | 0.77 ug/L | 47617 | 1,4-Difluorobenzene | 6.27 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Isobutane | 58 | C4H10 | 000075-28-5 | 59 |
| 2 | Isobutane | 58 | C4H10 | 000075-28-5 | 53 |
| 3 | Isobutane | 58 | C4H10 | 000075-28-5 | 36 |
| 4 | Isobutane | 58 | C4H10 | 000075-28-5 | 9 |
| 5 | Propane, 2-nitro- | 89 | C3H7NO2 | 000079-46-9 | 9 |



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR072120WMA.M
Quant Title : TRACE VOA SOM01.0

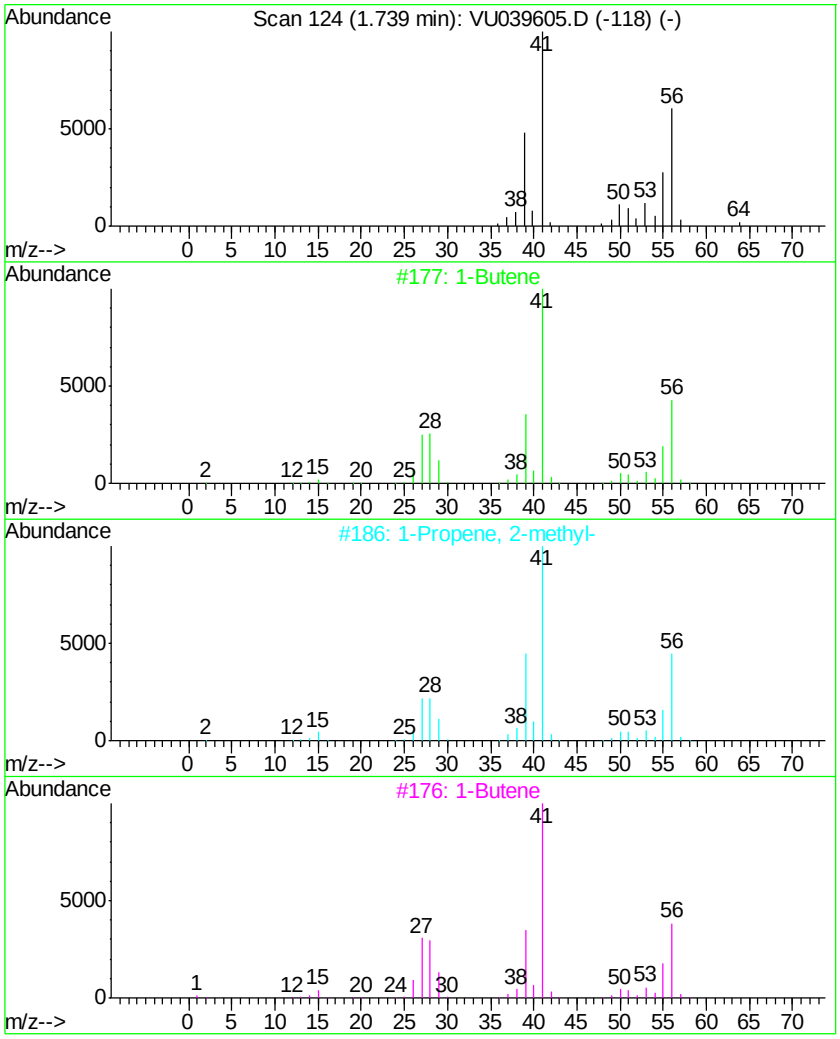
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

