

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
 Data File : VU037747.D  
 Acq On : 16 Apr 2020 12:07  
 Operator : JC/MD  
 Sample : L2296-03  
 Misc : 25mL/MSVOA U/WATER  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BFS89

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 1.617       | 76            | 86          | 98           | rBV2     | 817967         | 930949        | 3.18%           | 0.569%        |
| 2         | 2.594       | 376           | 390         | 420          | rVB3     | 236380         | 502584        | 1.71%           | 0.307%        |
| 3         | 3.073       | 522           | 539         | 541          | rBV      | 1242794        | 2043744       | 6.97%           | 1.250%        |
| 4         | 3.112       | 542           | 551         | 597          | rVB      | 4147140        | 8677852       | 29.61%          | 5.307%        |
| 5         | 3.395       | 617           | 639         | 669          | rBV      | 11179799       | 24654132      | 84.12%          | 15.079%       |
| 6         | 3.735       | 725           | 745         | 770          | rBV      | 7202849        | 15898078      | 54.25%          | 9.723%        |
| 7         | 3.909       | 773           | 799         | 822          | rBV2     | 374872         | 1389679       | 4.74%           | 0.850%        |
| 8         | 4.018       | 822           | 833         | 851          | rVB      | 245698         | 590851        | 2.02%           | 0.361%        |
| 9         | 4.131       | 851           | 868         | 882          | rBV      | 215783         | 539276        | 1.84%           | 0.330%        |
| 10        | 4.337       | 912           | 932         | 938          | rBV      | 366936         | 978445        | 3.34%           | 0.598%        |
| 11        | 4.472       | 960           | 974         | 986          | rBV2     | 316066         | 749838        | 2.56%           | 0.459%        |
| 12        | 4.571       | 986           | 1005        | 1022         | rVV      | 12339706       | 29307682      | 100.00%         | 17.925%       |
| 13        | 4.706       | 1023          | 1047        | 1075         | rVB2     | 1526916        | 4191179       | 14.30%          | 2.563%        |
| 14        | 5.102       | 1156          | 1170        | 1183         | rBV      | 164108         | 368004        | 1.26%           | 0.225%        |
| 15        | 5.353       | 1230          | 1248        | 1259         | rBV      | 1535719        | 3422828       | 11.68%          | 2.093%        |
| 16        | 5.423       | 1259          | 1270        | 1287         | rVB      | 1313327        | 2902448       | 9.90%           | 1.775%        |
| 17        | 5.761       | 1355          | 1375        | 1403         | rVB2     | 356227         | 929700        | 3.17%           | 0.569%        |
| 18        | 6.282       | 1522          | 1537        | 1553         | rBV      | 340889         | 632888        | 2.16%           | 0.387%        |
| 19        | 6.722       | 1660          | 1674        | 1689         | rBV      | 228060         | 442419        | 1.51%           | 0.271%        |
| 20        | 7.925       | 2026          | 2048        | 2057         | rBV      | 457488         | 790302        | 2.70%           | 0.483%        |
| 21        | 7.996       | 2057          | 2070        | 2084         | rVB      | 5355022        | 8885044       | 30.32%          | 5.434%        |
| 22        | 8.658       | 2260          | 2276        | 2314         | rBV      | 721418         | 1228784       | 4.19%           | 0.752%        |
| 23        | 9.468       | 2506          | 2528        | 2544         | rBV2     | 1462169        | 3150649       | 10.75%          | 1.927%        |
| 24        | 9.591       | 2553          | 2566        | 2586         | rVB      | 1186173        | 1878463       | 6.41%           | 1.149%        |
| 25        | 9.713       | 2590          | 2604        | 2634         | rVB      | 2004312        | 3273253       | 11.17%          | 2.002%        |
| 26        | 10.118      | 2717          | 2730        | 2756         | rVB      | 1105399        | 1747961       | 5.96%           | 1.069%        |
| 27        | 10.504      | 2834          | 2850        | 2862         | rBV      | 409424         | 643769        | 2.20%           | 0.394%        |
| 28        | 10.922      | 2960          | 2980        | 2996         | rBV      | 1247690        | 1906832       | 6.51%           | 1.166%        |
| 29        | 11.015      | 2996          | 3009        | 3014         | rBV      | 2754477        | 4527336       | 15.45%          | 2.769%        |
| 30        | 11.105      | 3027          | 3037        | 3054         | rVB      | 1536066        | 2343587       | 8.00%           | 1.433%        |
| 31        | 11.250      | 3064          | 3082        | 3091         | rBV      | 5905805        | 8721698       | 29.76%          | 5.334%        |
| 32        | 11.307      | 3091          | 3100        | 3124         | rVB      | 2185085        | 3411055       | 11.64%          | 2.086%        |
| 33        | 11.484      | 3132          | 3155        | 3171         | rBV      | 6082786        | 9224980       | 31.48%          | 5.642%        |
| 34        | 11.574      | 3171          | 3183        | 3193         | rBV      | 494783         | 761495        | 2.60%           | 0.466%        |

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Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

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\SOMUTR041420WMA.M  
Title : TRACE VOA SOM01.0

|    |        |      |      |      |      |         |         |       |        |
|----|--------|------|------|------|------|---------|---------|-------|--------|
| 35 | 11.761 | 3229 | 3241 | 3251 | rBV  | 461077  | 721893  | 2.46% | 0.442% |
| 36 | 11.851 | 3251 | 3269 | 3278 | rVV2 | 762769  | 1971255 | 6.73% | 1.206% |
| 37 | 11.909 | 3278 | 3287 | 3300 | rVB  | 1639715 | 2507759 | 8.56% | 1.534% |
| 38 | 12.108 | 3336 | 3349 | 3367 | rBV  | 1057556 | 1801588 | 6.15% | 1.102% |
| 39 | 12.208 | 3367 | 3380 | 3399 | rVB5 | 668133  | 1482461 | 5.06% | 0.907% |
| 40 | 12.590 | 3486 | 3499 | 3508 | rBV  | 326202  | 503085  | 1.72% | 0.308% |
| 41 | 12.996 | 3608 | 3625 | 3633 | rVV  | 207562  | 324353  | 1.11% | 0.198% |
| 42 | 13.050 | 3633 | 3642 | 3655 | rVB  | 361562  | 557218  | 1.90% | 0.341% |
| 43 | 13.481 | 3765 | 3776 | 3791 | rVB3 | 340389  | 585134  | 2.00% | 0.358% |
| 44 | 13.873 | 3886 | 3898 | 3914 | rBV2 | 293204  | 487565  | 1.66% | 0.298% |
| 45 | 14.121 | 3960 | 3975 | 3996 | rBV  | 558220  | 914226  | 3.12% | 0.559% |

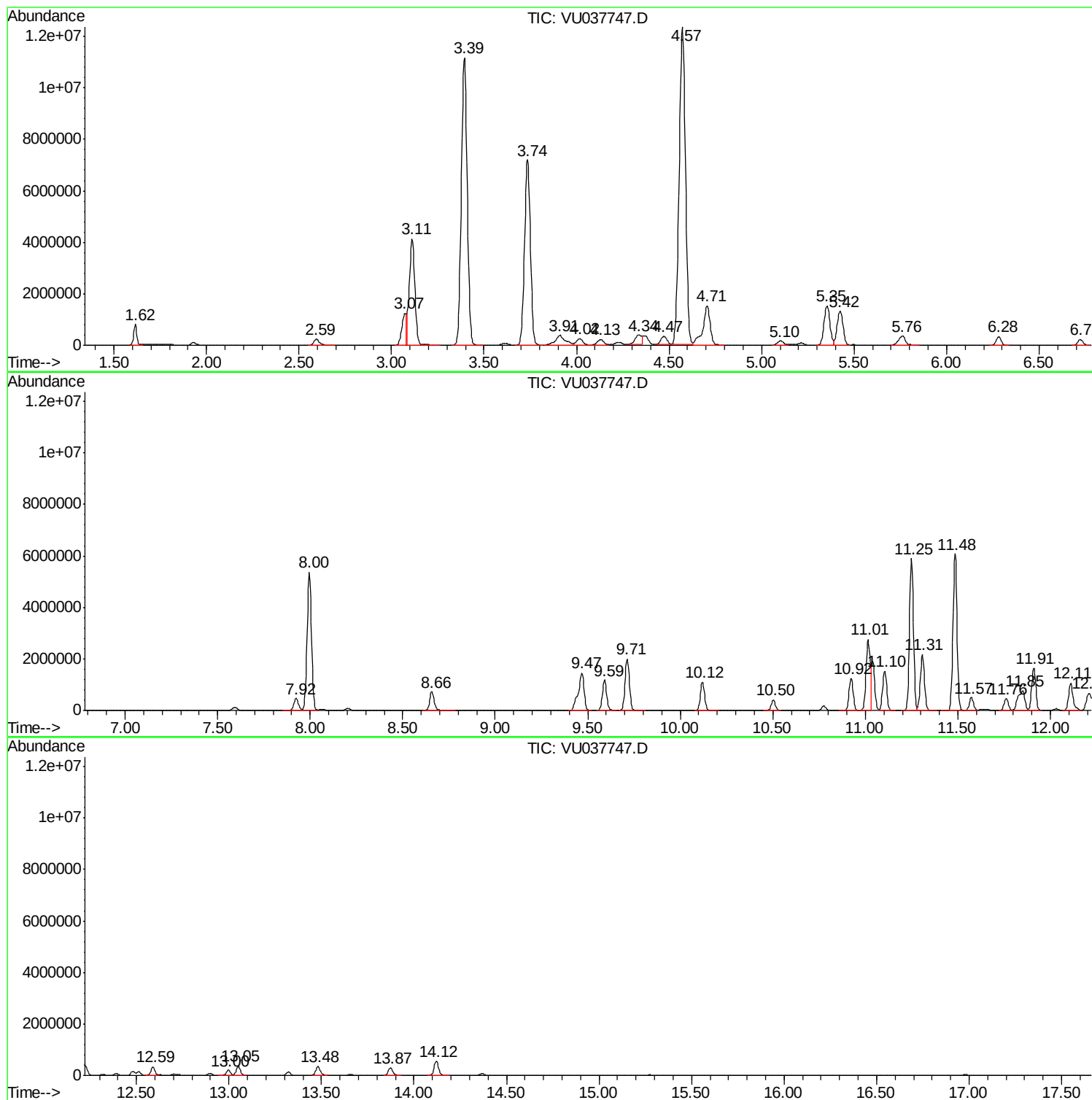
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Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

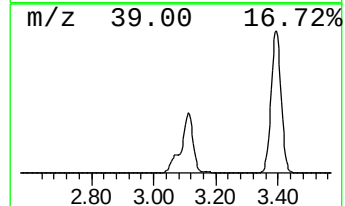
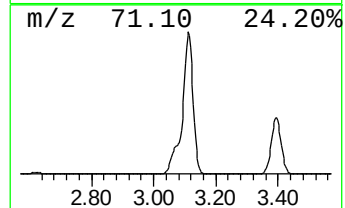
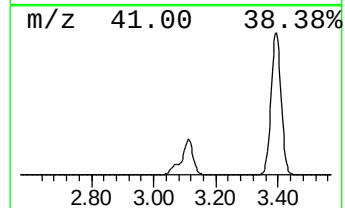
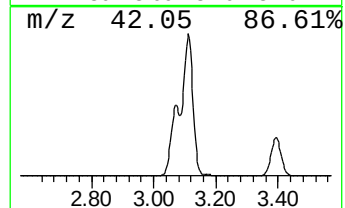
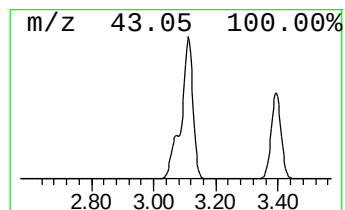
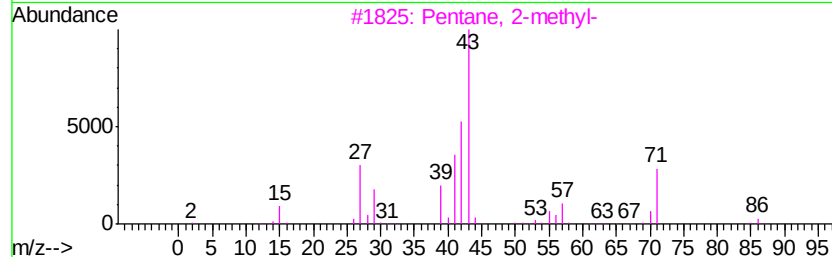
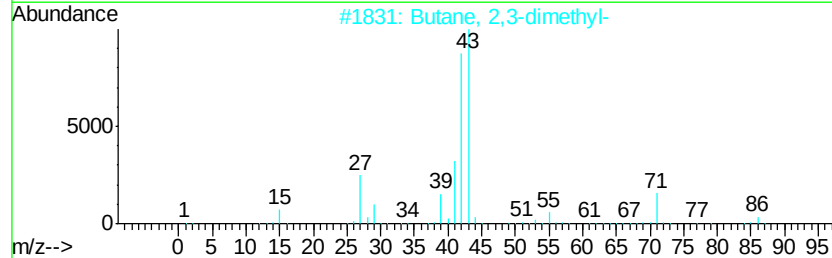
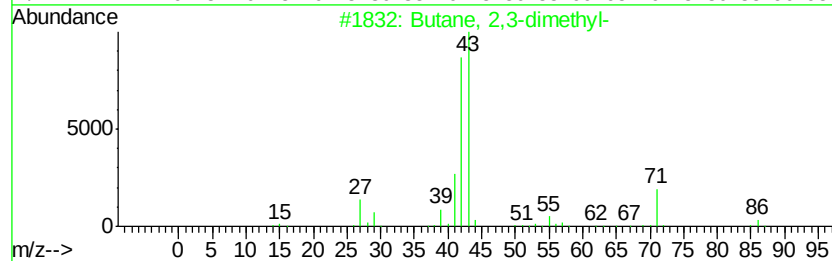
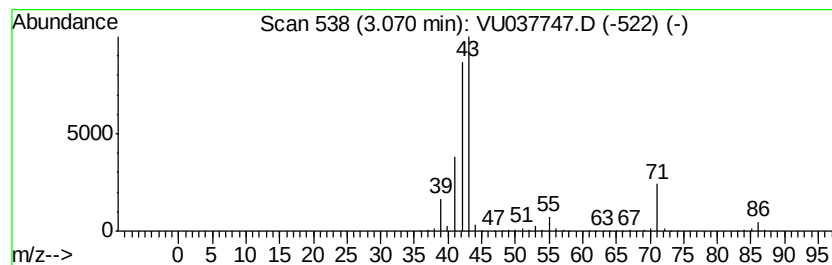
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T. | EstConc    | Area    | Relative to ISTD    | R.T. |
|------|------------|---------|---------------------|------|
| 3.07 | 16.15 ug/L | 2043740 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID          | MW | MolForm | CAS#        | Qual |
|------|----|-----------------------|----|---------|-------------|------|
| 1    | 5  | Butane, 2,3-dimethyl- | 86 | C6H14   | 000079-29-8 | 91   |
| 2    |    | Butane, 2,3-dimethyl- | 86 | C6H14   | 000079-29-8 | 90   |
| 3    |    | Pentane, 2-methyl-    | 86 | C6H14   | 000107-83-5 | 50   |
| 4    |    | Pentane, 2-methyl-    | 86 | C6H14   | 000107-83-5 | 50   |
| 5    |    | Butane, 2,3-dimethyl- | 86 | C6H14   | 000079-29-8 | 49   |



