

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU021220\  
 Data File : VU036761.D  
 Acq On : 12 Feb 2020 19:41  
 Operator : JC/MD  
 Sample : L1487-17  
 Misc : 25.0mL/MSVOA U/WATER  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 DBCW6

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\SOMUTR021220WMA.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         | 9.713       | 2587          | 2604        | 2635         | rVB      | 1294263        | 2122328       | 1.02%           | 0.534%        |
| 2         | 11.484      | 3136          | 3155        | 3167         | rBV      | 2705998        | 4182300       | 2.02%           | 1.052%        |
| 3         | 11.751      | 3223          | 3238        | 3249         | rBV3     | 1200237        | 2401972       | 1.16%           | 0.604%        |
| 4         | 12.118      | 3334          | 3352        | 3369         | rBV      | 21493155       | 33539658      | 16.19%          | 8.437%        |
| 5         | 12.340      | 3403          | 3421        | 3434         | rBV2     | 30454863       | 49030805      | 23.67%          | 12.334%       |
| 6         | 12.947      | 3599          | 3610        | 3619         | rBV      | 1561714        | 2498493       | 1.21%           | 0.629%        |
| 7         | 13.053      | 3632          | 3643        | 3653         | rBV      | 3756171        | 6801101       | 3.28%           | 1.711%        |
| 8         | 13.484      | 3756          | 3777        | 3787         | rBV2     | 1754962        | 2940375       | 1.42%           | 0.740%        |
| 9         | 13.549      | 3787          | 3797        | 3817         | rVB      | 1797069        | 2963938       | 1.43%           | 0.746%        |
| 10        | 13.661      | 3817          | 3832        | 3847         | rBV2     | 1710031        | 2975455       | 1.44%           | 0.748%        |
| 11        | 14.150      | 3957          | 3984        | 3993         | rBV6     | 63264367       | 207183899     | 100.00%         | 52.119%       |
| 12        | 14.266      | 4008          | 4020        | 4032         | rBV      | 7250055        | 10927678      | 5.27%           | 2.749%        |
| 13        | 15.278      | 4321          | 4335        | 4358         | rVV      | 31175367       | 51011296      | 24.62%          | 12.832%       |
| 14        | 15.462      | 4379          | 4392        | 4419         | rVB      | 11925030       | 18945361      | 9.14%           | 4.766%        |

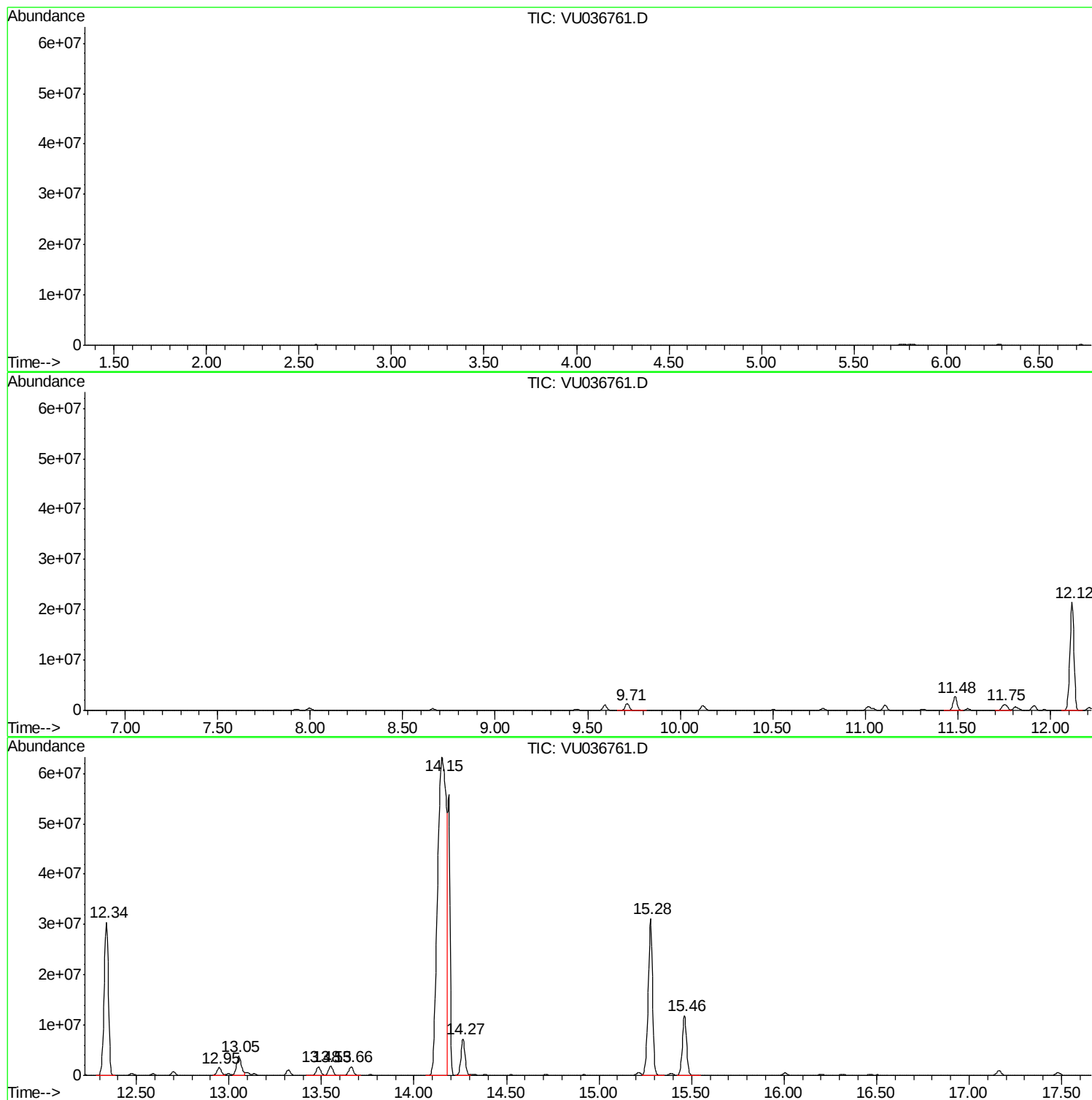
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Instrument :  
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ClientSampleId :  
DBCW6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.M  
Quant Title : TRACE VOA SOM01.0

