

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN070622\  
 Data File : VN073351.D  
 Acq On : 06 Jul 2022 17:38  
 Operator : JC\MD  
 Sample : N3586-04 10X  
 Misc : 1.11g/10mL/100uL/5.00mL/MSVOA\_N/MEOH  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 MSVOA\_N  
 ClientSampleId :  
 31457

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N062822W.M  
 Title : SW846 8260

Signal : TIC: VN073351.D\data.ms

| 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         | 7.998       | 1020          | 1039        | 1049         | rBV3     | 1559583        | 4891107       | 17.49%          | 8.405%        |
| 2         | 8.098       | 1049          | 1056        | 1067         | rVV      | 468766         | 1194372       | 4.27%           | 2.052%        |
| 3         | 8.222       | 1067          | 1077        | 1094         | rVV      | 1864163        | 4429957       | 15.84%          | 7.613%        |
| 4         | 8.369       | 1094          | 1102        | 1112         | rVV      | 242141         | 557837        | 1.99%           | 0.959%        |
| 5         | 8.486       | 1113          | 1122        | 1130         | rVV4     | 250300         | 664462        | 2.38%           | 1.142%        |
| 6         | 8.639       | 1142          | 1148        | 1158         | rVB      | 147226         | 328225        | 1.17%           | 0.564%        |
| 7         | 8.757       | 1159          | 1168        | 1180         | rBV      | 981974         | 1965663       | 7.03%           | 3.378%        |
| 8         | 8.898       | 1184          | 1192        | 1206         | rVB      | 620556         | 1279754       | 4.57%           | 2.199%        |
| 9         | 9.392       | 1261          | 1276        | 1291         | rBV      | 509500         | 1119924       | 4.00%           | 1.925%        |
| 10        | 10.380      | 1435          | 1444        | 1448         | rBV      | 856141         | 1609572       | 5.75%           | 2.766%        |
| 11        | 10.445      | 1448          | 1455        | 1470         | rVB      | 14734315       | 27973134      | 100.00%         | 48.071%       |
| 12        | 11.680      | 1658          | 1665        | 1676         | rBV      | 895525         | 1609276       | 5.75%           | 2.766%        |
| 13        | 11.892      | 1692          | 1701        | 1711         | rBV3     | 372468         | 768726        | 2.75%           | 1.321%        |
| 14        | 12.216      | 1745          | 1756        | 1763         | rBV4     | 139013         | 385785        | 1.38%           | 0.663%        |
| 15        | 12.669      | 1828          | 1833        | 1839         | rBV      | 759564         | 1247443       | 4.46%           | 2.144%        |
| 16        | 12.921      | 1869          | 1876        | 1879         | rBV      | 264236         | 444153        | 1.59%           | 0.763%        |
| 17        | 12.986      | 1879          | 1887        | 1894         | rVB3     | 536836         | 1084255       | 3.88%           | 1.863%        |
| 18        | 13.304      | 1934          | 1941        | 1954         | rVB      | 503418         | 1041699       | 3.72%           | 1.790%        |
| 19        | 13.610      | 1987          | 1993        | 2003         | rVB2     | 994771         | 1791762       | 6.41%           | 3.079%        |
| 20        | 13.933      | 2043          | 2048        | 2053         | rBV      | 470851         | 713816        | 2.55%           | 1.227%        |
| 21        | 14.786      | 2188          | 2193        | 2199         | rVB2     | 406863         | 596764        | 2.13%           | 1.026%        |
| 22        | 14.863      | 2200          | 2206        | 2210         | rVV4     | 168327         | 403082        | 1.44%           | 0.693%        |
| 23        | 14.904      | 2210          | 2213        | 2221         | rVB3     | 148481         | 284733        | 1.02%           | 0.489%        |
| 24        | 15.398      | 2291          | 2297        | 2307         | rBV2     | 196860         | 477615        | 1.71%           | 0.821%        |
| 25        | 15.627      | 2332          | 2336        | 2346         | rVB      | 321107         | 560632        | 2.00%           | 0.963%        |
| 26        | 16.245      | 2434          | 2441        | 2448         | rBV5     | 163088         | 471862        | 1.69%           | 0.811%        |
| 27        | 16.592      | 2495          | 2500        | 2511         | rVB      | 131744         | 295499        | 1.06%           | 0.508%        |