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Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

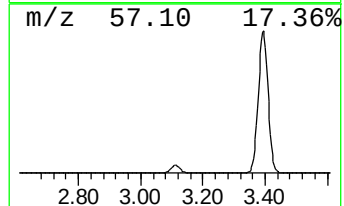
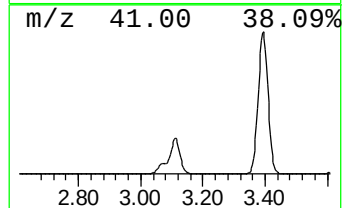
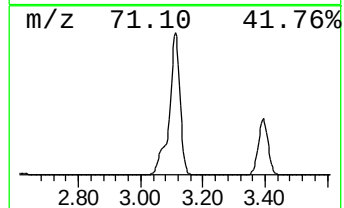
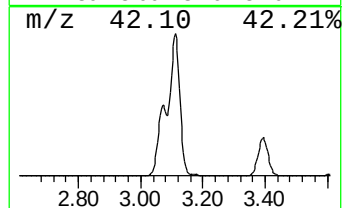
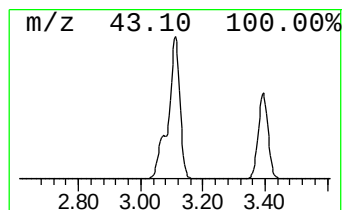
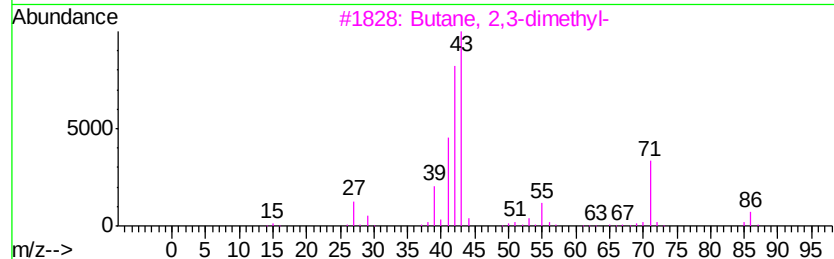
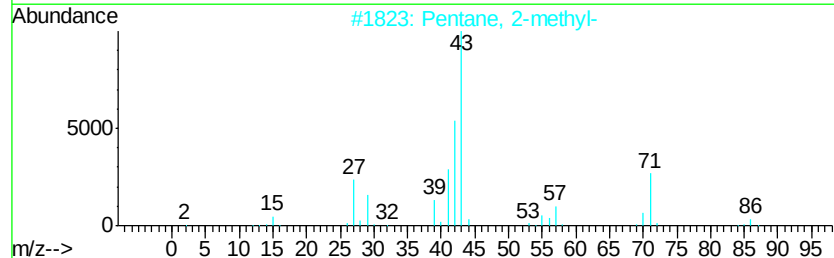
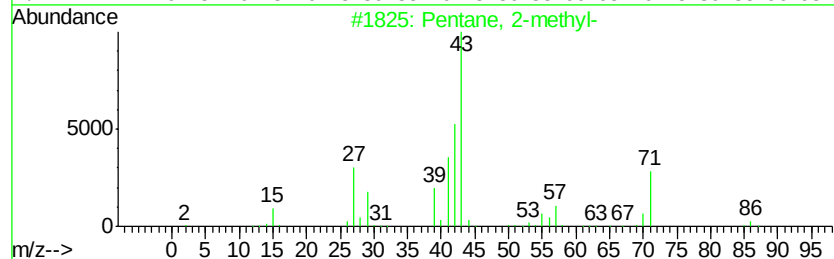
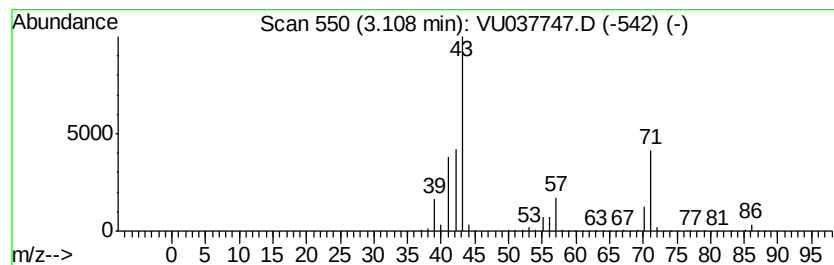
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T. | EstConc    | Area    | Relative to ISTD    | R.T. |
|------|------------|---------|---------------------|------|
| 3.11 | 68.56 ug/L | 8677850 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | 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 | 47   |
| 4    |    | Hexane, 2,3-dimethyl-  | 114 | C8H18   | 000584-94-1 | 28   |
| 5    |    | Pentane, 3,3-dimethyl- | 100 | C7H16   | 000562-49-2 | 23   |



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Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

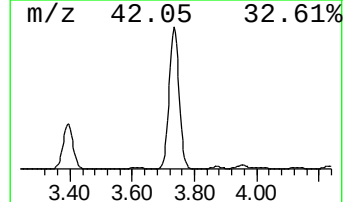
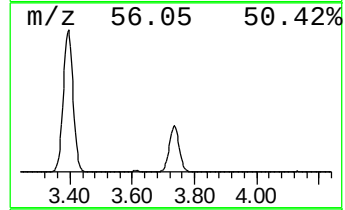
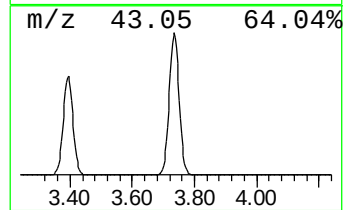
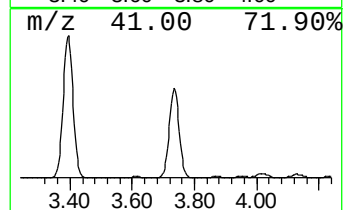
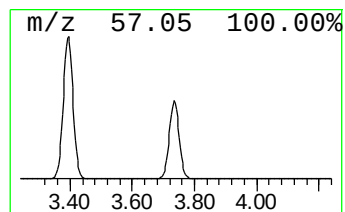
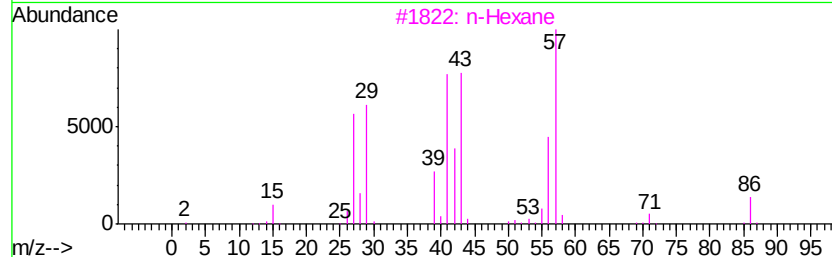
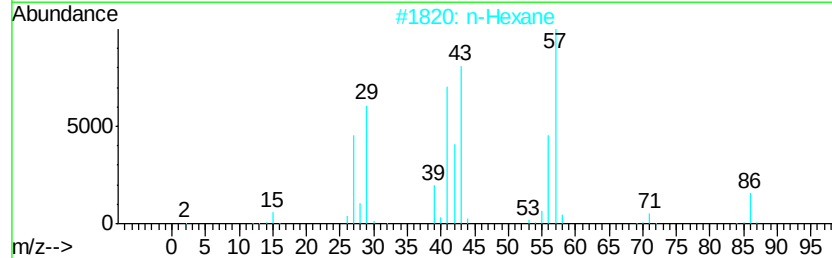
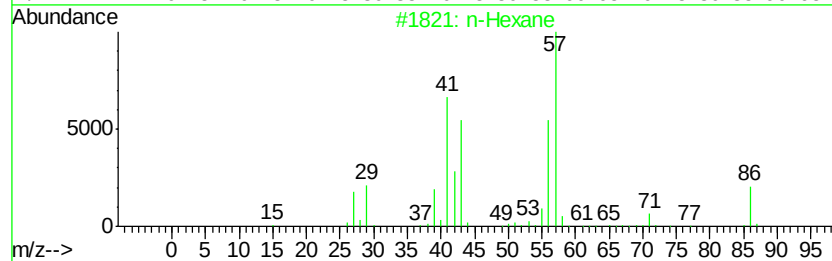
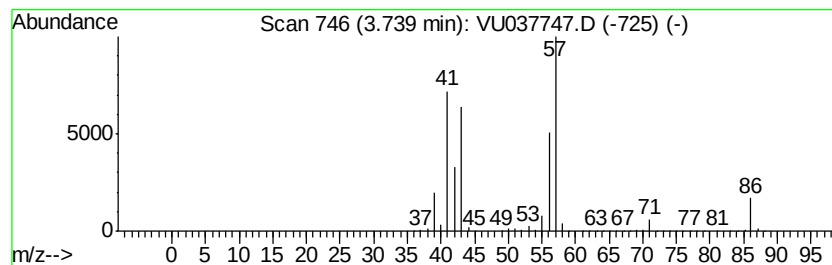
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T. | EstConc     | Area     | Relative to ISTD    | R.T. |
|------|-------------|----------|---------------------|------|
| 3.74 | 125.60 ug/L | 15898100 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | n-Hexane                      | 86  | C6H14   | 000110-54-3 | 94   |
| 2    |    | n-Hexane                      | 86  | C6H14   | 000110-54-3 | 91   |
| 3    |    | n-Hexane                      | 86  | C6H14   | 000110-54-3 | 91   |
| 4    |    | Pentane, 2,2,3,4-tetramethyl- | 128 | C9H20   | 001186-53-4 | 42   |
| 5    |    | Methane, isocyanato-          | 57  | C2H3NO  | 000624-83-9 | 37   |



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Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