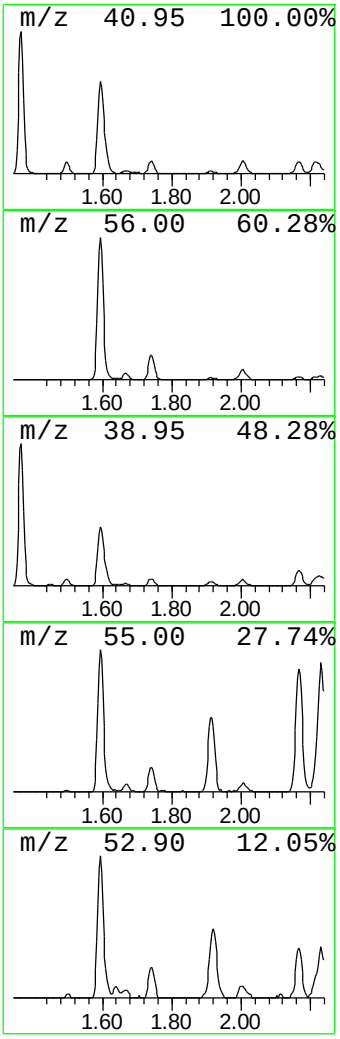
**********

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 1.74 | 0.63 ug/L | 38627 | 1,4-Difluorobenzene | 6.27 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Butene | 56 | C4H8 | 000106-98-9 | 72 |
| 2 | 1-Propene, 2-methyl- | 56 | C4H8 | 000115-11-7 | 72 |
| 3 | 1-Butene | 56 | C4H8 | 000106-98-9 | 72 |
| 4 | 2-Butene, (Z)- | 56 | C4H8 | 000590-18-1 | 64 |
| 5 | 2-Butene, (E)- | 56 | C4H8 | 000624-64-6 | 64 |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR072120WMA.M
Quant Title : TRACE VOA SOM01.0

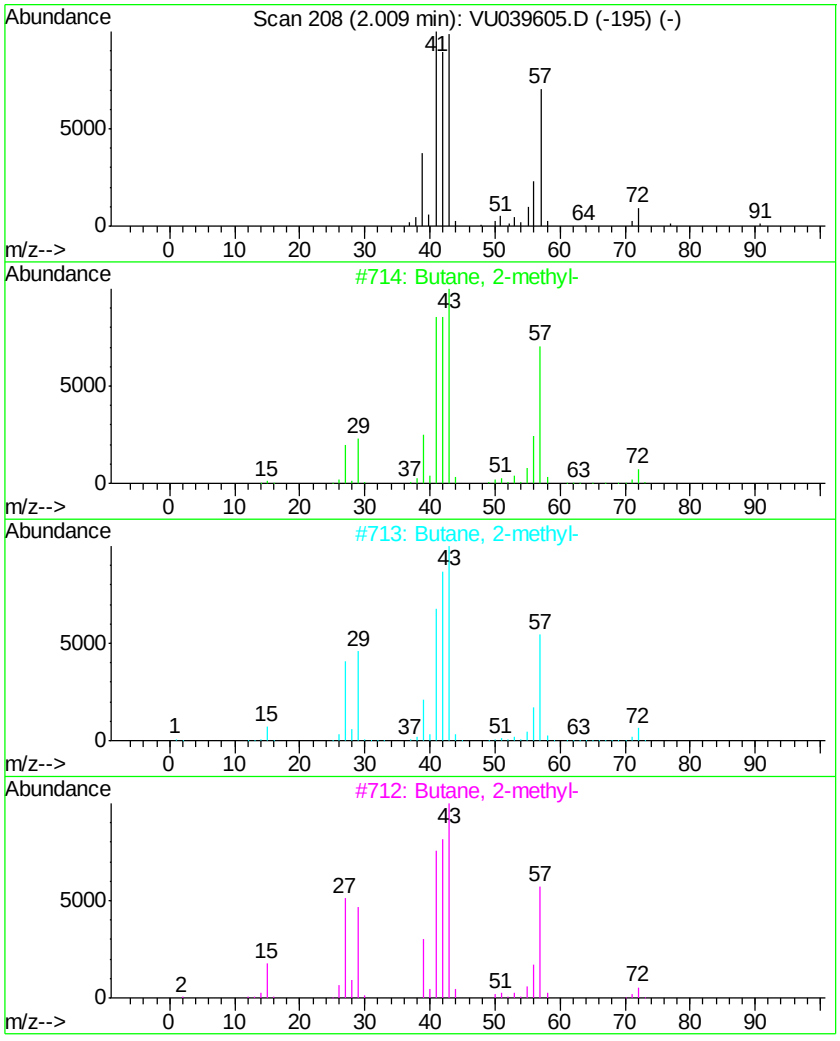
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