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



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ClientSampleId :  
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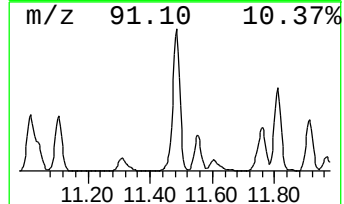
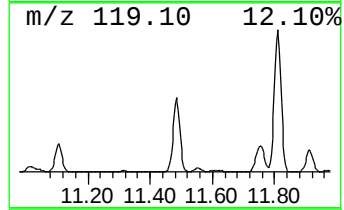
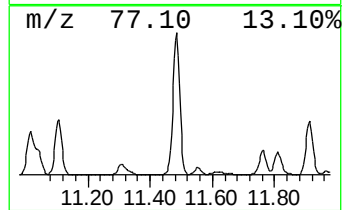
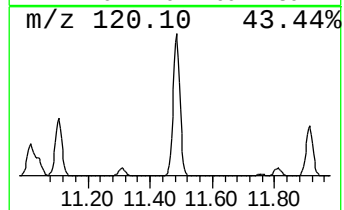
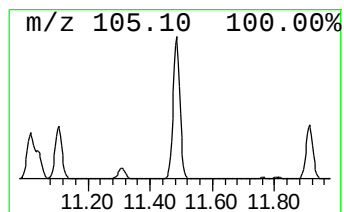
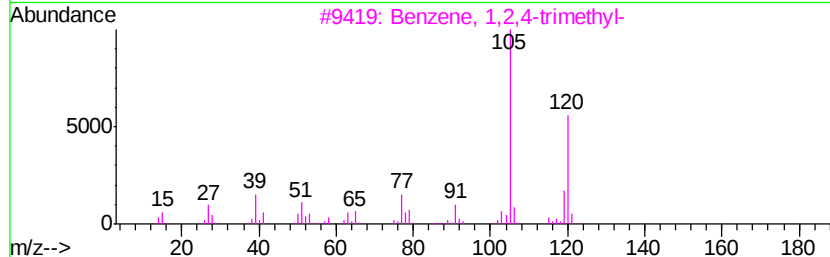
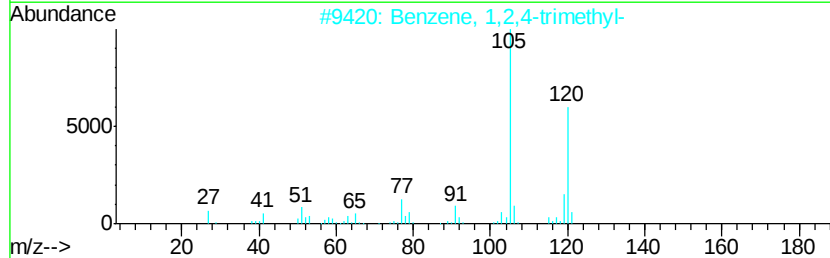
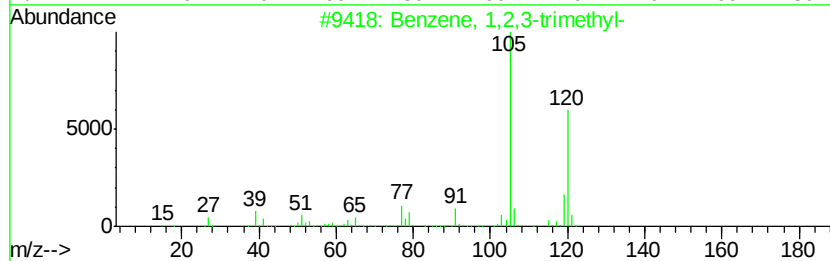
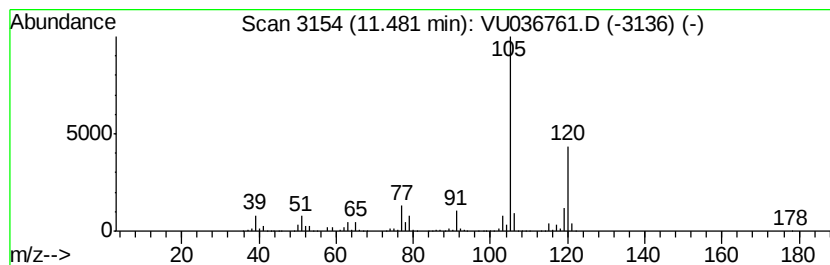
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Quant Title : TRACE VOA SOM01.0

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

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Peak Number 1 Benzene, 1,2,3-trimethyl- Concentration Rank 8

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 11.48 | 8.71 ug/L | 4182300 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | 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 | 95   |
| 4    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 5    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 94   |



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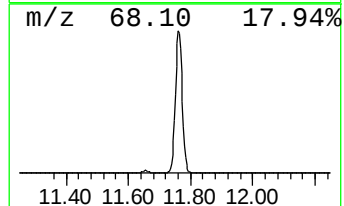
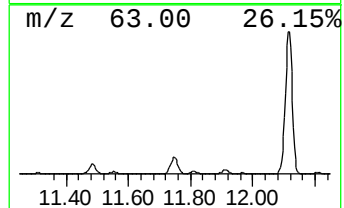
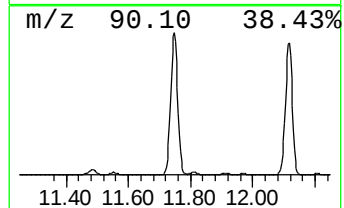
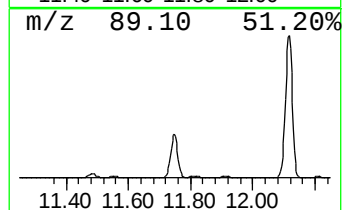
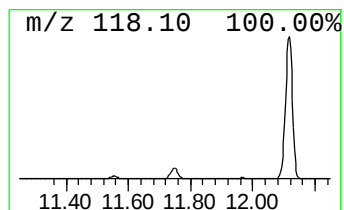
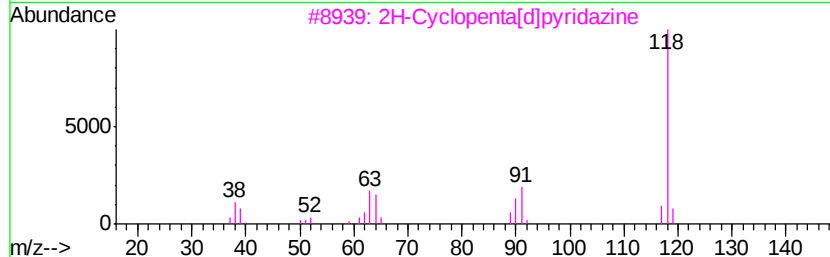
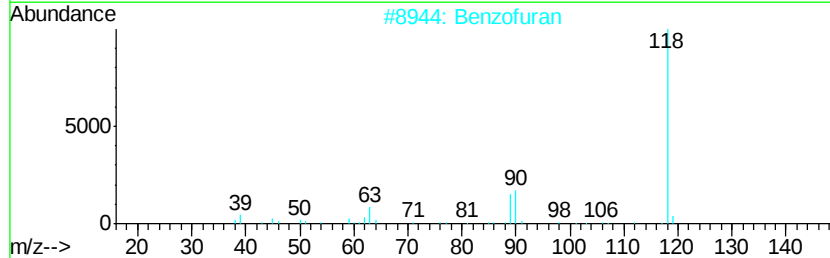
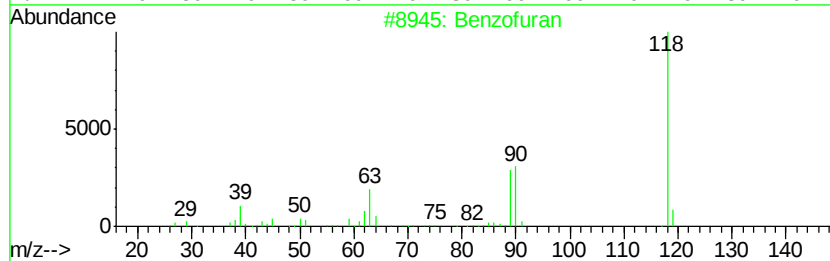
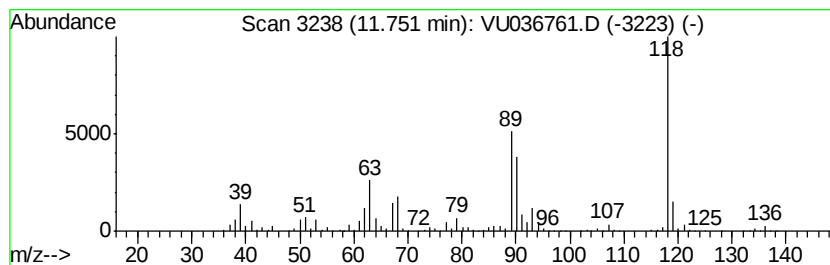
Instrument :  
MSVOA\_U  
ClientSampleId :  
DBCW6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.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 Benzofuran Concentration Rank 13

| R.T.    | EstConc   | Area                       | Relative to ISTD       | R.T.           |
|---------|-----------|----------------------------|------------------------|----------------|
| 11.75   | 5.00 ug/L | 2401970                    | 1,4-Dichlorobenzene-d4 | 11.83          |
| Hit# of | 5         | Tentative ID               | MW MolForm             | CAS# Qual      |
| 1       |           | Benzofuran                 | 118 C8H6O              | 000271-89-6 81 |
| 2       |           | Benzofuran                 | 118 C8H6O              | 000271-89-6 80 |
| 3       |           | 2H-Cyclopenta[d]pyridazine | 118 C7H6N2             | 000270-64-4 53 |
| 4       |           | 1H-Benzimidazole           | 118 C7H6N2             | 000051-17-2 49 |
| 5       |           | 3-Hydroxyphenylacetylene   | 118 C8H6O              | 010401-11-3 43 |