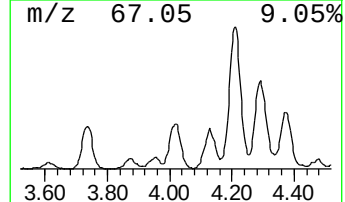
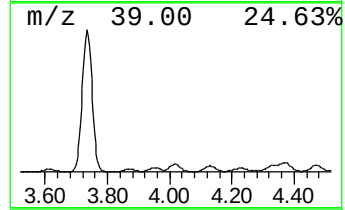
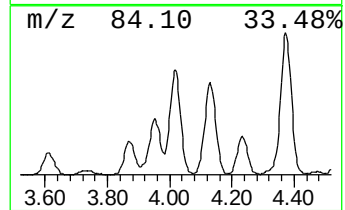
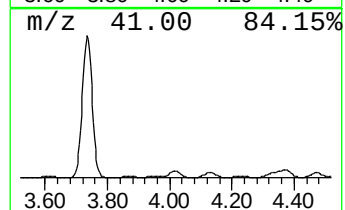
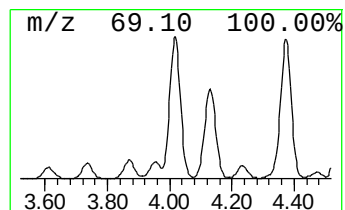
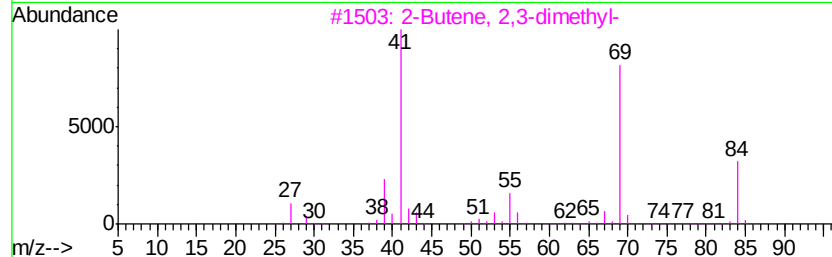
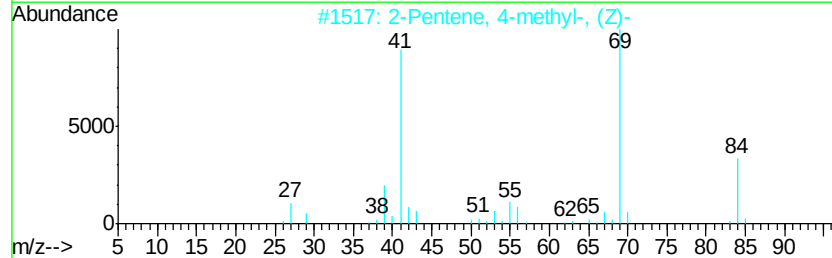
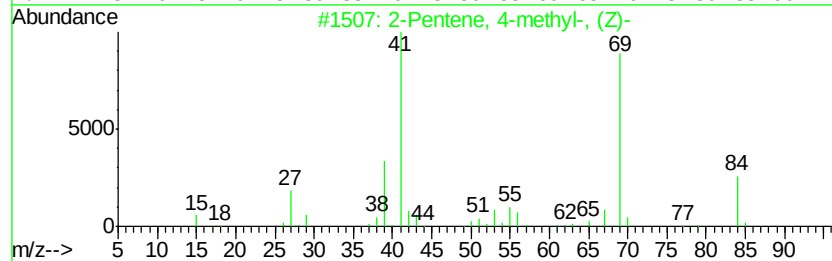
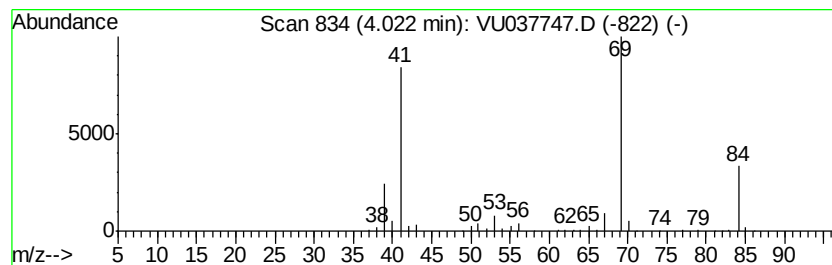
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 4.02 | 4.67 ug/L | 590851 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID               | MW | MolForm | CAS#        | Qual |
|------|----|----------------------------|----|---------|-------------|------|
| 1    | 5  | 2-Pentene, 4-methyl-, (Z)- | 84 | C6H12   | 000691-38-3 | 90   |
| 2    |    | 2-Pentene, 4-methyl-, (Z)- | 84 | C6H12   | 000691-38-3 | 78   |
| 3    |    | 2-Butene, 2,3-dimethyl-    | 84 | C6H12   | 000563-79-1 | 78   |
| 4    |    | 2-Butene, 2,3-dimethyl-    | 84 | C6H12   | 000563-79-1 | 78   |
| 5    |    | 2-Pentene, 2-methyl-       | 84 | C6H12   | 000625-27-4 | 78   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

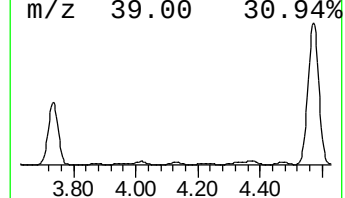
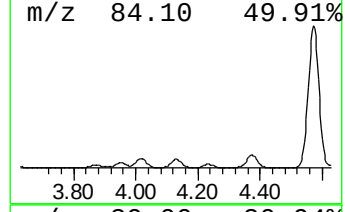
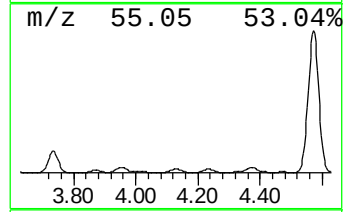
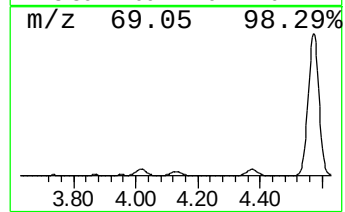
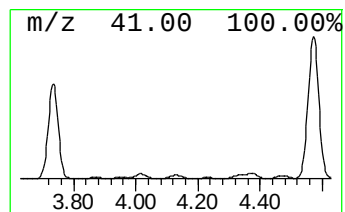
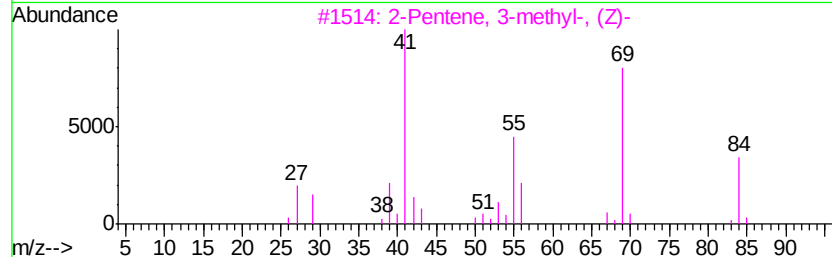
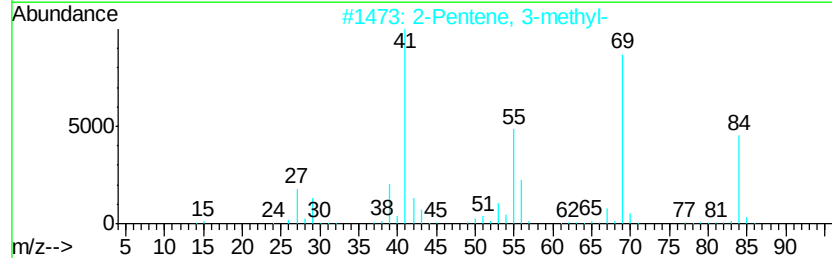
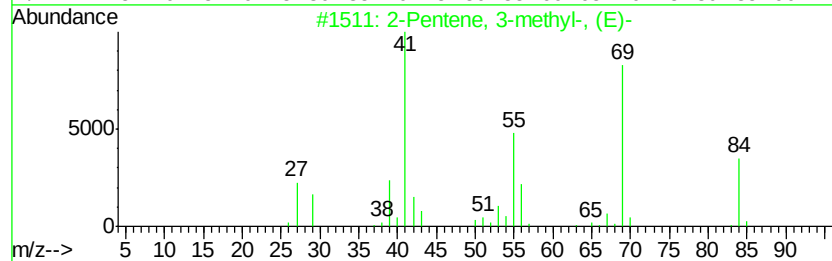
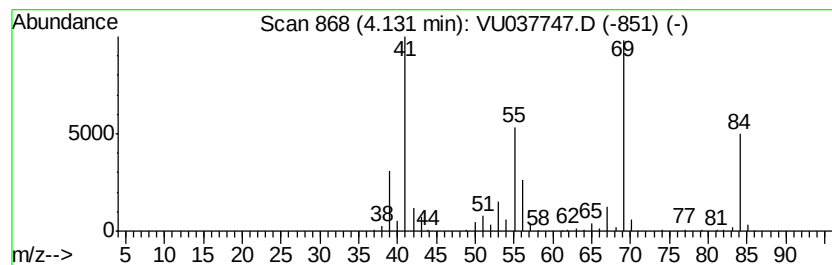
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 4.13 | 4.26 ug/L | 539276 | 1,4-Difluorobenzene | 6.28 |

| Hit# of 5 | Tentative ID               | MW | MolForm | CAS#        | Qual |
|-----------|----------------------------|----|---------|-------------|------|
| 1         | 2-Pentene, 3-methyl-, (E)- | 84 | C6H12   | 000616-12-6 | 94   |
| 2         | 2-Pentene, 3-methyl-       | 84 | C6H12   | 000922-61-2 | 94   |
| 3         | 2-Pentene, 3-methyl-, (Z)- | 84 | C6H12   | 000922-62-3 | 91   |
| 4         | 2-Pentene, 3-methyl-, (Z)- | 84 | C6H12   | 000922-62-3 | 91   |
| 5         | 2-Pentene, 3-methyl-, (E)- | 84 | C6H12   | 000616-12-6 | 91   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

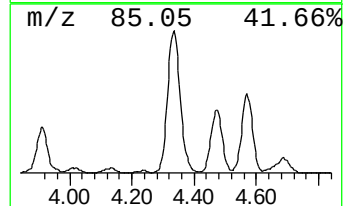
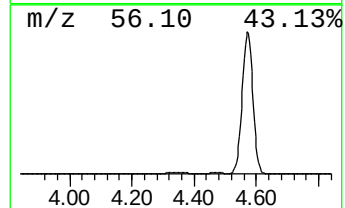
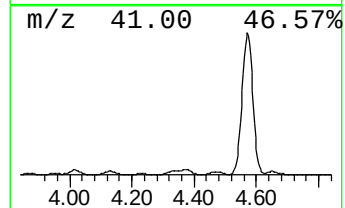
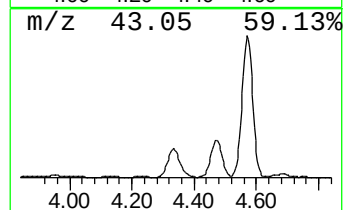
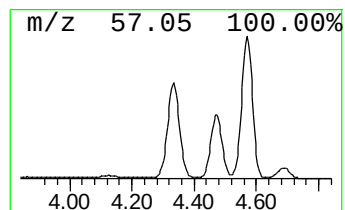
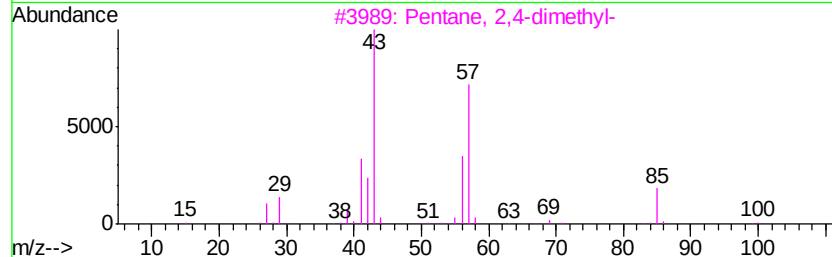
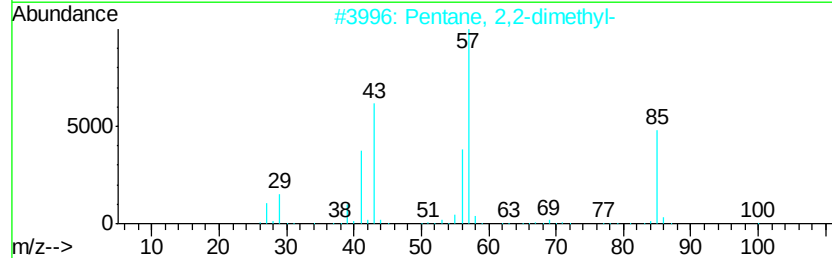
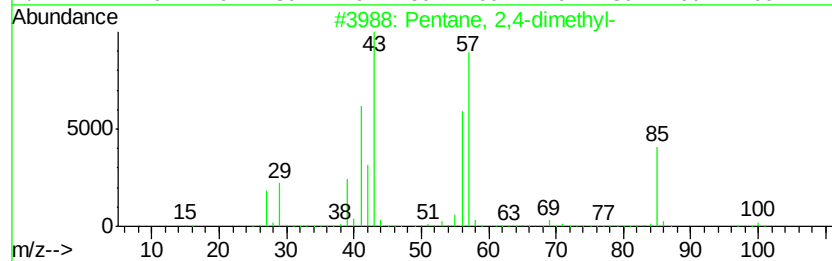
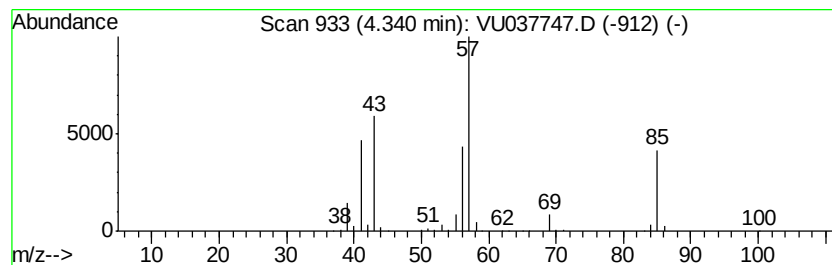
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 4.34 | 7.73 ug/L | 978445 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentane, 2,4-dimethyl-   | 100 | C7H16   | 000108-08-7 | 83   |
| 2    |    | Pentane, 2,2-dimethyl-   | 100 | C7H16   | 000590-35-2 | 83   |
| 3    |    | Pentane, 2,4-dimethyl-   | 100 | C7H16   | 000108-08-7 | 78   |
| 4    |    | Pentane, 2,2-dimethyl-   | 100 | C7H16   | 000590-35-2 | 78   |
| 5    |    | Butane, 2,2,3-trimethyl- | 100 | C7H16   | 000464-06-2 | 78   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