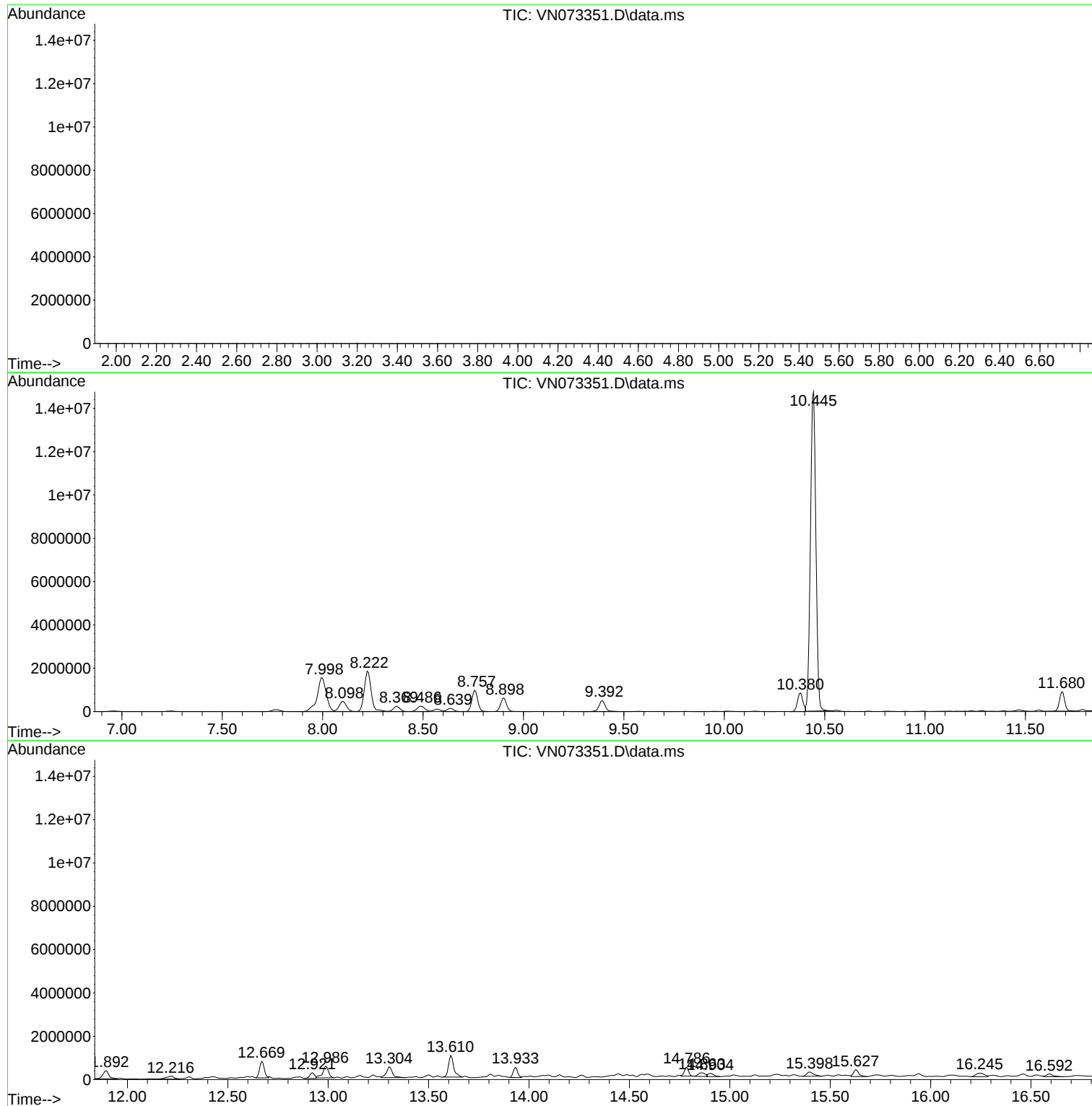
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Instrument :  
MSVOA\_N  
ClientSampleId :  
31457

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N062822W.M  
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P