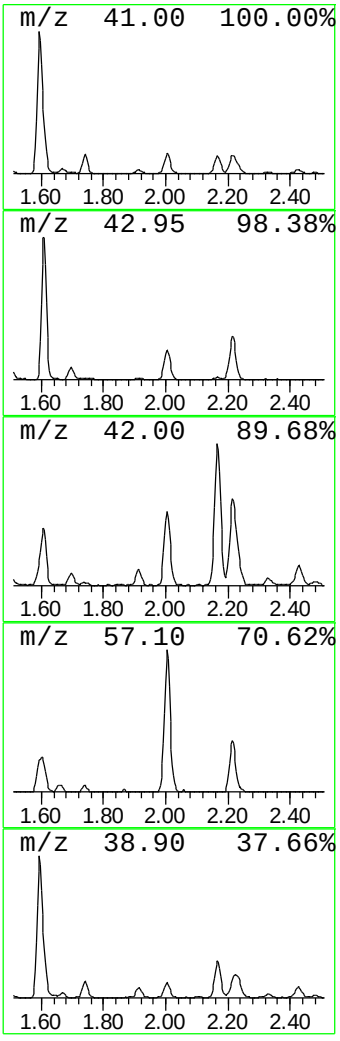
**********

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.01 | 1.21 ug/L | 74348 | 1,4-Difluorobenzene | 6.27 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2-methyl- | 72 | C5H12 | 000078-78-4 | 81 |
| 2 | Butane, 2-methyl- | 72 | C5H12 | 000078-78-4 | 72 |
| 3 | Butane, 2-methyl- | 72 | C5H12 | 000078-78-4 | 64 |
| 4 | 3-Buten-1-ol | 72 | C4H8O | 000627-27-0 | 33 |
| 5 | Pentane | 72 | C5H12 | 000109-66-0 | 25 |



# Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR072120WMA.M
Quant Title : TRACE VOA SOM01.0

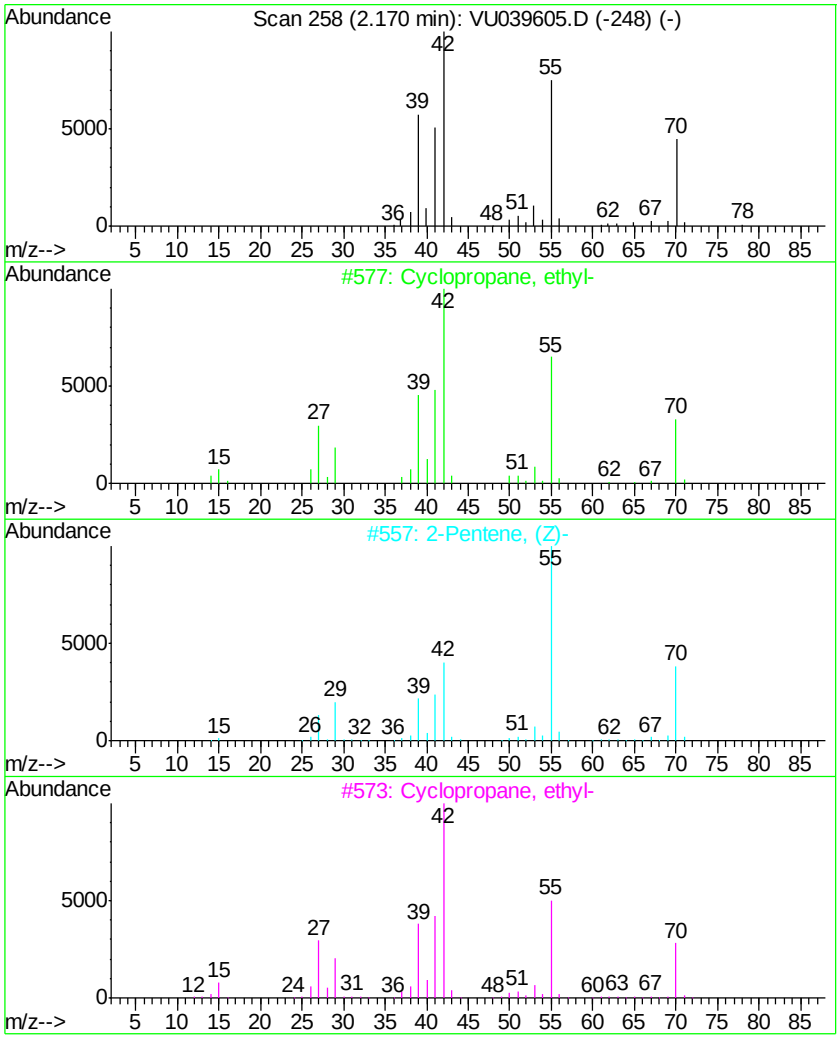
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