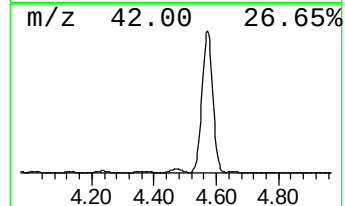
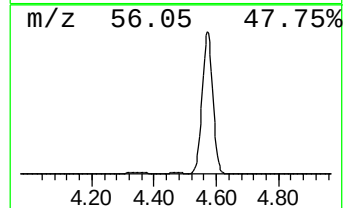
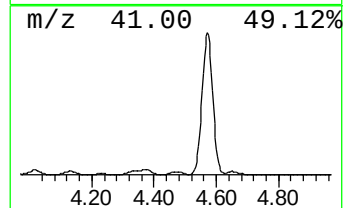
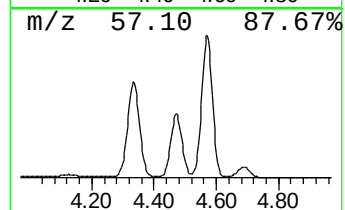
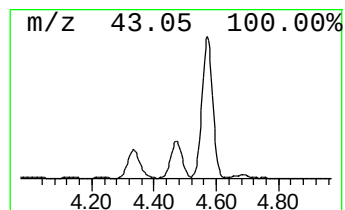
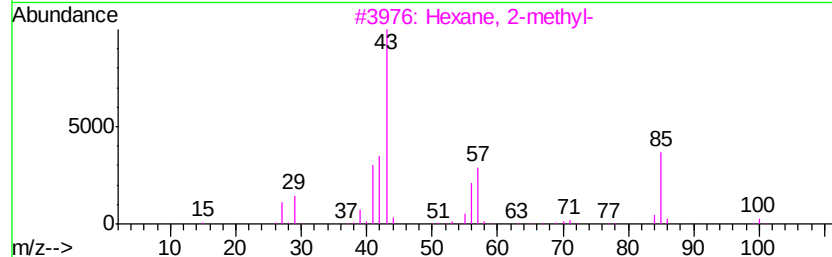
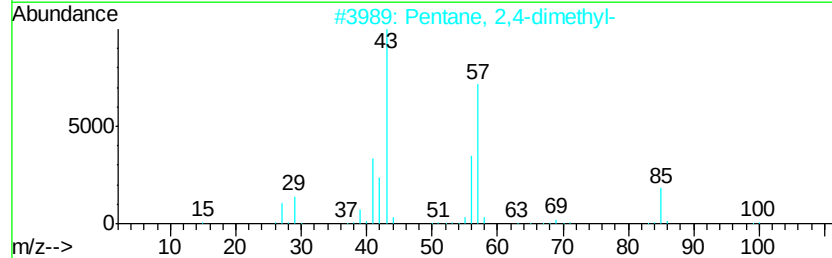
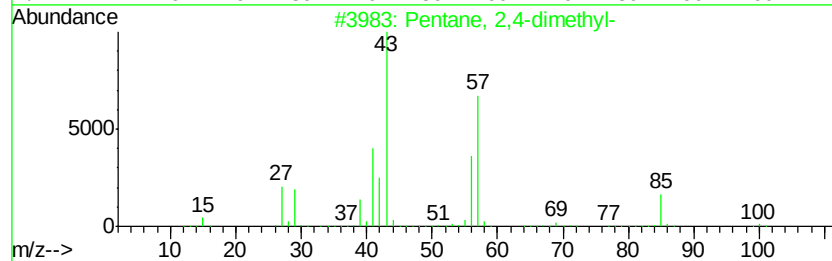
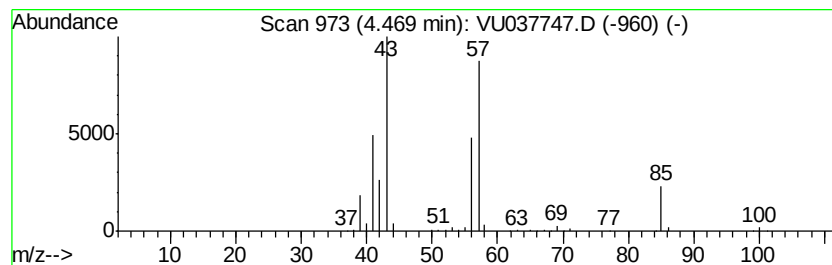
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 4.47 | 5.92 ug/L | 749838 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentane, 2,4-dimethyl-   | 100 | C7H16   | 000108-08-7 | 91   |
| 2    |    |   | Pentane, 2,4-dimethyl-   | 100 | C7H16   | 000108-08-7 | 83   |
| 3    |    |   | Hexane, 2-methyl-        | 100 | C7H16   | 000591-76-4 | 72   |
| 4    |    |   | Pentane, 2,4-dimethyl-   | 100 | C7H16   | 000108-08-7 | 72   |
| 5    |    |   | Butane, 2,2,3-trimethyl- | 100 | C7H16   | 000464-06-2 | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

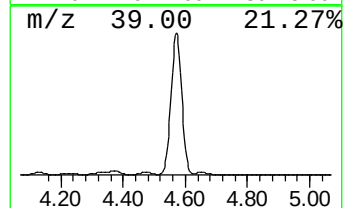
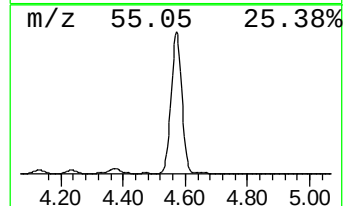
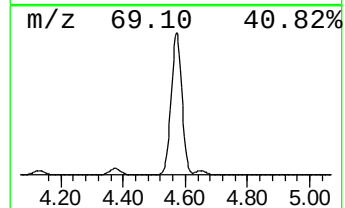
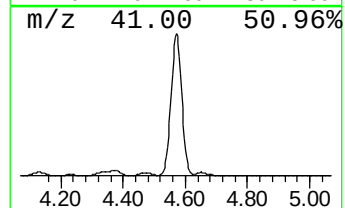
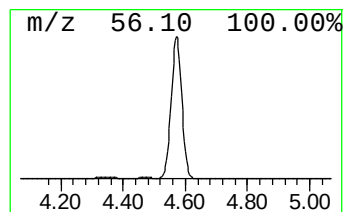
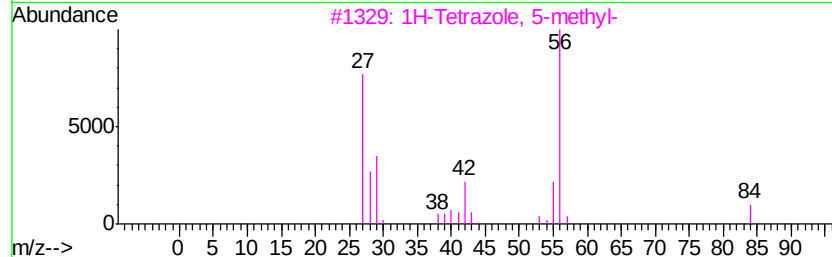
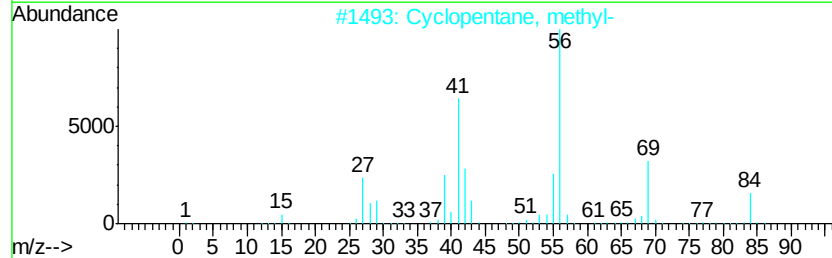
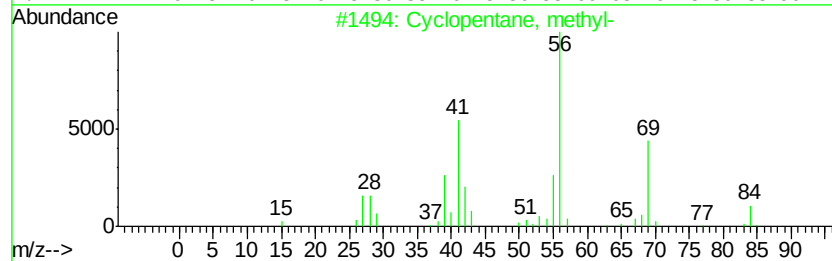
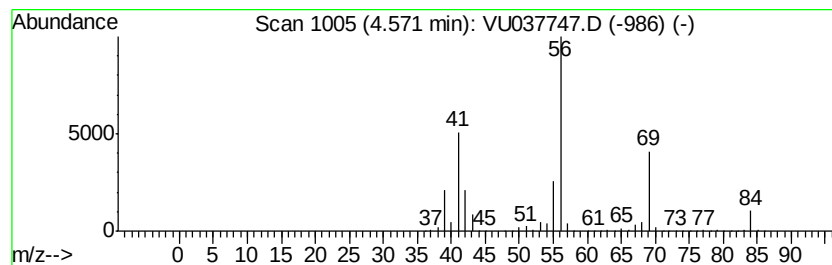
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 (DEL) Alkane: Cyclic4.57 Concentration Rank 1

| R.T. | EstConc     | Area     | Relative to ISTD    | R.T. |
|------|-------------|----------|---------------------|------|
| 4.57 | 231.54 ug/L | 29307700 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID            | MW | MolForm | CAS#        | Qual |
|------|----|-------------------------|----|---------|-------------|------|
| 1    | 5  | Cyclopentane, methyl-   | 84 | C6H12   | 000096-37-7 | 94   |
| 2    |    | Cyclopentane, methyl-   | 84 | C6H12   | 000096-37-7 | 91   |
| 3    |    | 1H-Tetrazole, 5-methyl- | 84 | C2H4N4  | 004076-36-2 | 80   |
| 4    |    | Cyclopropane, propyl-   | 84 | C6H12   | 002415-72-7 | 78   |
| 5    |    | Cyclopentane, methyl-   | 84 | C6H12   | 000096-37-7 | 78   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

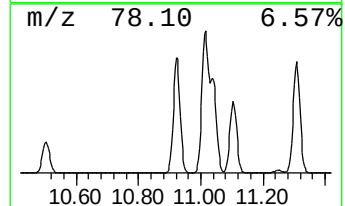
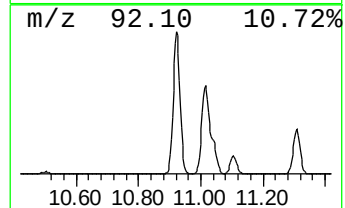
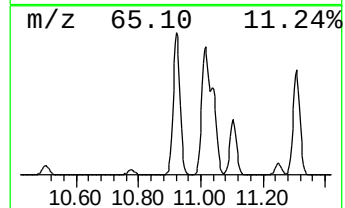
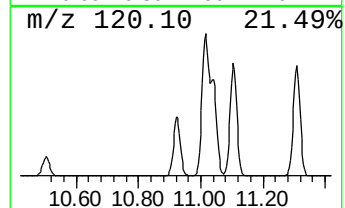
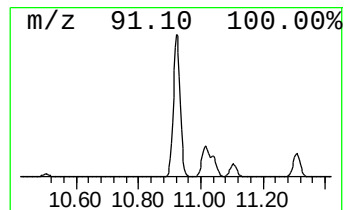
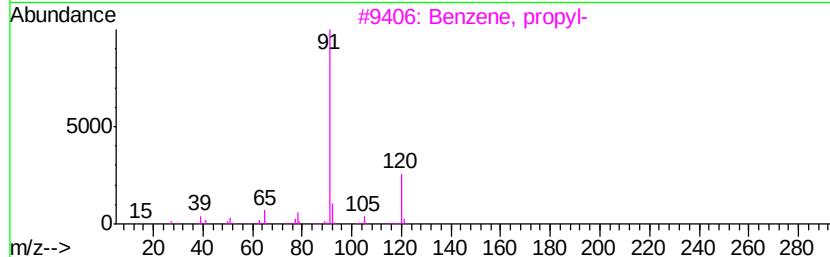
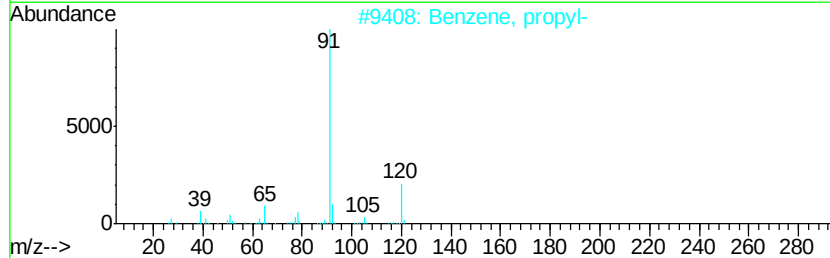
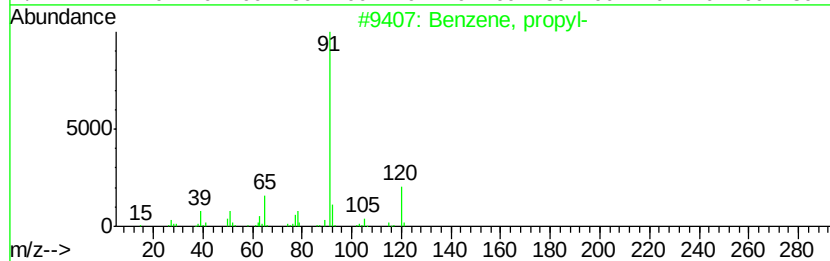
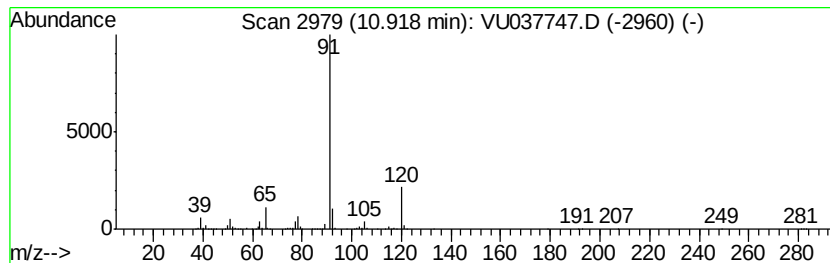
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Benzene, propyl- Concentration Rank 13