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ALS Vial : 41 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
DBCW6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.M  
Quant Title : TRACE VOA SOM01.0

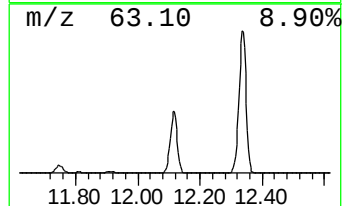
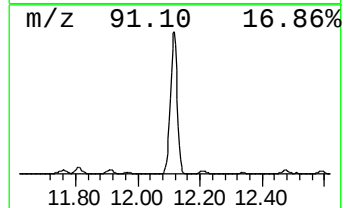
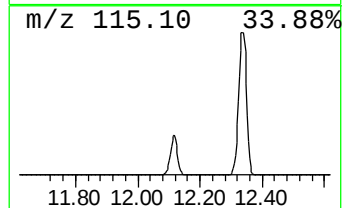
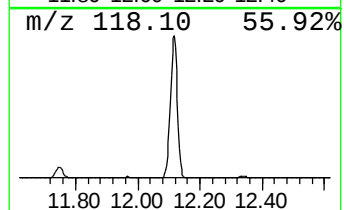
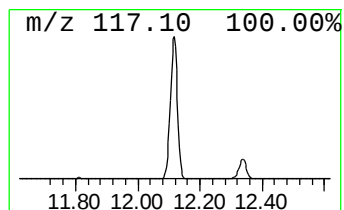
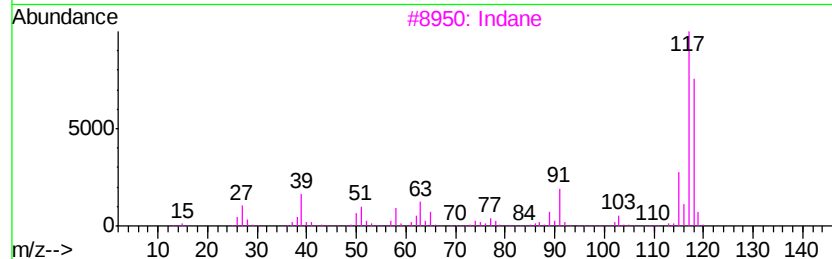
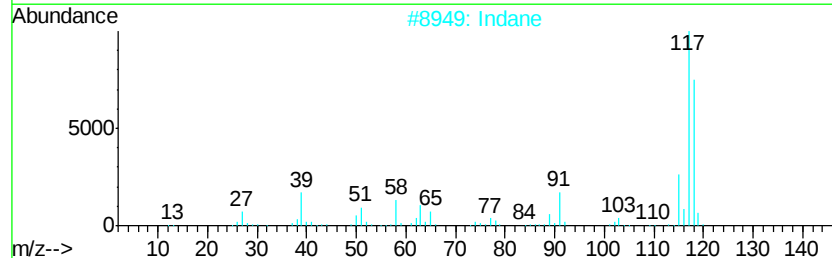
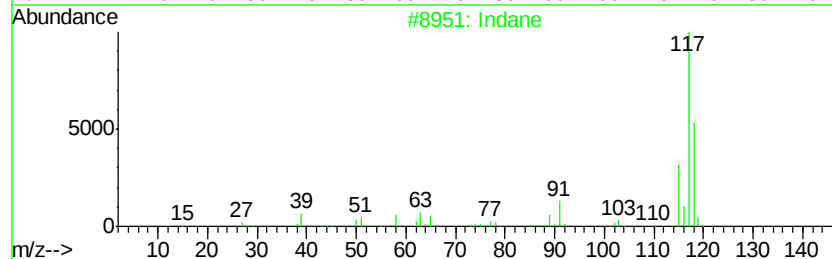
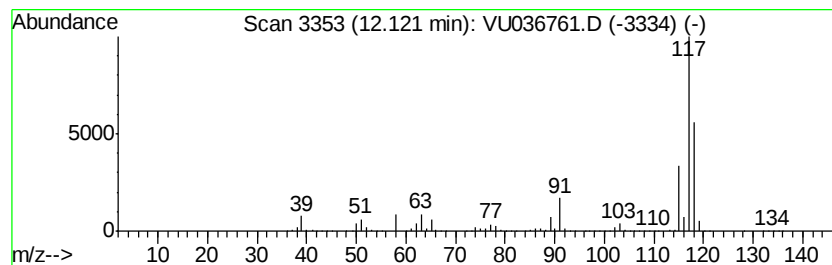
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Indane Concentration Rank 4

| R.T.  | EstConc    | Area     | Relative to ISTD       | R.T.  |
|-------|------------|----------|------------------------|-------|
| 12.12 | 69.82 ug/L | 33539700 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# of | 5                     | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-----------------------|--------------|-----|---------|-------------|------|
| 1       | Indane                |              | 118 | C9H10   | 000496-11-7 | 93   |
| 2       | Indane                |              | 118 | C9H10   | 000496-11-7 | 93   |
| 3       | Indane                |              | 118 | C9H10   | 000496-11-7 | 90   |
| 4       | Indane                |              | 118 | C9H10   | 000496-11-7 | 81   |
| 5       | Benzene, cyclopropyl- |              | 118 | C9H10   | 000873-49-4 | 74   |



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ALS Vial : 41 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
DBCW6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.M  
Quant Title : TRACE VOA SOM01.0

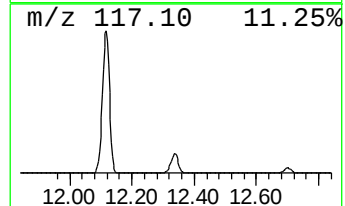
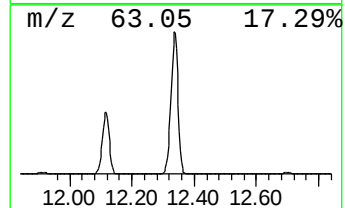
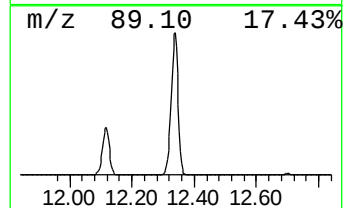
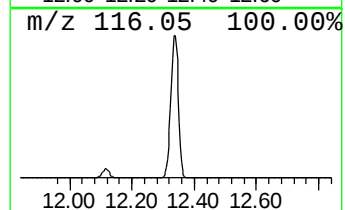
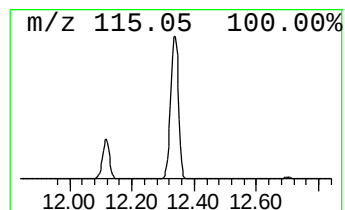
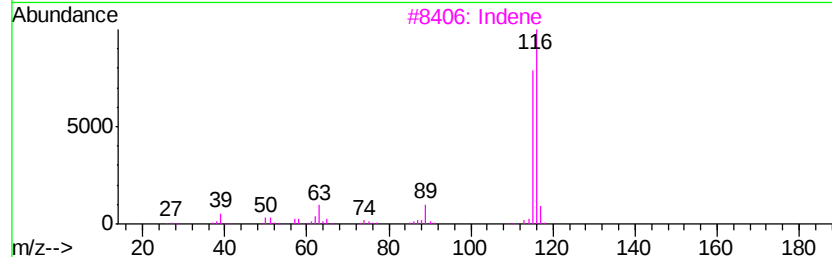
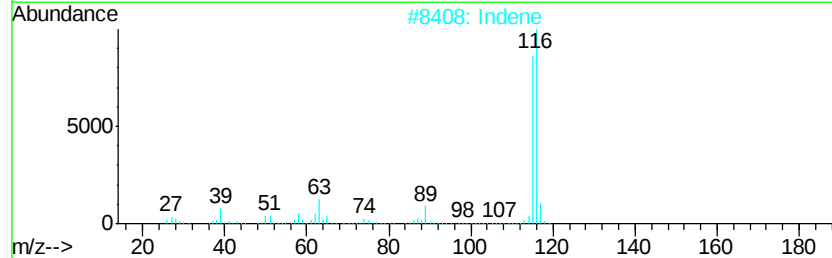
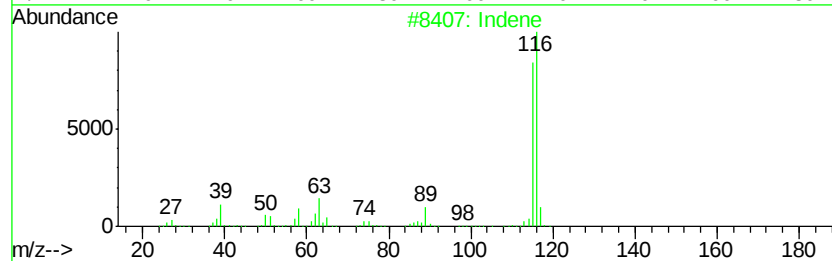
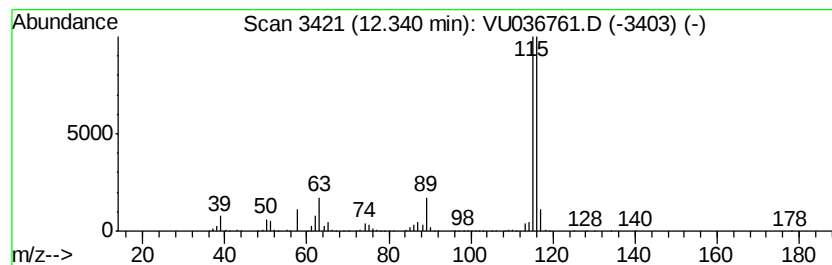
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Indene Concentration Rank 3