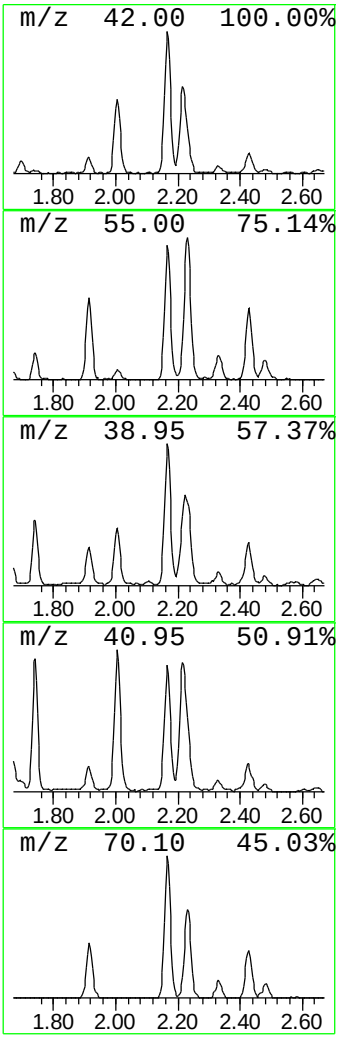
**********

Peak Number 5 (DEL) Alkane: Cyclic2.17 Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.17 | 1.69 ug/L | 103894 | 1,4-Difluorobenzene | 6.27 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclopropane, ethyl- | 70 | C5H10 | 001191-96-4 | 91 |
| 2 | 2-Pentene, (Z)- | 70 | C5H10 | 000627-20-3 | 83 |
| 3 | Cyclopropane, ethyl- | 70 | C5H10 | 001191-96-4 | 68 |
| 4 | 1-Pentene | 70 | C5H10 | 000109-67-1 | 64 |
| 5 | Cyclobutane, methyl- | 70 | C5H10 | 000598-61-8 | 64 |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