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 10.92 | 4.84 ug/L | 1906830 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Benzene, propyl-                    | 120 | C9H12    | 000103-65-1 | 91   |
| 2    |    | Benzene, propyl-                    | 120 | C9H12    | 000103-65-1 | 91   |
| 3    |    | Benzene, propyl-                    | 120 | C9H12    | 000103-65-1 | 91   |
| 4    |    | 1,2-Ethanediamine, N,N'-bis(phen... | 240 | C16H20N2 | 000140-28-3 | 64   |
| 5    |    | N-Benzyl-2-phenethylamine           | 211 | C15H17N  | 003647-71-0 | 56   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

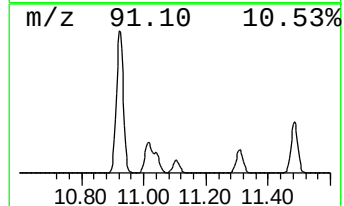
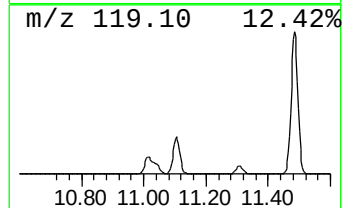
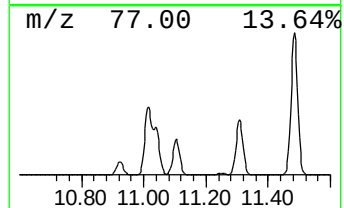
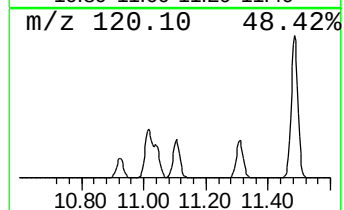
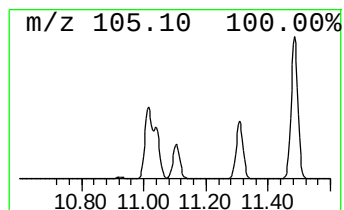
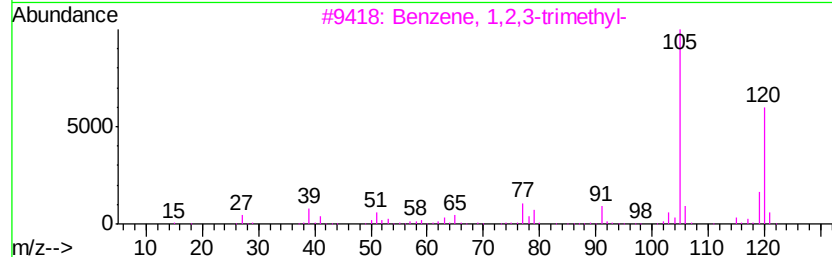
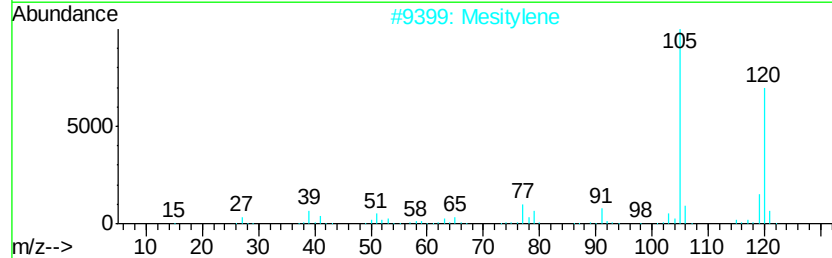
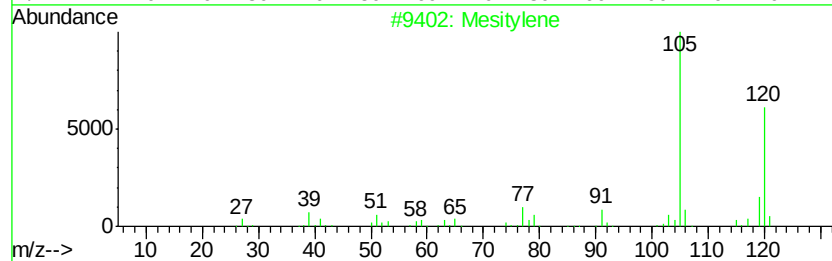
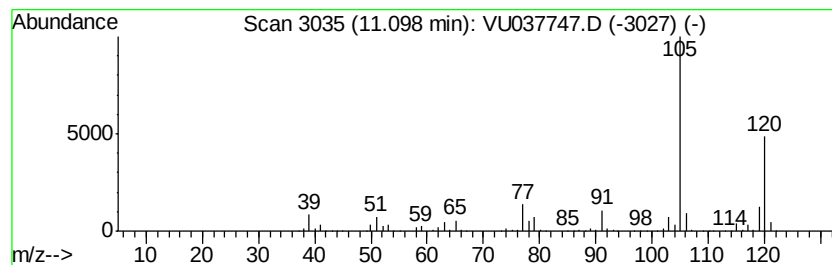
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Mesitylene Concentration Rank 11

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 11.10 | 5.94 ug/L | 2343590 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Mesitylene                | 120 | C9H12   | 000108-67-8 | 97   |
| 2    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 97   |
| 3    |    | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 97   |
| 4    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 95   |
| 5    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

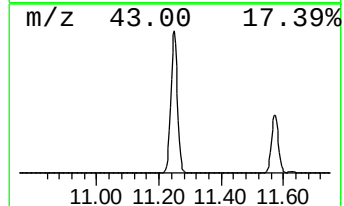
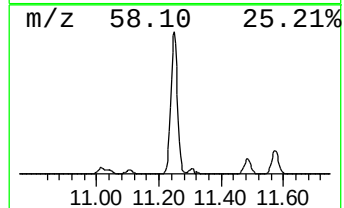
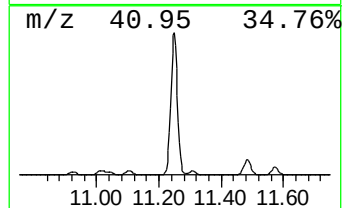
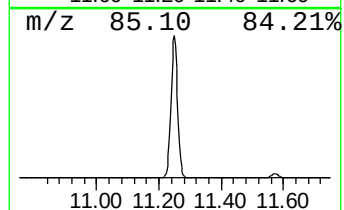
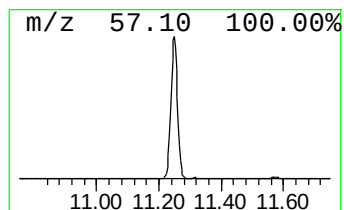
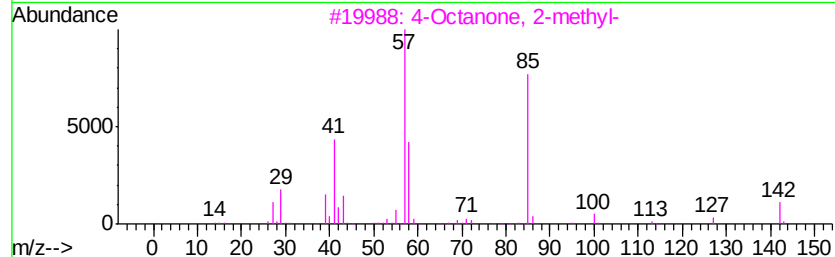
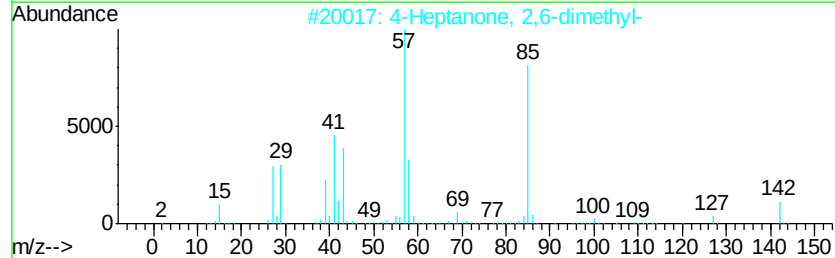
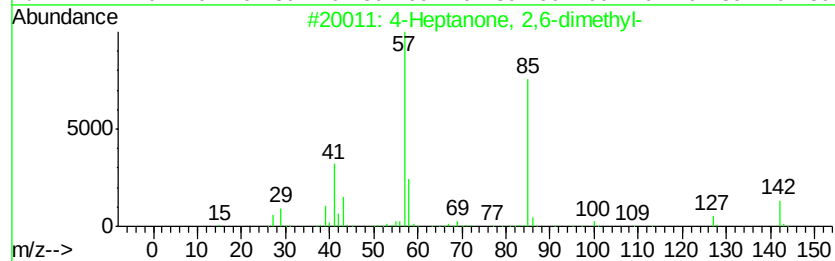
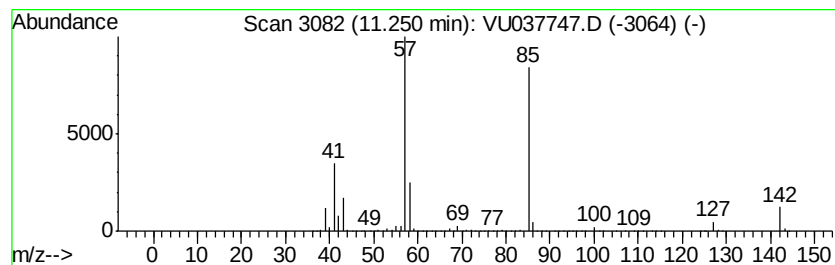
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 4-Heptanone, 2,6-dimethyl- Concentration Rank 5

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 11.25 | 22.12 ug/L | 8721700 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | 4-Heptanone, 2,6-dimethyl-     | 142 | C9H18O  | 000108-83-8 | 94   |
| 2    |    | 4-Heptanone, 2,6-dimethyl-     | 142 | C9H18O  | 000108-83-8 | 72   |
| 3    |    | 4-Octanone, 2-methyl-          | 142 | C9H18O  | 007492-38-8 | 64   |
| 4    |    | 5-Nonanone                     | 142 | C9H18O  | 000502-56-7 | 64   |
| 5    |    | 2,4-Hexanedione, 5,5-dimethyl- | 142 | C8H14O2 | 007307-04-2 | 59   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