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 12.34 | 102.06 ug/L | 49030800 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# of | 5                            | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|------------------------------|--------------|-----|---------|-------------|------|
| 1       | Indene                       |              | 116 | C9H8    | 000095-13-6 | 97   |
| 2       | Indene                       |              | 116 | C9H8    | 000095-13-6 | 96   |
| 3       | Indene                       |              | 116 | C9H8    | 000095-13-6 | 94   |
| 4       | Benzene, 1-propynyl-         |              | 116 | C9H8    | 000673-32-5 | 91   |
| 5       | Benzene, 1-ethynyl-4-methyl- |              | 116 | C9H8    | 000766-97-2 | 80   |



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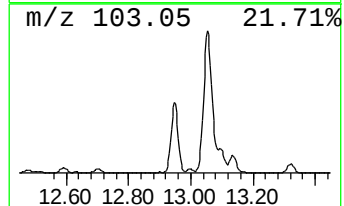
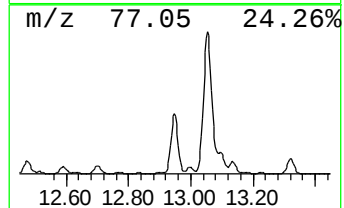
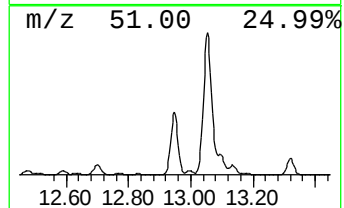
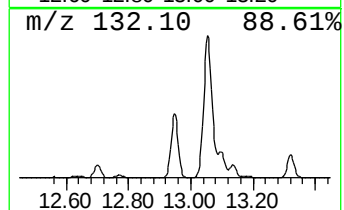
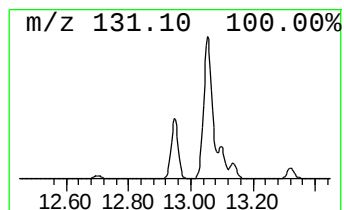
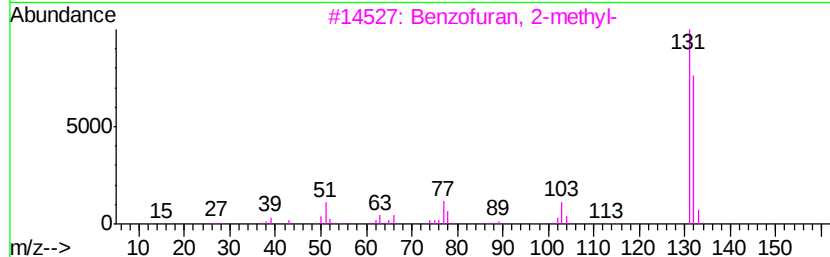
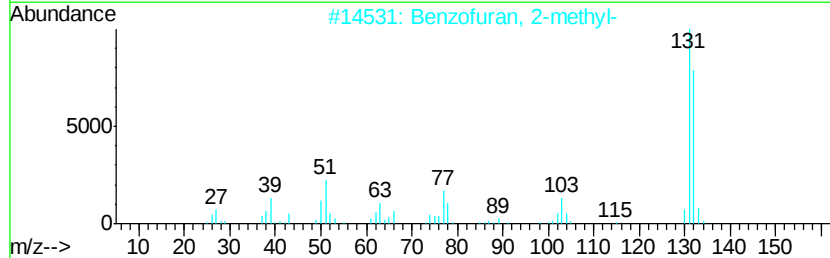
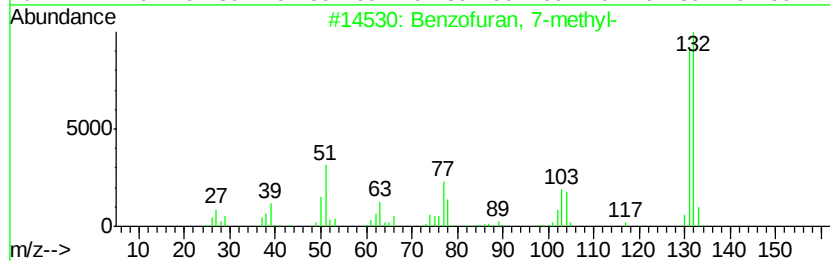
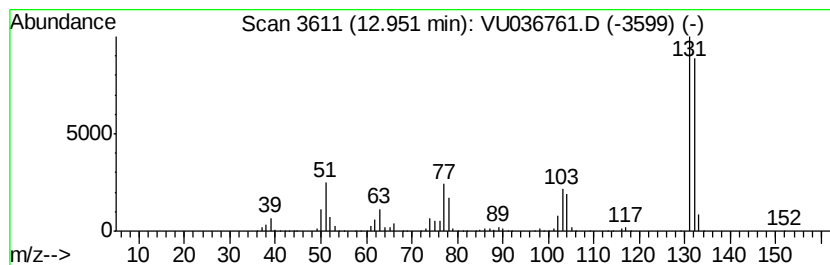
Instrument :  
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Quant Title : TRACE VOA SOM01.0

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

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Peak Number 5 Benzofuran, 7-methyl- Concentration Rank 12

| R.T.    | EstConc   | Area                   | Relative to ISTD       | R.T.           |
|---------|-----------|------------------------|------------------------|----------------|
| 12.95   | 5.20 ug/L | 2498490                | 1,4-Dichlorobenzene-d4 | 11.83          |
| Hit# of | 5         | Tentative ID           | MW MolForm             | CAS# Qual      |
| 1       |           | Benzofuran, 7-methyl-  | 132 C9H8O              | 017059-52-8 97 |
| 2       |           | Benzofuran, 2-methyl-  | 132 C9H8O              | 004265-25-2 94 |
| 3       |           | Benzofuran, 2-methyl-  | 132 C9H8O              | 004265-25-2 94 |
| 4       |           | 3-Phenyl-2-propyn-1-ol | 132 C9H8O              | 001504-58-1 91 |
| 5       |           | Benzofuran, 2-methyl-  | 132 C9H8O              | 004265-25-2 91 |