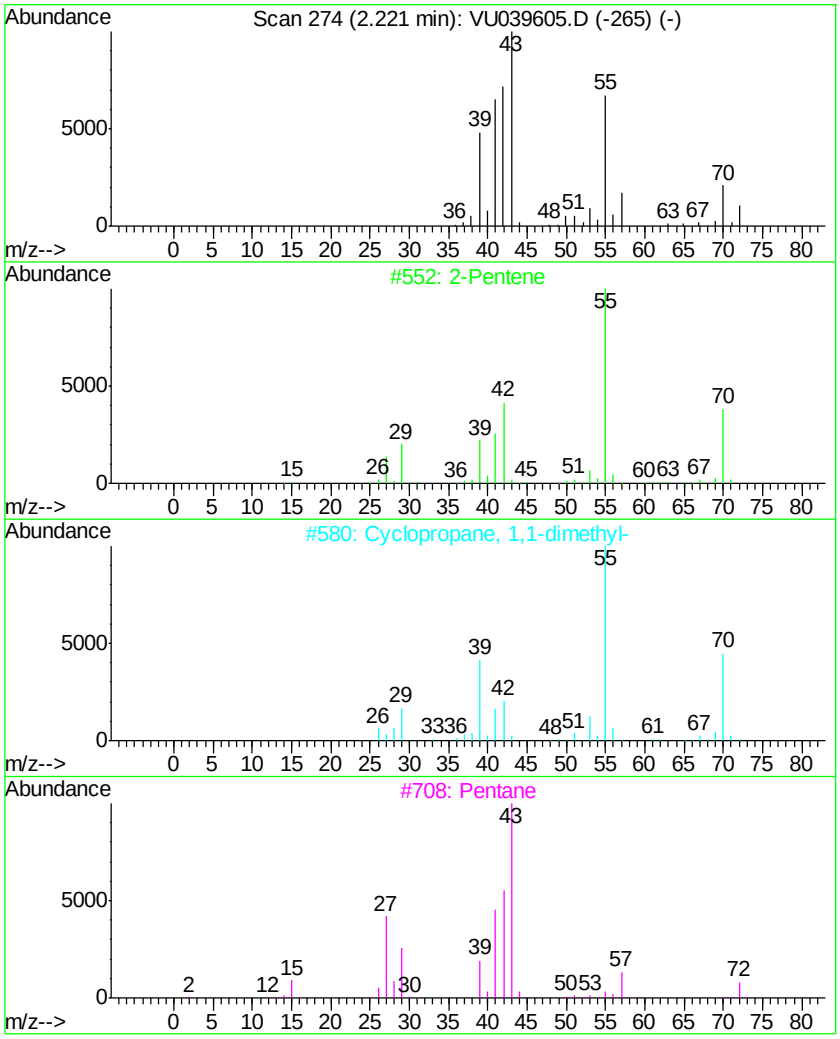
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR072120WMA.M
Quant Title : TRACE VOA SOM01.0

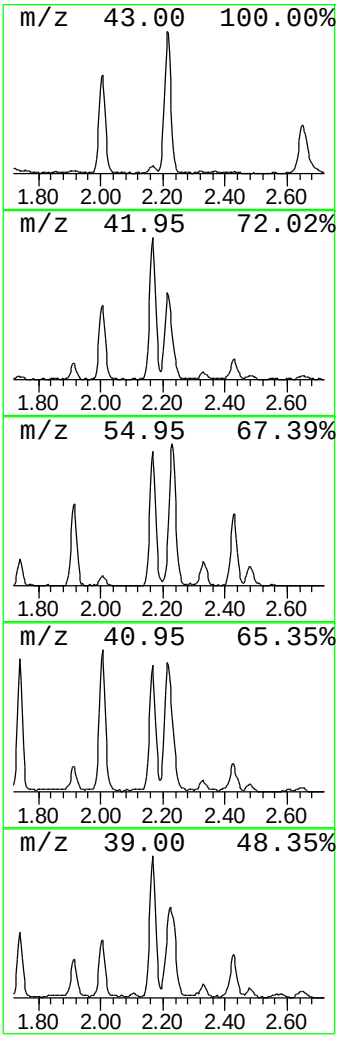
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.22 | 2.12 ug/L | 130669 | 1,4-Difluorobenzene | 6.27 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentene | 70 | C5H10 | 000109-68-2 | 52 |
| 2 | Cyclopropane, 1,1-dimethyl- | 70 | C5H10 | 001630-94-0 | 47 |
| 3 | Pentane | 72 | C5H12 | 000109-66-0 | 43 |
| 4 | Cyclopropane, ethyl- | 70 | C5H10 | 001191-96-4 | 38 |
| 5 | Pentane, 1-chloro- | 106 | C5H11Cl | 000543-59-9 | 38 |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR072120WMA.M
Quant Title : TRACE VOA SOM01.0