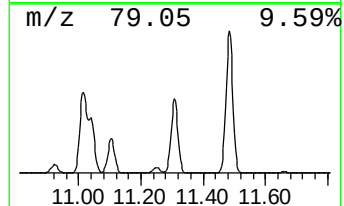
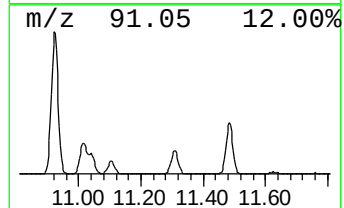
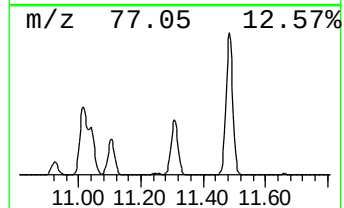
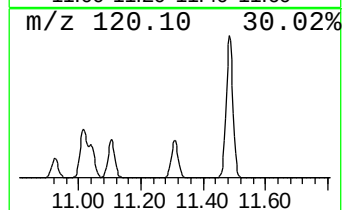
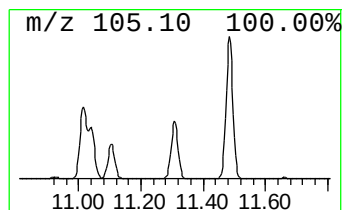
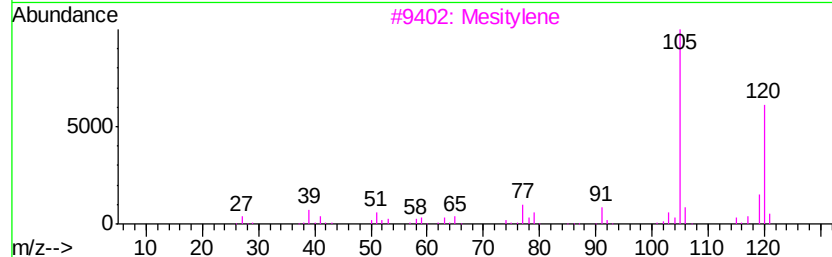
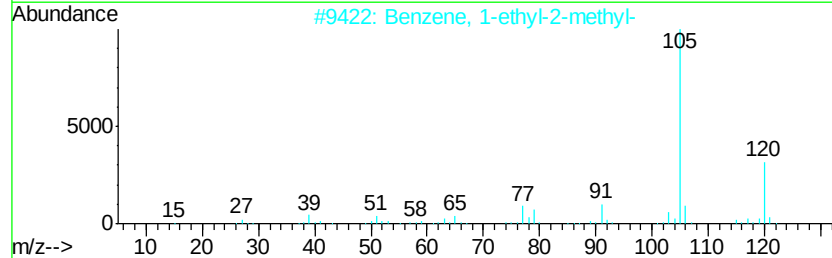
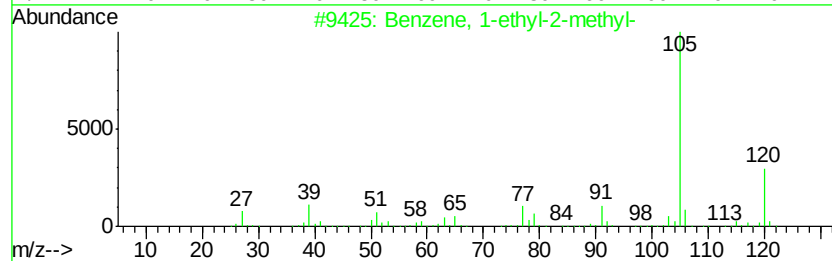
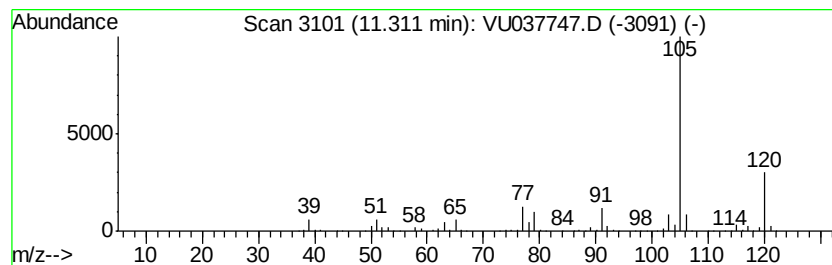
Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 13 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

| R.T.  | EstConc   | Area    | Relative to ISTD           |     | R.T.    |             |      |
|-------|-----------|---------|----------------------------|-----|---------|-------------|------|
| 11.31 | 8.65 ug/L | 3411060 | 1,4-Dichlorobenzene-d4     |     | 11.83   |             |      |
| Hit#  | of        | 5       | Tentative ID               | MW  | MolForm | CAS#        | Qual |
| 1     |           |         | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 2     |           |         | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 3     |           |         | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |
| 4     |           |         | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |
| 5     |           |         | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

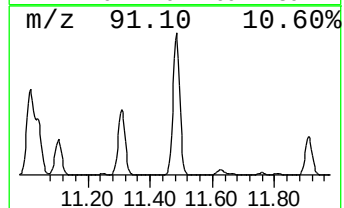
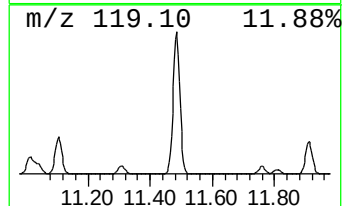
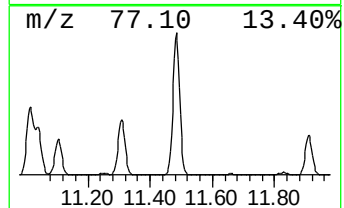
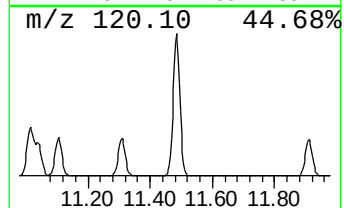
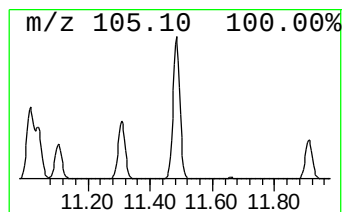
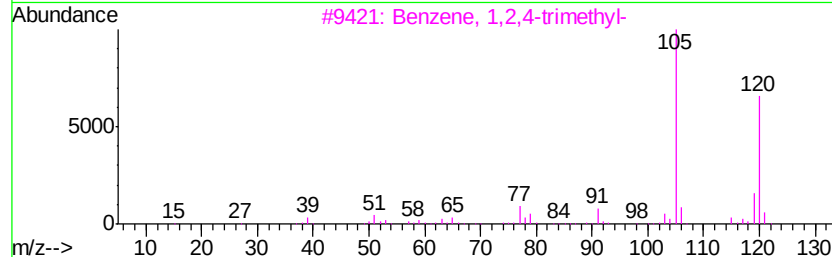
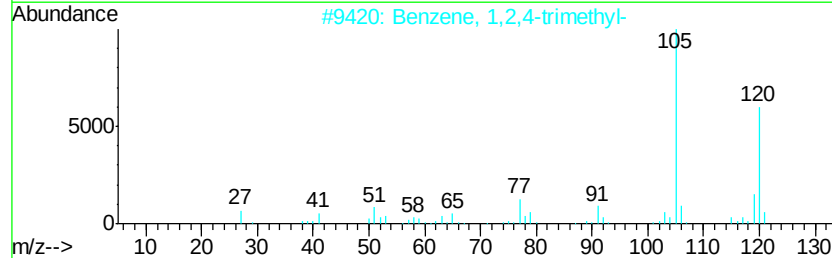
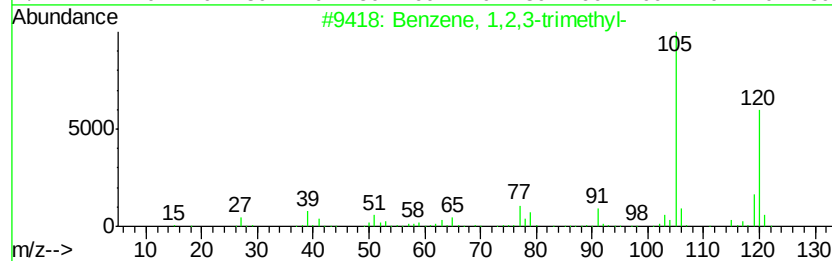
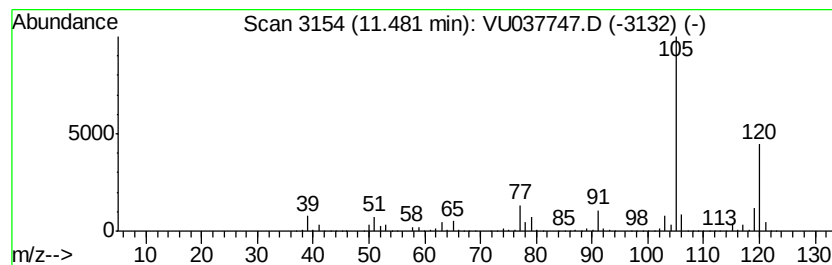
Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

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

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Peak Number 14 Benzene, 1,2,3-trimethyl- Concentration Rank 4

| R.T.    | EstConc    | Area                      | Relative to ISTD       | R.T.    |             |      |
|---------|------------|---------------------------|------------------------|---------|-------------|------|
| 11.48   | 23.40 ug/L | 9224980                   | 1,4-Dichlorobenzene-d4 | 11.83   |             |      |
| Hit# of | 5          | Tentative ID              | MW                     | MolForm | CAS#        | Qual |
| 1       |            | Benzene, 1,2,3-trimethyl- | 120                    | C9H12   | 000526-73-8 | 97   |
| 2       |            | Benzene, 1,2,4-trimethyl- | 120                    | C9H12   | 000095-63-6 | 95   |
| 3       |            | Benzene, 1,2,4-trimethyl- | 120                    | C9H12   | 000095-63-6 | 94   |
| 4       |            | Benzene, 1,2,4-trimethyl- | 120                    | C9H12   | 000095-63-6 | 94   |
| 5       |            | Benzene, 1,2,4-trimethyl- | 120                    | C9H12   | 000095-63-6 | 93   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

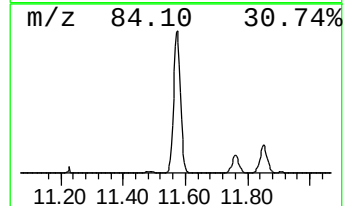
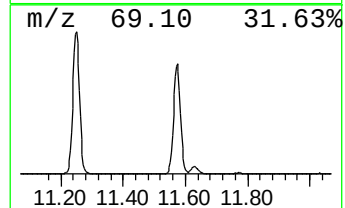
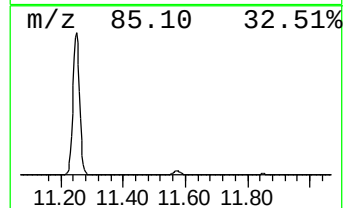
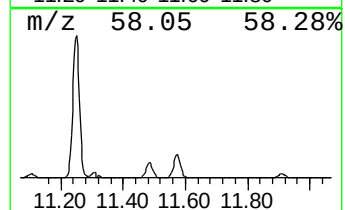
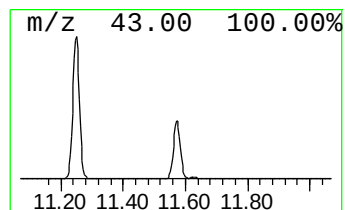
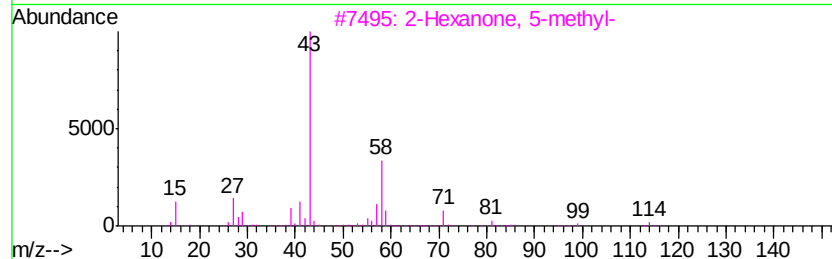
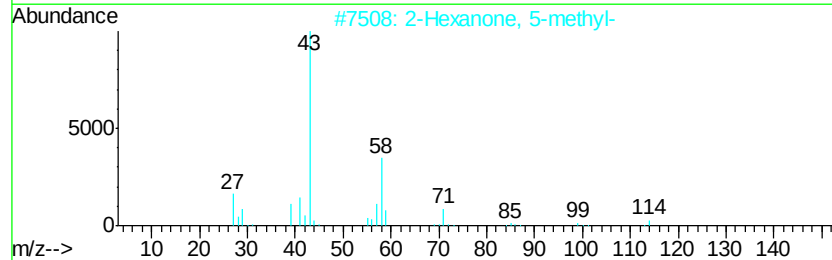
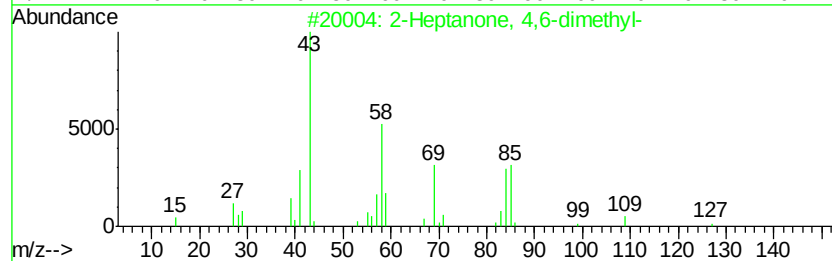
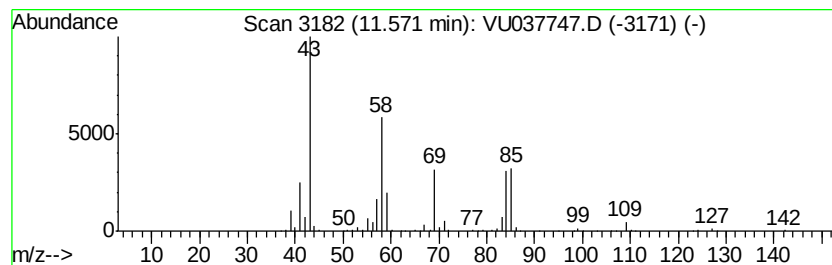
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 2-Heptanone, 4,6-dimethyl- Concentration Rank 18

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 11.57 | 1.93 ug/L | 761495 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Heptanone, 4,6-dimethyl- | 142 | C9H18O  | 019549-80-5 | 90   |
| 2    |    |   | 2-Hexanone, 5-methyl-      | 114 | C7H14O  | 000110-12-3 | 47   |
| 3    |    |   | 2-Hexanone, 5-methyl-      | 114 | C7H14O  | 000110-12-3 | 43   |
| 4    |    |   | Hexane, 3-ethyl-           | 114 | C8H18   | 000619-99-8 | 38   |
| 5    |    |   | Hexane, 3-ethyl-           | 114 | C8H18   | 000619-99-8 | 38   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