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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.M  
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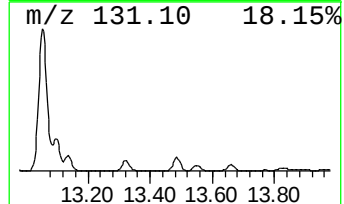
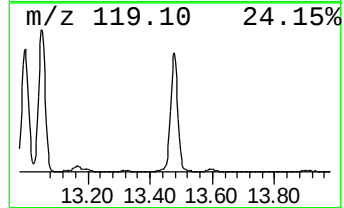
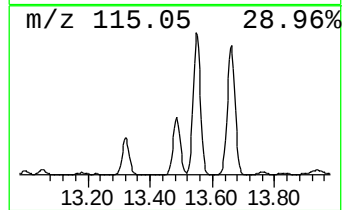
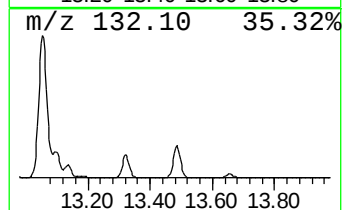
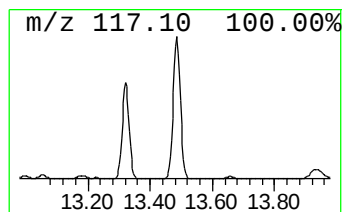
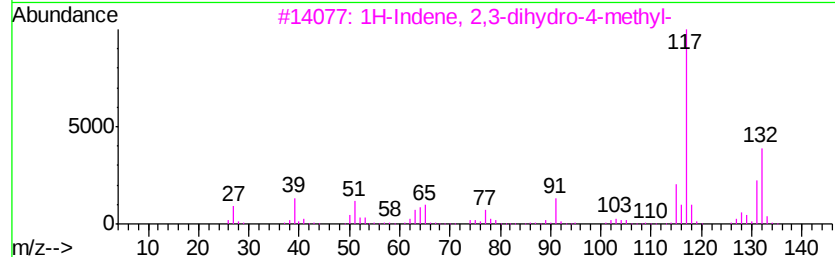
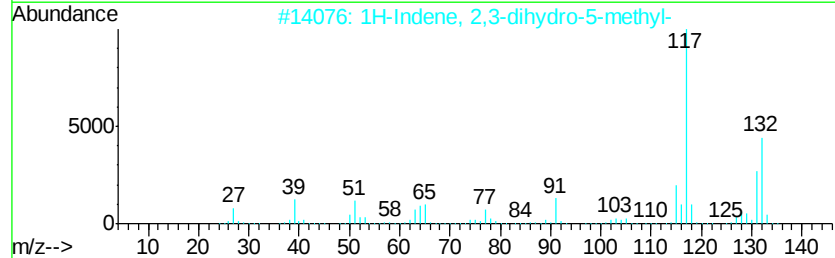
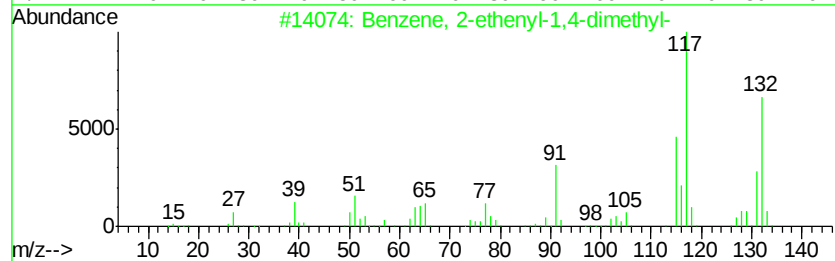
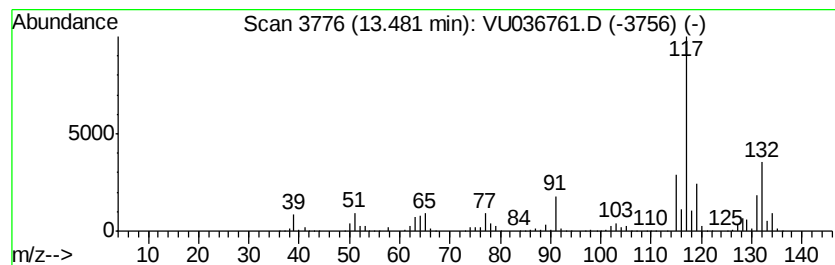
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 11

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 13.48 | 6.12 ug/L | 2940380 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|---------|---|----------------------------------|-----|---------|-------------|------|
| 1       |   | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12  | 002039-89-6 | 96   |
| 2       |   | 1H-Indene, 2,3-dihydro-5-methyl- | 132 | C10H12  | 000874-35-1 | 95   |
| 3       |   | 1H-Indene, 2,3-dihydro-4-methyl- | 132 | C10H12  | 000824-22-6 | 93   |
| 4       |   | 1-Phenyl-1-butene                | 132 | C10H12  | 000824-90-8 | 93   |
| 5       |   | 3-Phenylbut-1-ene                | 132 | C10H12  | 000934-10-1 | 90   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU021220\  
Data File : VU036761.D  
Acq On : 12 Feb 2020 19:41  
Operator : JC/MD  
Sample : L1487-17  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
DBCW6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.M  
Quant Title : TRACE VOA SOM01.0

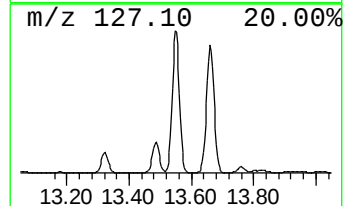
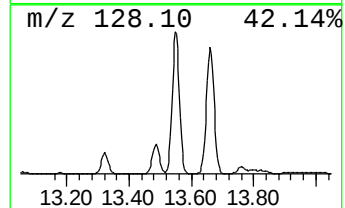
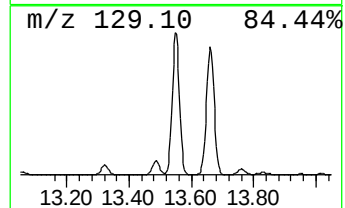
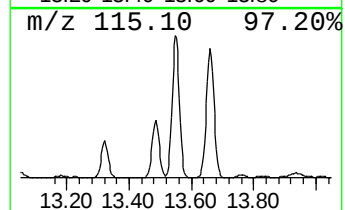
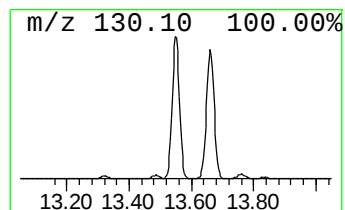
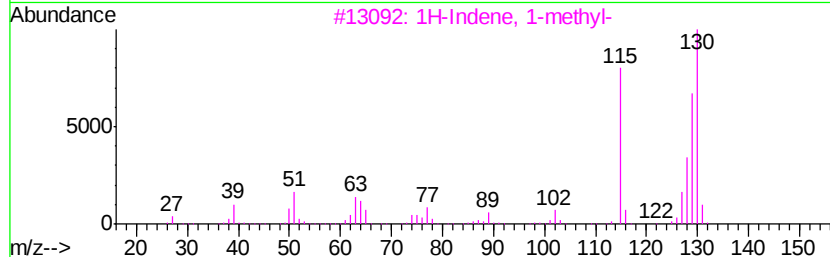
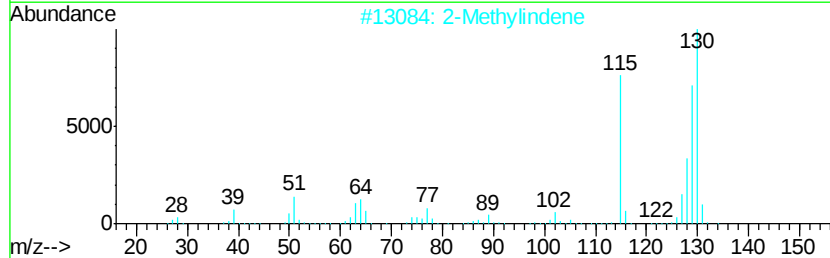
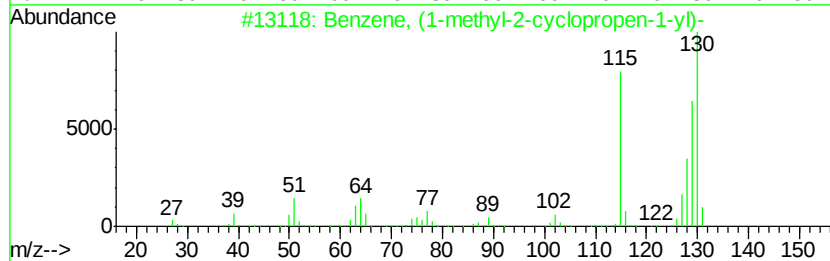
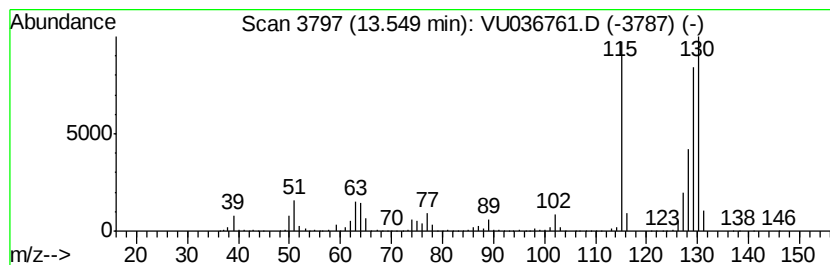
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Benzene, (1-methyl-2-cyclop... Concentration Rank 10