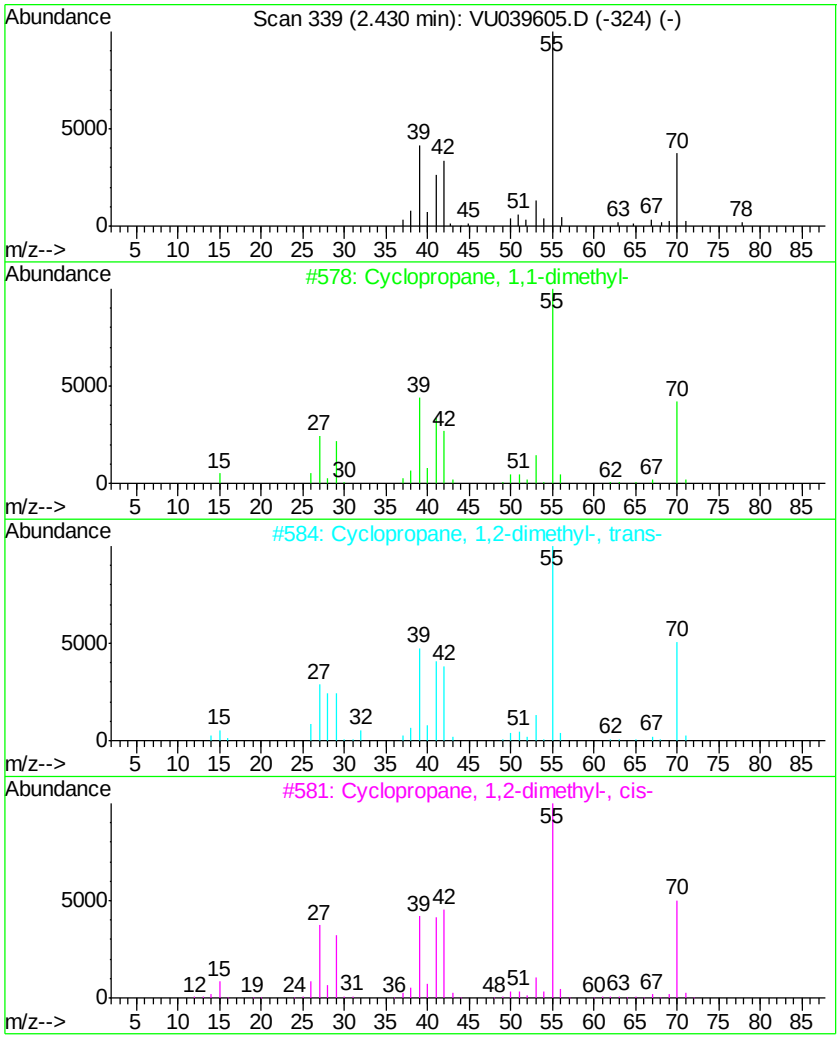
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