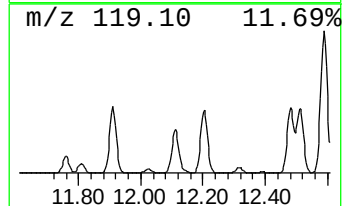
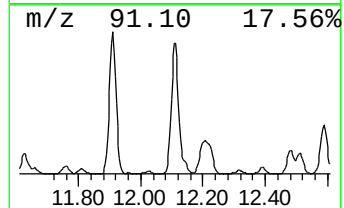
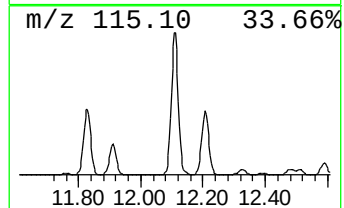
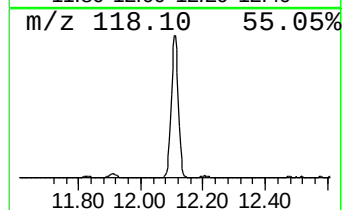
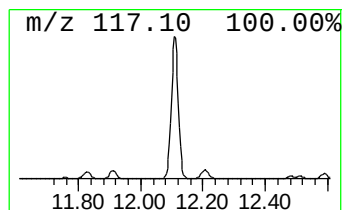
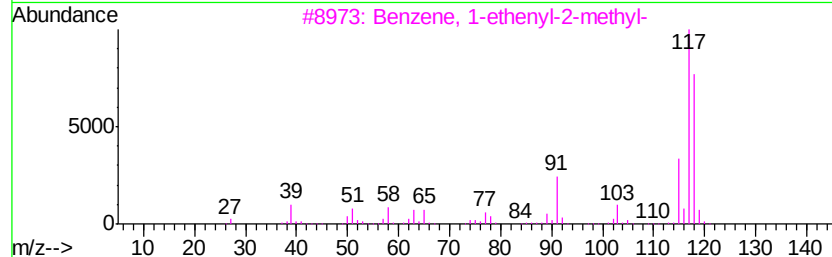
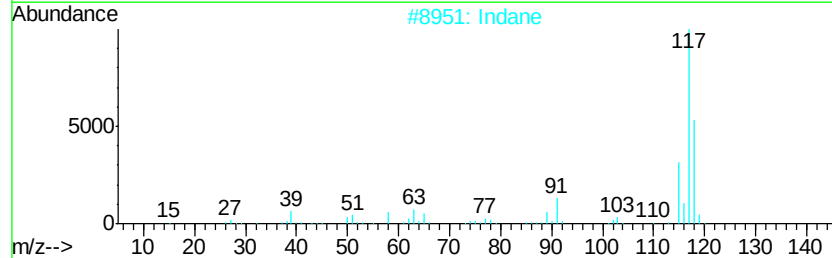
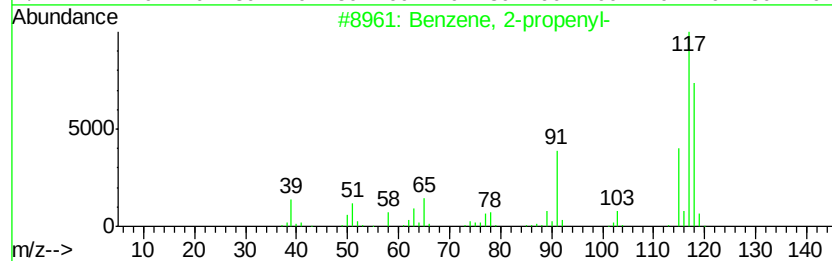
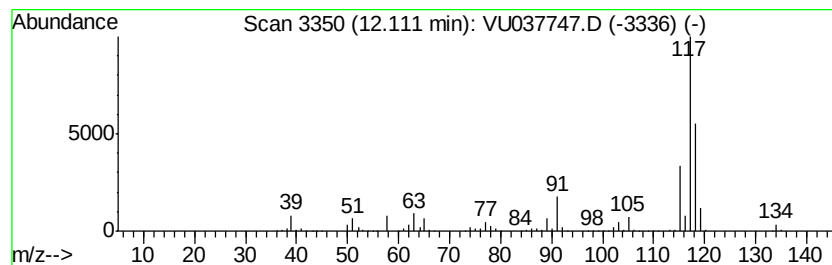
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Benzene, 2-propenyl- Concentration Rank 15

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 12.11 | 4.57 ug/L | 1801590 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-propenyl-         | 118 | C9H10   | 000300-57-2 | 87   |
| 2    |    |   | Indane                       | 118 | C9H10   | 000496-11-7 | 87   |
| 3    |    |   | Benzene, 1-ethenyl-2-methyl- | 118 | C9H10   | 000611-15-4 | 76   |
| 4    |    |   | Indane                       | 118 | C9H10   | 000496-11-7 | 68   |
| 5    |    |   | Benzene, 2-propenyl-         | 118 | C9H10   | 000300-57-2 | 64   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

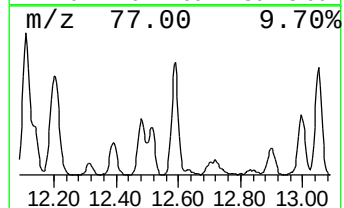
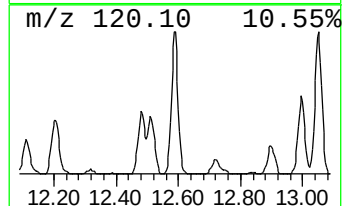
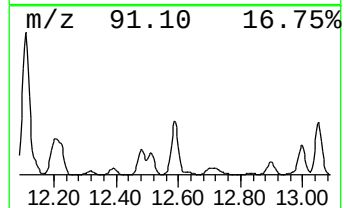
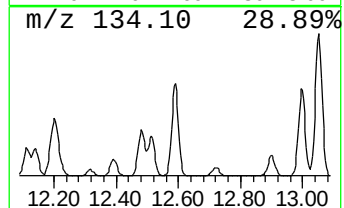
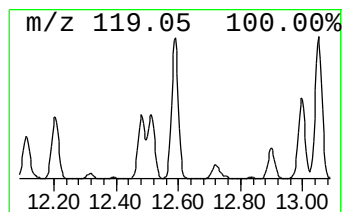
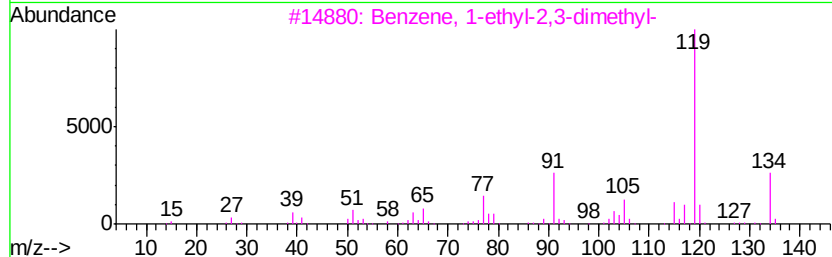
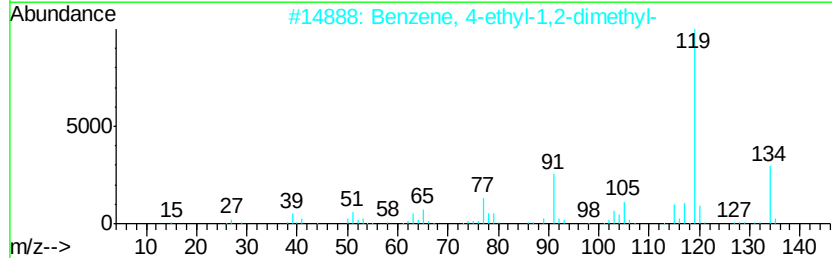
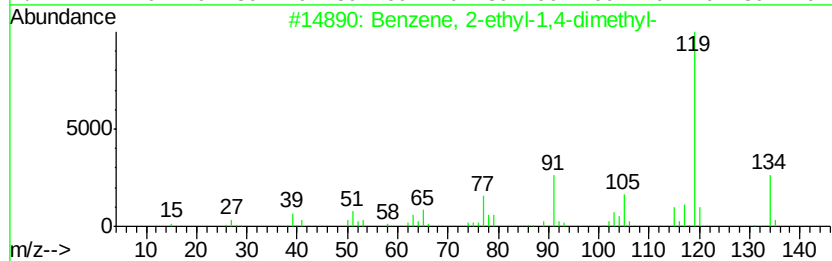
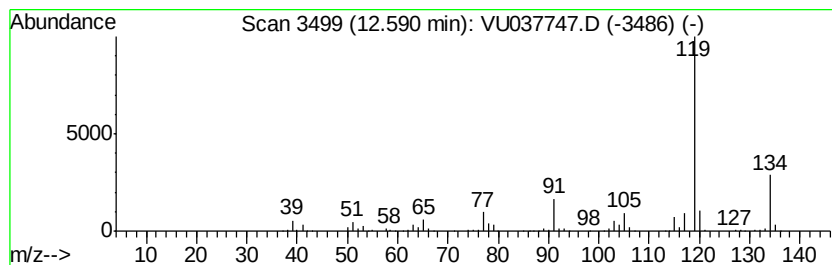
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 21

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.59 | 1.28 ug/L | 503085 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 96   |
| 2    |    | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 95   |
| 3    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 95   |
| 4    |    | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 94   |
| 5    |    | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

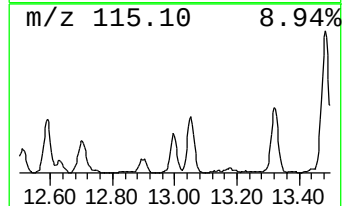
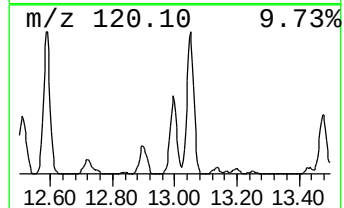
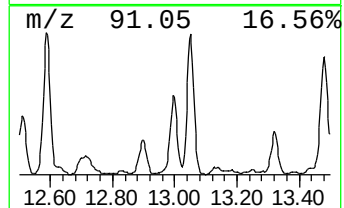
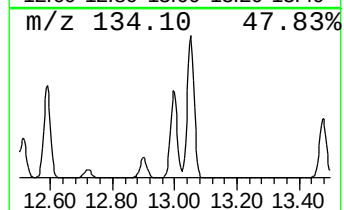
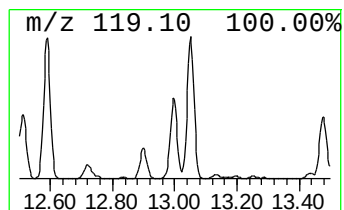
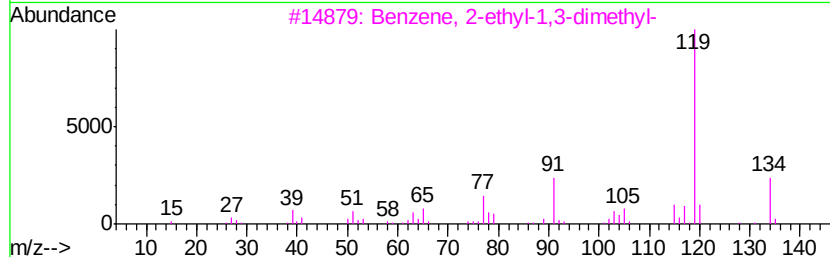
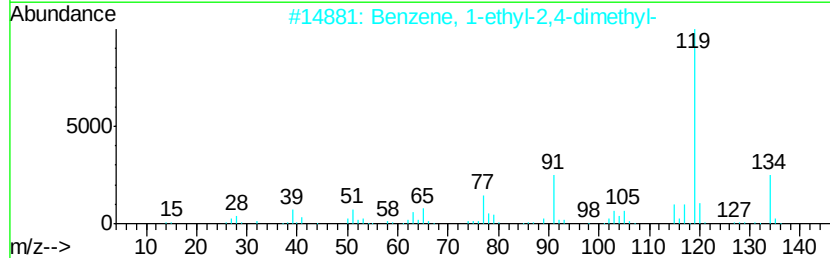
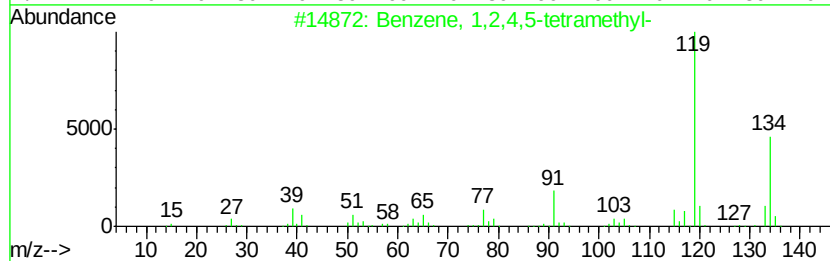
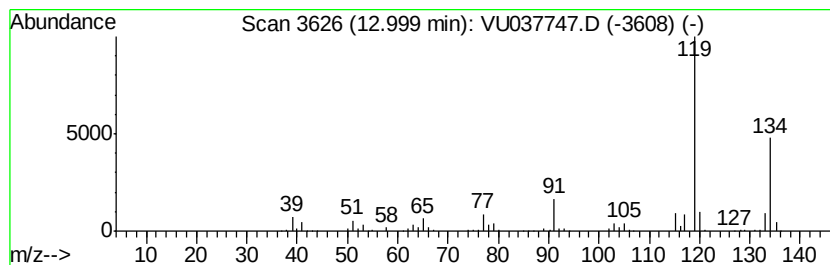
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 22