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 13.55 | 6.17 ug/L | 2963940 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (1-methyl-2-cyclopropen... | 130 | C10H10  | 065051-83-4 | 97   |
| 2    |    | 2-Methylindene                      | 130 | C10H10  | 002177-47-1 | 97   |
| 3    |    | 1H-Indene, 1-methyl-                | 130 | C10H10  | 000767-59-9 | 96   |
| 4    |    | Cycloprop[alindene, 1,1a,6,6a-te... | 130 | C10H10  | 015677-15-3 | 93   |
| 5    |    | 1H-Indene, 3-methyl-                | 130 | C10H10  | 000767-60-2 | 93   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU021220\  
Data File : VU036761.D  
Acq On : 12 Feb 2020 19:41  
Operator : JC/MD  
Sample : L1487-17  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
DBCW6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.M  
Quant Title : TRACE VOA SOM01.0

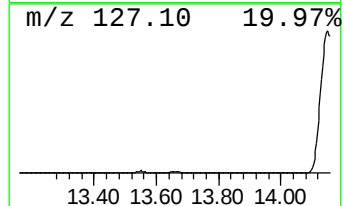
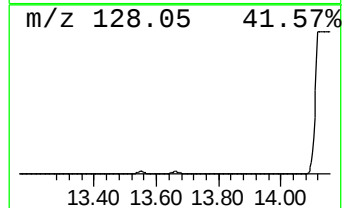
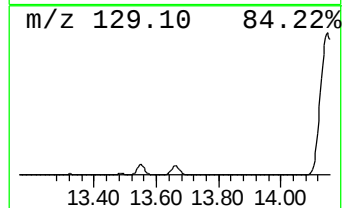
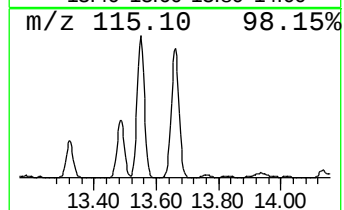
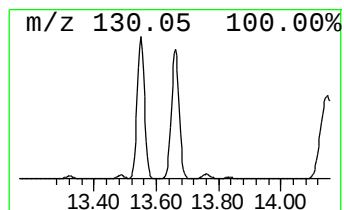
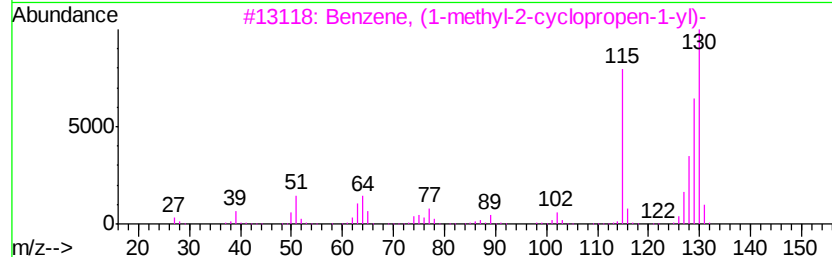
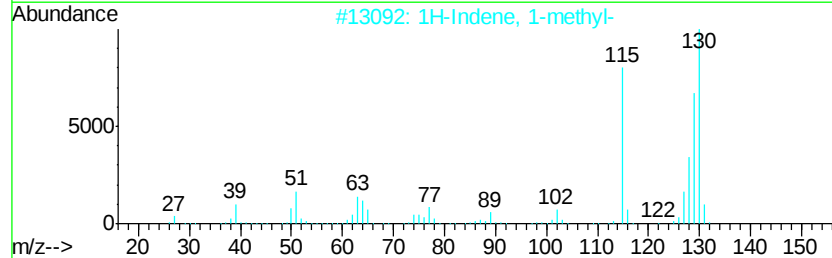
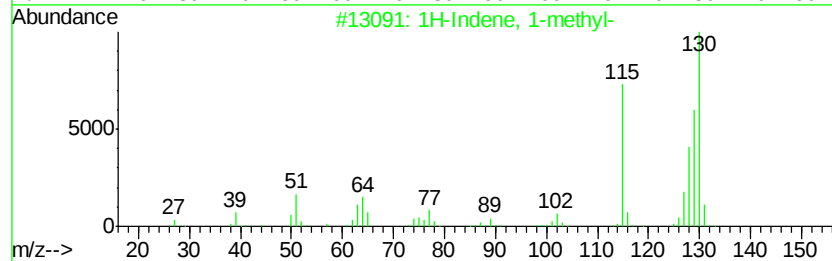
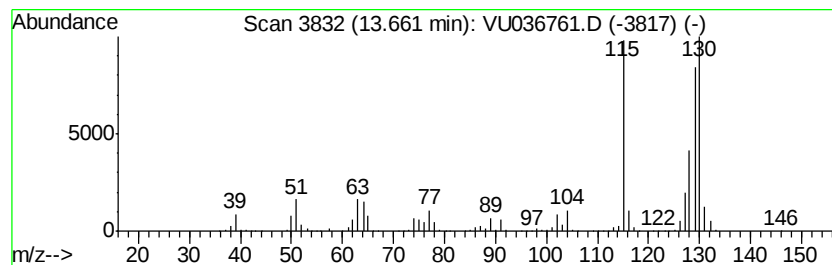
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 1H-Indene, 1-methyl- Concentration Rank 9

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 13.66 | 6.19 ug/L | 2975460 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# of | 5                                   | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------------------------------|--------------|-----|---------|-------------|------|
| 1       | 1H-Indene, 1-methyl-                |              | 130 | C10H10  | 000767-59-9 | 96   |
| 2       | 1H-Indene, 1-methyl-                |              | 130 | C10H10  | 000767-59-9 | 96   |
| 3       | Benzene, (1-methyl-2-cyclopropen... |              | 130 | C10H10  | 065051-83-4 | 96   |
| 4       | 2-Methylindene                      |              | 130 | C10H10  | 002177-47-1 | 96   |
| 5       | Benzene, 1-butynyl-                 |              | 130 | C10H10  | 000622-76-4 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU021220\  
Data File : VU036761.D  
Acq On : 12 Feb 2020 19:41  
Operator : JC/MD  
Sample : L1487-17  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
DBCW6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.M  
Quant Title : TRACE VOA SOM01.0