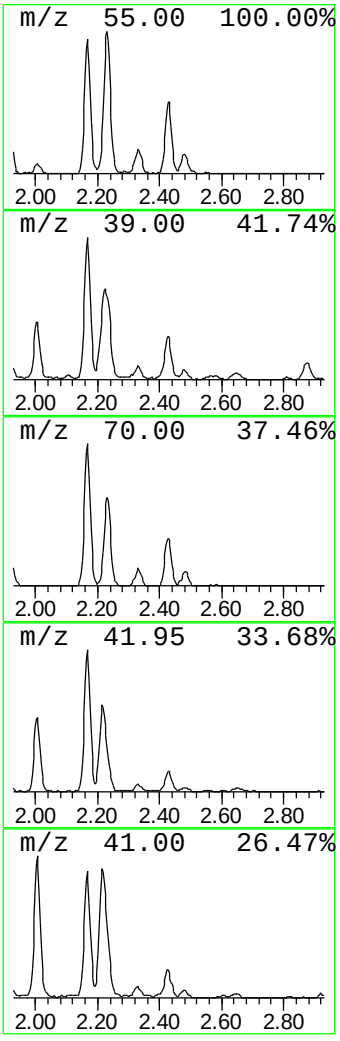
**********

Peak Number 7 (DEL) Alkane: Cyclic2.43 Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.43 | 0.64 ug/L | 39136 | 1,4-Difluorobenzene | 6.27 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclopropane, 1,1-dimethyl- | 70 | C5H10 | 001630-94-0 | 87 |
| 2 | Cyclopropane, 1,2-dimethyl-, trans- | 70 | C5H10 | 002402-06-4 | 87 |
| 3 | Cyclopropane, 1,2-dimethyl-, cis- | 70 | C5H10 | 000930-18-7 | 83 |
| 4 | Cyclopropane, 1,1-dimethyl- | 70 | C5H10 | 001630-94-0 | 83 |
| 5 | 2-Butene, 2-methyl- | 70 | C5H10 | 000513-35-9 | 80 |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR072120WMA.M
Quant Title : TRACE VOA SOM01.0

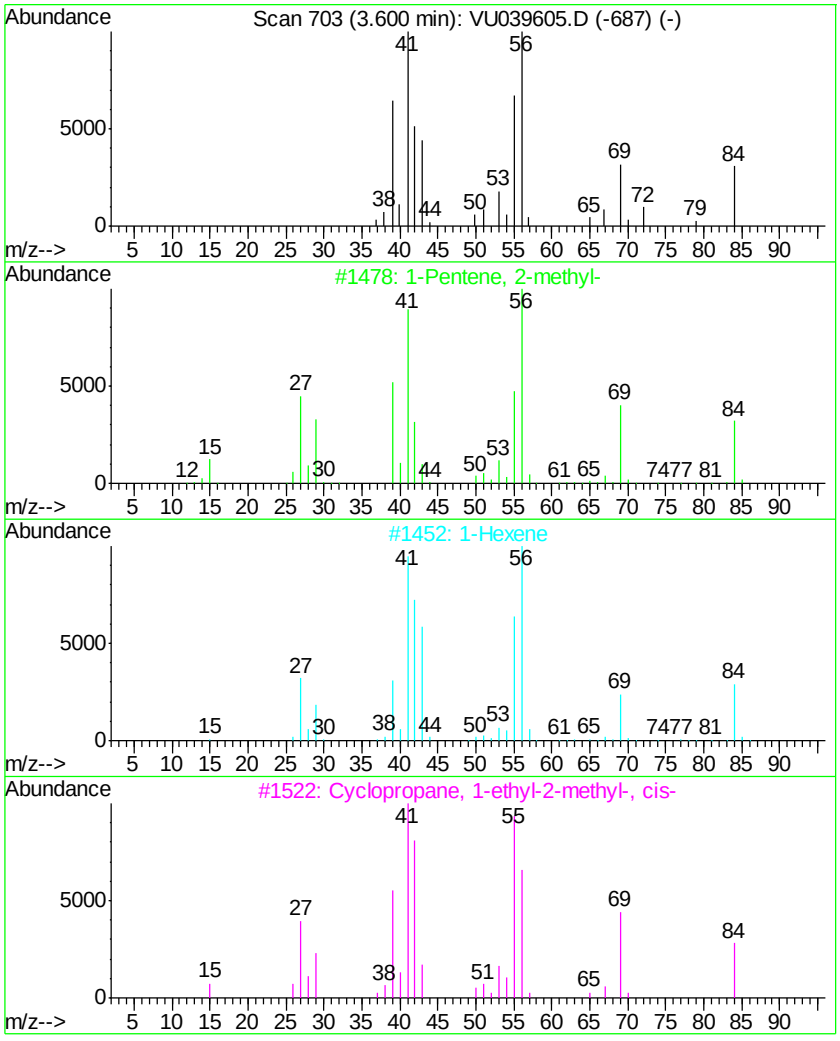
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