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 13.00 | 0.82 ug/L | 324353 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 97   |
| 2    |    | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 95   |
| 3    |    | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 95   |
| 4    |    | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 94   |
| 5    |    | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

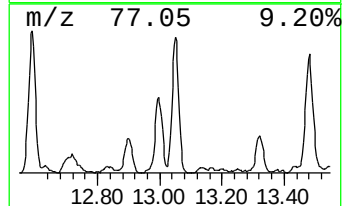
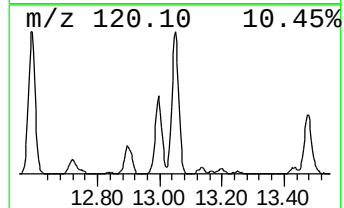
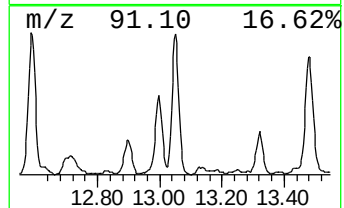
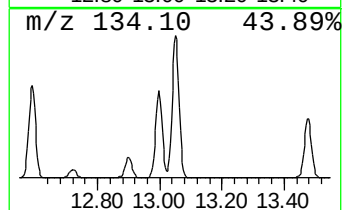
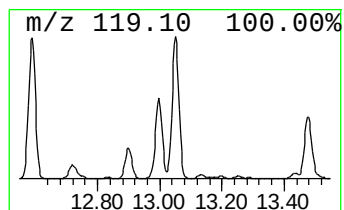
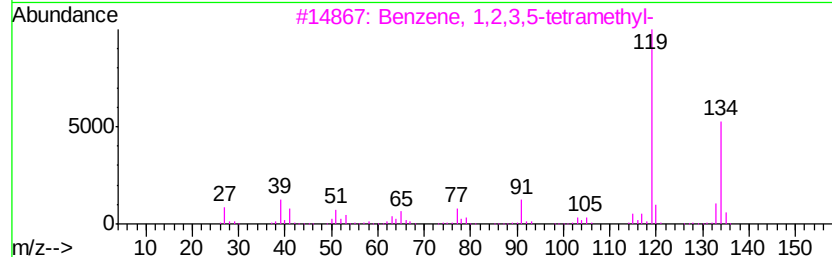
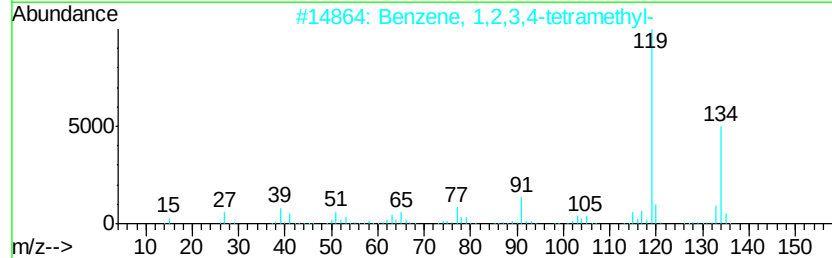
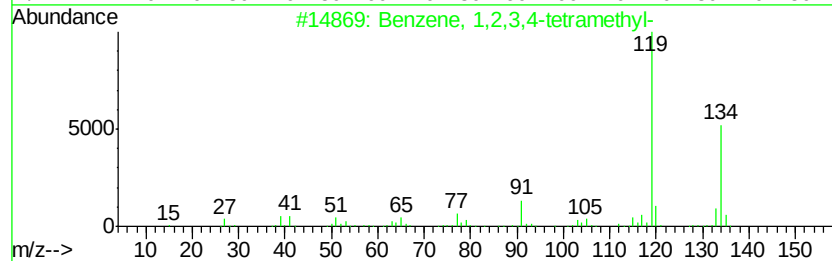
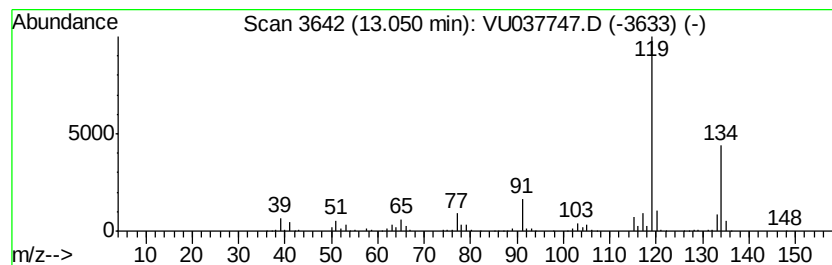
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 20

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 13.05 | 1.41 ug/L | 557218 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 97   |
| 2    |    |   | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 97   |
| 3    |    |   | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 95   |
| 4    |    |   | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 95   |
| 5    |    |   | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU041620\  
Data File : VU037747.D  
Acq On : 16 Apr 2020 12:07  
Operator : JC/MD  
Sample : L2296-03  
Misc : 25mL/MSVOA U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFS89

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
Quant Title : TRACE VOA SOM01.0

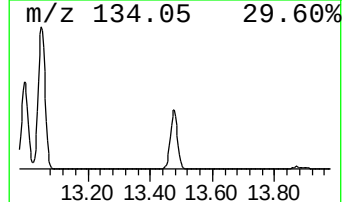
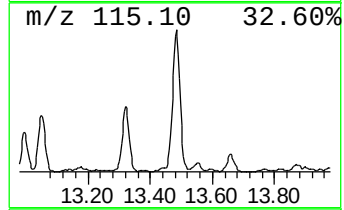
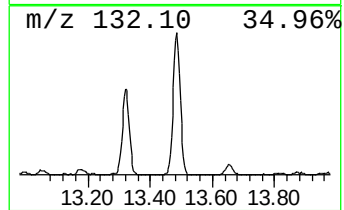
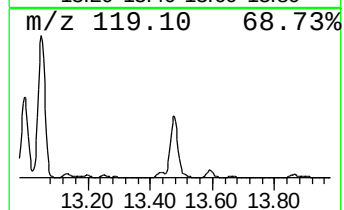
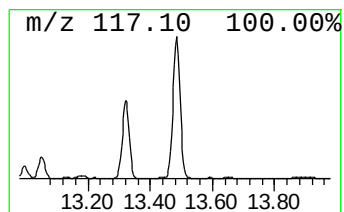
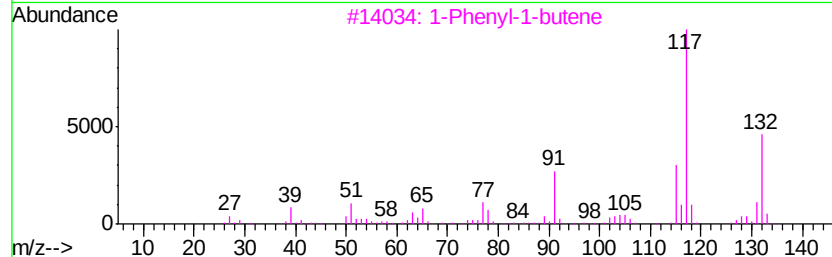
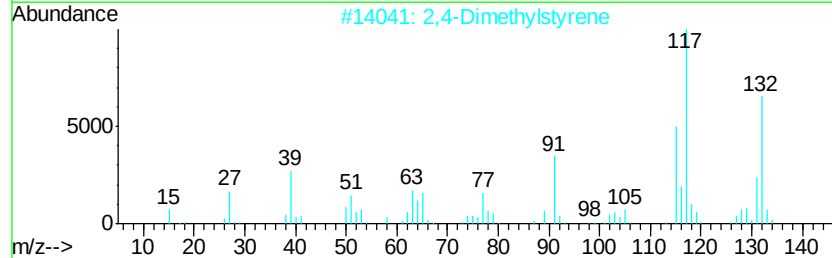
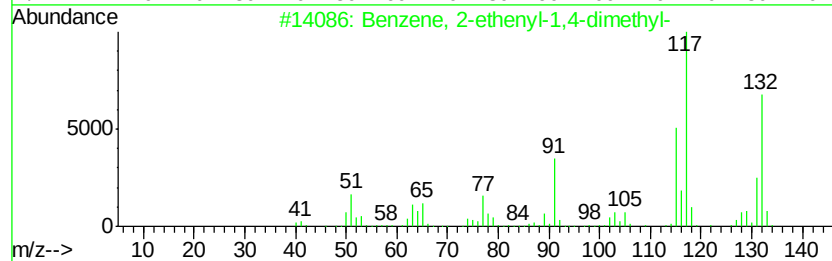
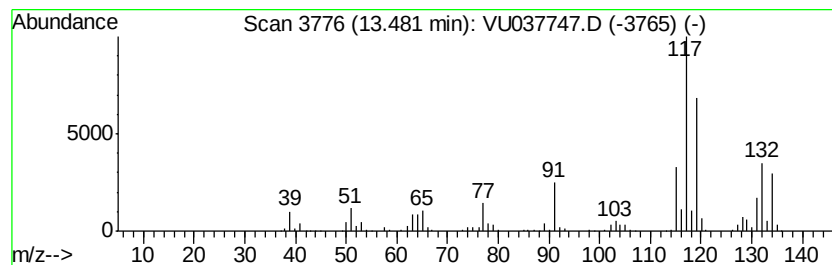
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 19

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 13.48 | 1.48 ug/L | 585134 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12  | 002039-89-6 | 95   |
| 2    |    | 2,4-Dimethylstyrene              | 132 | C10H12  | 002234-20-0 | 95   |
| 3    |    | 1-Phenyl-1-butene                | 132 | C10H12  | 000824-90-8 | 93   |
| 4    |    | Benzene, 2-butenyl-              | 132 | C10H12  | 001560-06-1 | 93   |
| 5    |    | 3-Phenylbut-1-ene                | 132 | C10H12  | 000934-10-1 | 92   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU041620\  
 Data File : VU037747.D  
 Acq On : 16 Apr 2020 12:07  
 Operator : JC/MD  
 Sample : L2296-03  
 Misc : 25mL/MSVOA\_U/WATER  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BFS89

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR041420WMA.M  
 Quant Title : TRACE VOA SOM01.0

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| (DEL) Alkane: Str... | 3.07  | 16.1    | ug/L  | 2043740  | 1                     | 6.28  | 632888  | 5.0  |
| (DEL) Alkane: Str... | 3.11  | 68.6    | ug/L  | 8677850  | 1                     | 6.28  | 632888  | 5.0  |
| (DEL) Alkane: Str... | 3.74  | 125.6   | ug/L  | 15898100 | 1                     | 6.28  | 632888  | 5.0  |
| (DEL) Alkane: Str... | 4.02  | 4.7     | ug/L  | 590851   | 1                     | 6.28  | 632888  | 5.0  |
| (DEL) Alkane: Str... | 4.13  | 4.3     | ug/L  | 539276   | 1                     | 6.28  | 632888  | 5.0  |
| (DEL) Alkane: Str... | 4.34  | 7.7     | ug/L  | 978445   | 1                     | 6.28  | 632888  | 5.0  |
| (DEL) Alkane: Str... | 4.47  | 5.9     | ug/L  | 749838   | 1                     | 6.28  | 632888  | 5.0  |
| (DEL) Alkane: Cyc... | 4.57  | 231.5   | ug/L  | 29307700 | 1                     | 6.28  | 632888  | 5.0  |
| Benzene, propyl-     | 10.92 | 4.8     | ug/L  | 1906830  | 3                     | 11.83 | 1971260 | 5.0  |
| Mesitylene           | 11.10 | 5.9     | ug/L  | 2343590  | 3                     | 11.83 | 1971260 | 5.0  |
| 4-Heptanone, 2,6-... | 11.25 | 22.1    | ug/L  | 8721700  | 3                     | 11.83 | 1971260 | 5.0  |
| Benzene, 1-ethyl-... | 11.31 | 8.7     | ug/L  | 3411060  | 3                     | 11.83 | 1971260 | 5.0  |
| Benzene, 1,2,3-tr... | 11.48 | 23.4    | ug/L  | 9224980  | 3                     | 11.83 | 1971260 | 5.0  |
| 2-Heptanone, 4,6-... | 11.57 | 1.9     | ug/L  | 761495   | 3                     | 11.83 | 1971260 | 5.0  |
| Benzene, 2-propenyl- | 12.11 | 4.6     | ug/L  | 1801590  | 3                     | 11.83 | 1971260 | 5.0  |
| Benzene, 2-ethyl-... | 12.59 | 1.3     | ug/L  | 503085   | 3                     | 11.83 | 1971260 | 5.0  |
| Benzene, 1,2,4,5-... | 13.00 | 0.8     | ug/L  | 324353   | 3                     | 11.83 | 1971260 | 5.0  |
| Benzene, 1,2,3,4-... | 13.05 | 1.4     | ug/L  | 557218   | 3                     | 11.83 | 1971260 | 5.0  |
| Benzene, 2-etheny... | 13.48 | 1.5     | ug/L  | 585134   | 3                     | 11.83 | 1971260 | 5.0  |