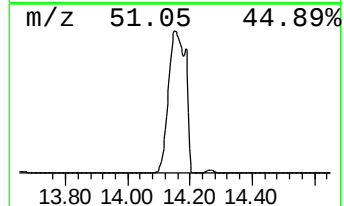
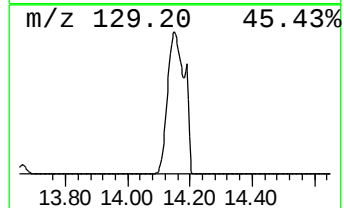
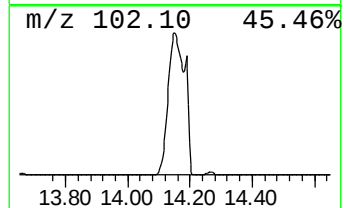
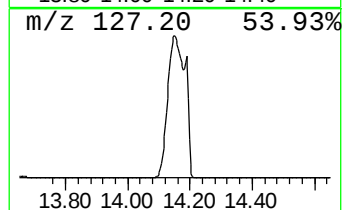
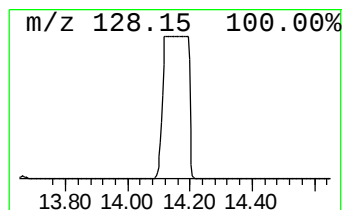
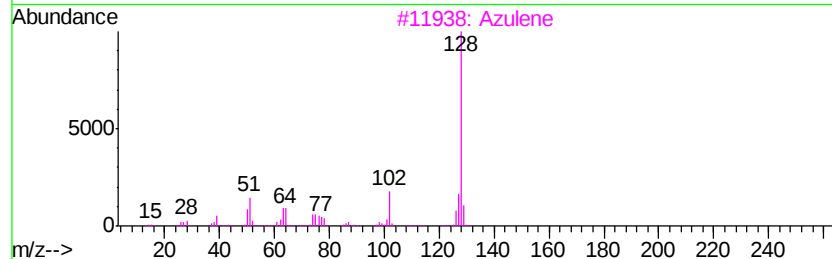
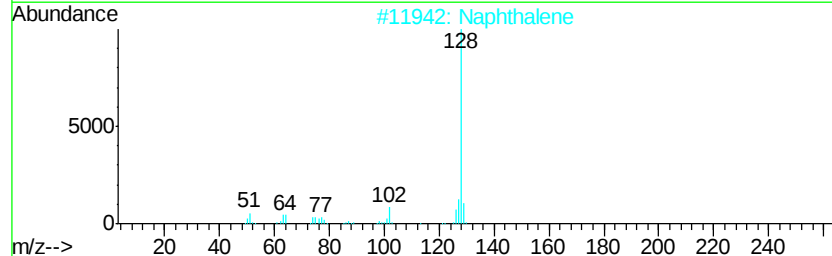
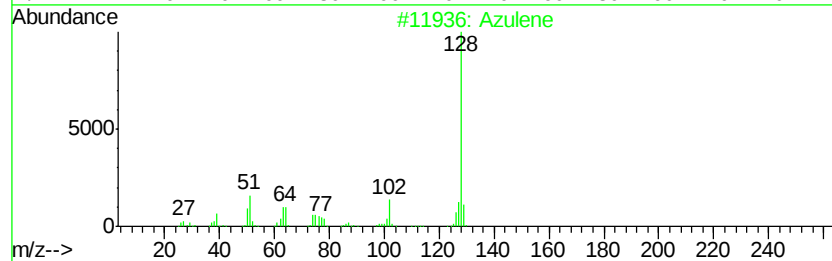
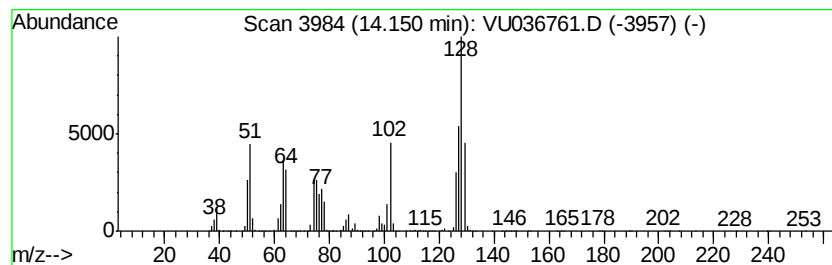
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Azulene Concentration Rank 1

| R.T.  | EstConc     | Area      | Relative to ISTD       | R.T.  |
|-------|-------------|-----------|------------------------|-------|
| 14.15 | 431.28 ug/L | 207184000 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# of | 5                       | Tentative ID | MW  | MolForm | CAS#         | Qual |
|---------|-------------------------|--------------|-----|---------|--------------|------|
| 1       | Azulene                 |              | 128 | C10H8   | 000275-51-4  | 95   |
| 2       | Naphthalene             |              | 128 | C10H8   | 000091-20-3  | 95   |
| 3       | Azulene                 |              | 128 | C10H8   | 000275-51-4  | 90   |
| 4       | 4-Phenylbut-3-ene-1-yne |              | 128 | C10H8   | 1000222-12-0 | 89   |
| 5       | Azulene                 |              | 128 | C10H8   | 000275-51-4  | 76   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU021220\  
Data File : VU036761.D  
Acq On : 12 Feb 2020 19:41  
Operator : JC/MD  
Sample : L1487-17  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
DBCW6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.M  
Quant Title : TRACE VOA SOM01.0

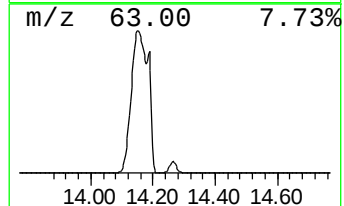
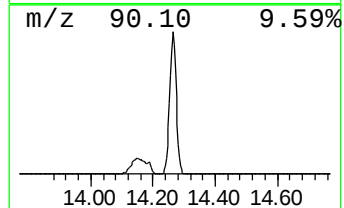
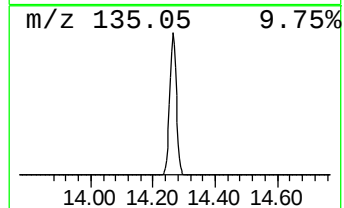
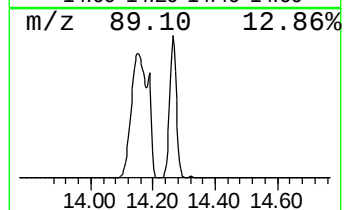
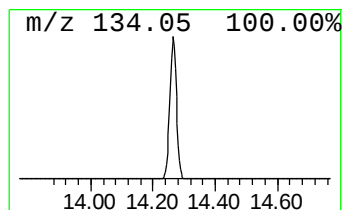
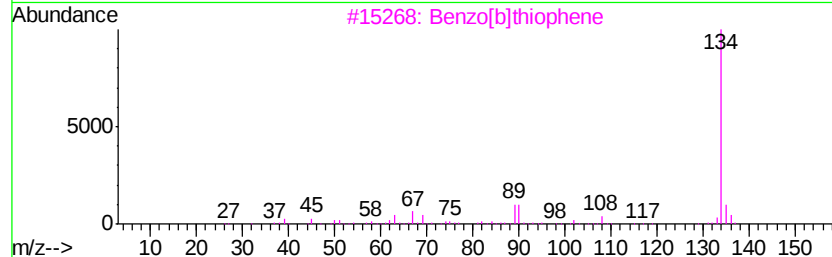
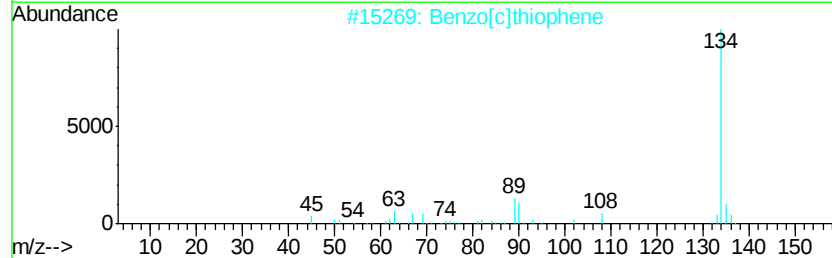
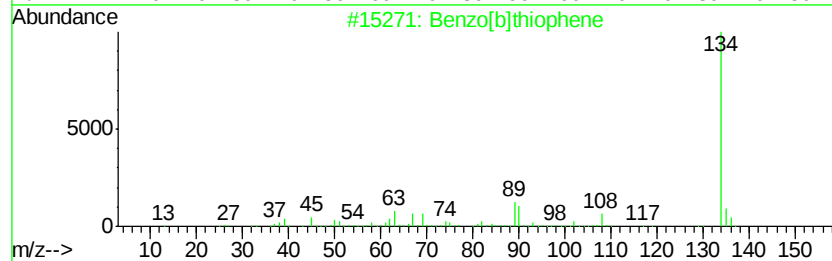
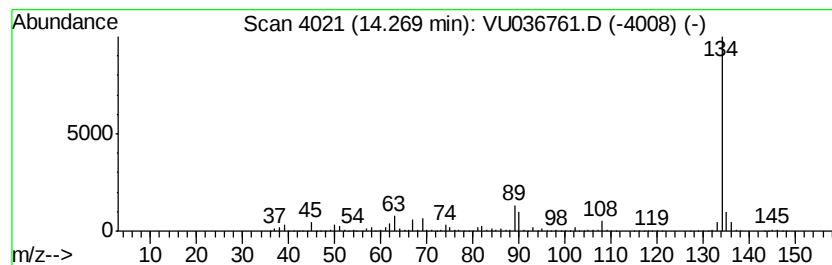
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Benzo[b]thiophene Concentration Rank 6