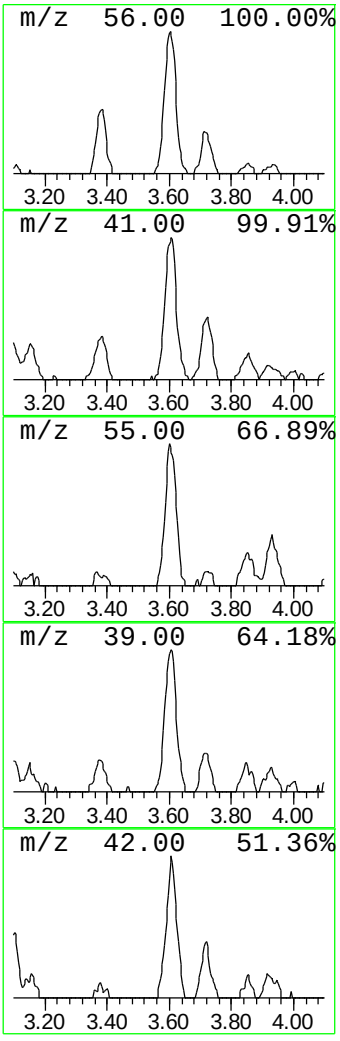
**********

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.60 | 1.08 ug/L | 66144 | 1,4-Difluorobenzene | 6.27 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Pentene, 2-methyl- | 84 | C6H12 | 000763-29-1 | 80 |
| 2 | 1-Hexene | 84 | C6H12 | 000592-41-6 | 58 |
| 3 | Cyclopropane, 1-ethyl-2-methyl-,... | 84 | C6H12 | 019781-68-1 | 53 |
| 4 | 3-Hexene, (E)- | 84 | C6H12 | 013269-52-8 | 52 |
| 5 | 1-Hexene | 84 | C6H12 | 000592-41-6 | 49 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072120\
Data File : VU039605.D
Acq On : 21 Jul 2020 18:29
Operator : SY/MD
Sample : L3347-08
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 1 Sample Multiplier: 1

**Instrument :** MSVOA_U
**ClientSampleId :** GAY02

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR072120WMA.M
Quant Title : TRACE VOA SOM01.0

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: Str... | 1.37 | 7.2 | ug/L | 443489 | 1 | 6.27 | 307497 | 5.0 |
| (DEL) Alkane: Str... | 1.50 | 0.8 | ug/L | 47617 | 1 | 6.27 | 307497 | 5.0 |
| (DEL) Alkane: Str... | 1.74 | 0.6 | ug/L | 38627 | 1 | 6.27 | 307497 | 5.0 |
| (DEL) Alkane: Str... | 2.01 | 1.2 | ug/L | 74348 | 1 | 6.27 | 307497 | 5.0 |
| (DEL) Alkane: Cyc... | 2.17 | 1.7 | ug/L | 103894 | 1 | 6.27 | 307497 | 5.0 |
| (DEL) Alkane: Str... | 2.22 | 2.1 | ug/L | 130669 | 1 | 6.27 | 307497 | 5.0 |
| (DEL) Alkane: Cyc... | 2.43 | 0.6 | ug/L | 39136 | 1 | 6.27 | 307497 | 5.0 |
| (DEL) Alkane: Str... | 3.60 | 1.1 | ug/L | 66144 | 1 | 6.27 | 307497 | 5.0 |