| R.T.  | EstConc    | Area     | Relative to ISTD       | R.T.  |
|-------|------------|----------|------------------------|-------|
| 14.27 | 22.75 ug/L | 10927700 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[b]thiophene      | 134 | C8H6S   | 000095-15-8 | 97   |
| 2    |    | Benzo[c]thiophene      | 134 | C8H6S   | 000270-82-6 | 97   |
| 3    |    | Benzo[b]thiophene      | 134 | C8H6S   | 000095-15-8 | 95   |
| 4    |    | Benzo[b]thiophene      | 134 | C8H6S   | 000095-15-8 | 95   |
| 5    |    | Cyclopenta[c]thiapyran | 134 | C8H6S   | 000270-63-3 | 94   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU021220\  
Data File : VU036761.D  
Acq On : 12 Feb 2020 19:41  
Operator : JC/MD  
Sample : L1487-17  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
DBCW6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.M  
Quant Title : TRACE VOA SOM01.0

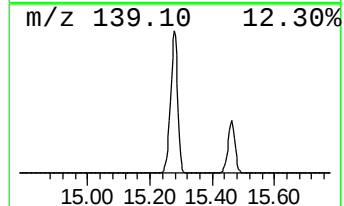
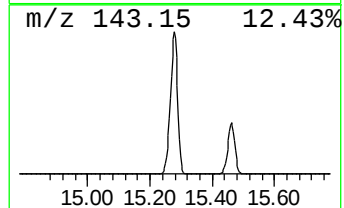
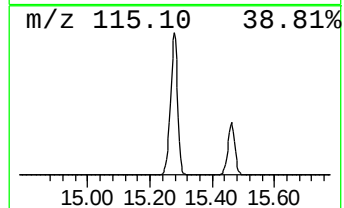
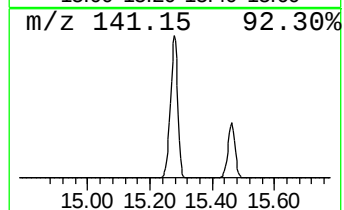
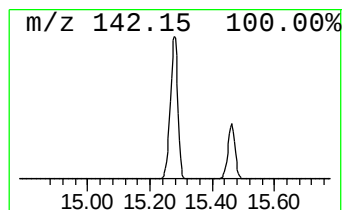
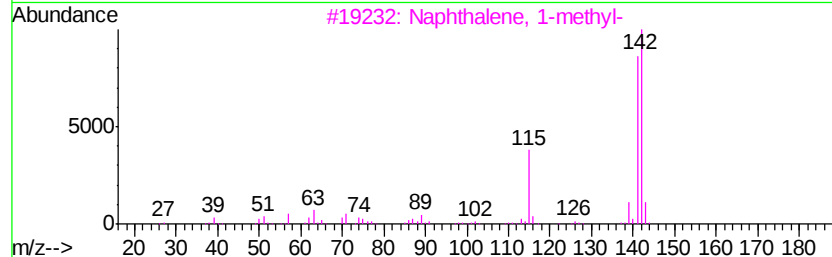
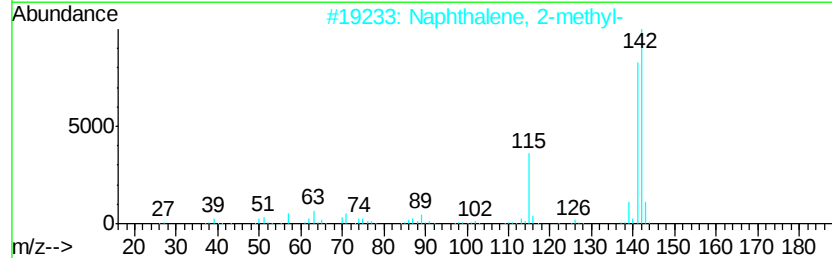
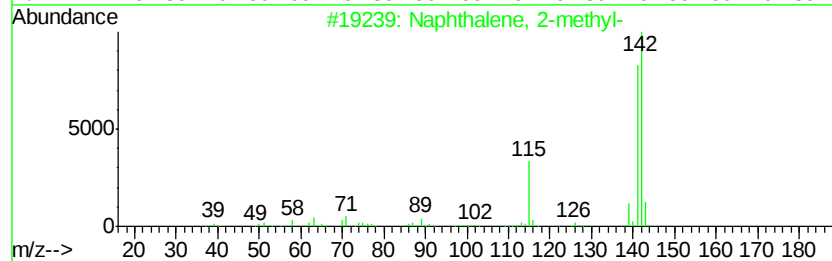
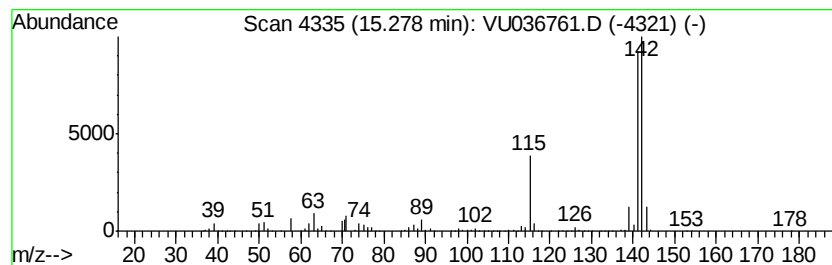
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Naphthalene, 1-methyl- Concentration Rank 2

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 15.28 | 106.19 ug/L | 51011300 | 1,4-Dichlorobenzene-d4 | 11.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 96   |
| 2    |    | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 96   |
| 3    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 96   |
| 4    |    | 1,4-Methanonaphthalene, 1,4-dihy... | 142 | C11H10  | 004453-90-1 | 93   |
| 5    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 93   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU021220\  
 Data File : VU036761.D  
 Acq On : 12 Feb 2020 19:41  
 Operator : JC/MD  
 Sample : L1487-17  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 DBCW6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR021220WMA.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 |
| Benzene, 1,2,3-tr... | 11.48 | 8.7     | ug/L  | 4182300   | 3                     | 11.83 | 2401970 | 5.0  |
| Benzofuran           | 11.75 | 5.0     | ug/L  | 2401970   | 3                     | 11.83 | 2401970 | 5.0  |
| Indane               | 12.12 | 69.8    | ug/L  | 33539700  | 3                     | 11.83 | 2401970 | 5.0  |
| Indene               | 12.34 | 102.1   | ug/L  | 49030800  | 3                     | 11.83 | 2401970 | 5.0  |
| Benzofuran, 7-met... | 12.95 | 5.2     | ug/L  | 2498490   | 3                     | 11.83 | 2401970 | 5.0  |
| Benzene, 2-etheny... | 13.48 | 6.1     | ug/L  | 2940380   | 3                     | 11.83 | 2401970 | 5.0  |
| Benzene, (1-methy... | 13.55 | 6.2     | ug/L  | 2963940   | 3                     | 11.83 | 2401970 | 5.0  |
| 1H-Indene, 1-methyl- | 13.66 | 6.2     | ug/L  | 2975460   | 3                     | 11.83 | 2401970 | 5.0  |
| Azulene              | 14.15 | 431.3   | ug/L  | 207184000 | 3                     | 11.83 | 2401970 | 5.0  |
| Benzo[b]thiophene    | 14.27 | 22.8    | ug/L  | 10927700  | 3                     | 11.83 | 2401970 | 5.0  |
| Naphthalene, 1-me... | 15.28 | 106.2   | ug/L  | 51011300  | 3                     | 11.83 | 2401970 | 5.0  |