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Instrument :  
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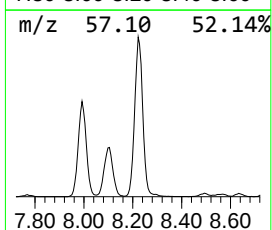
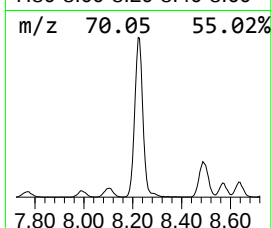
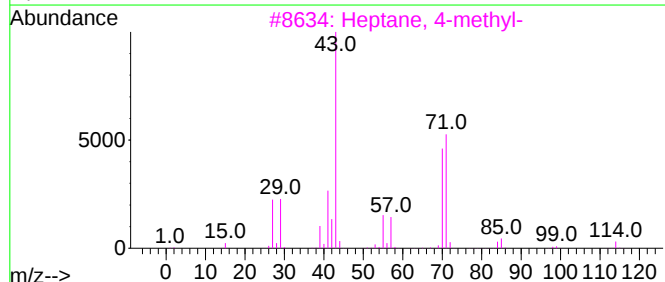
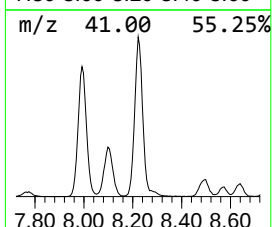
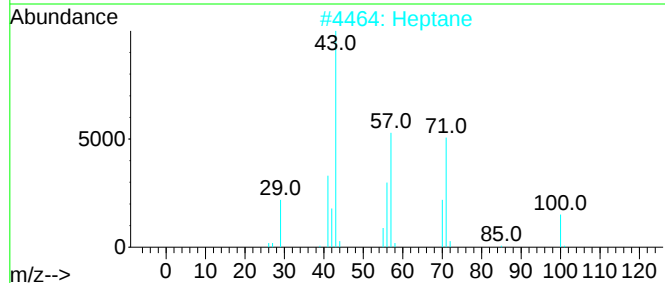
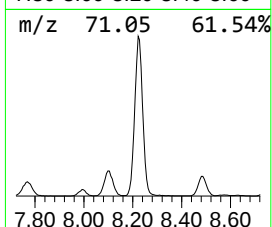
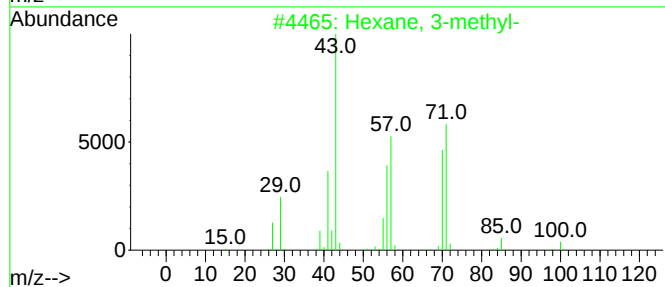
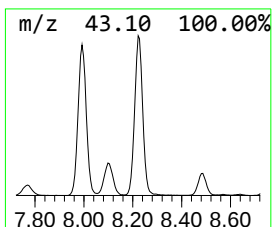
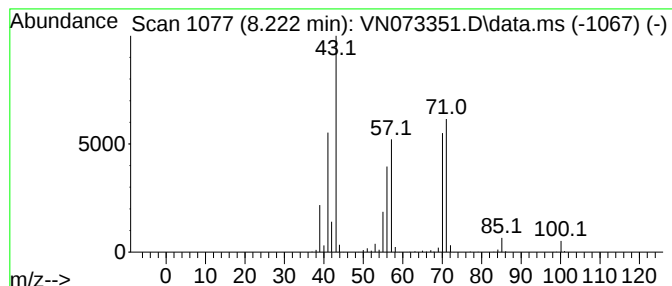
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N062822W.M  
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Hexane, 3-methyl- Concentration Rank 2

| R.T.  | EstConc    | Area    | Relative to ISTD   | R.T.  |
|-------|------------|---------|--------------------|-------|
| 8.222 | 45.29 ug/l | 4429960 | Pentafluorobenzene | 8.016 |

| Hit# | of 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------------|-----|---------|-------------|------|
| 1    |      | Hexane, 3-methyl-         | 100 | C7H16   | 000589-34-4 | 94   |
| 2    |      | Heptane                   | 100 | C7H16   | 000142-82-5 | 80   |
| 3    |      | Heptane, 4-methyl-        | 114 | C8H18   | 000589-53-7 | 59   |
| 4    |      | 1-Butanol, 2-ethyl-       | 102 | C6H14O  | 000097-95-0 | 53   |
| 5    |      | Pentane, 2,3,4-trimethyl- | 114 | C8H18   | 000565-75-3 | 53   |



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ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_N  
ClientSampleId :  
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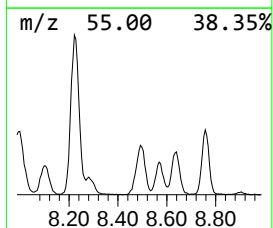
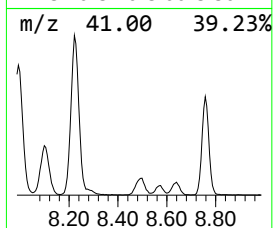
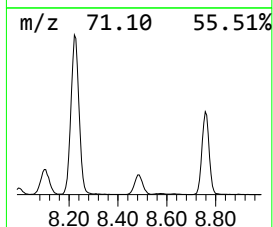
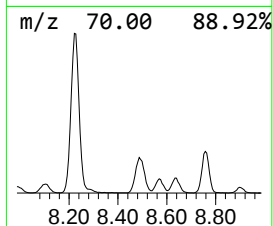
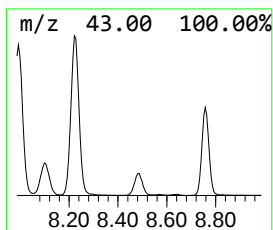
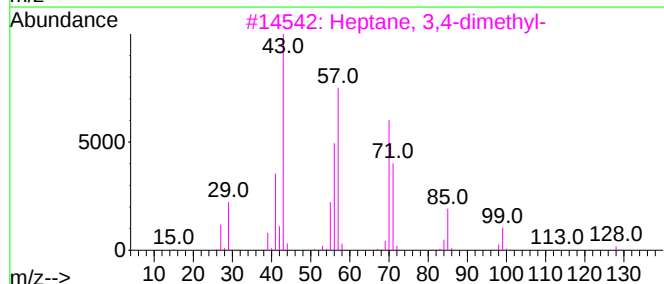
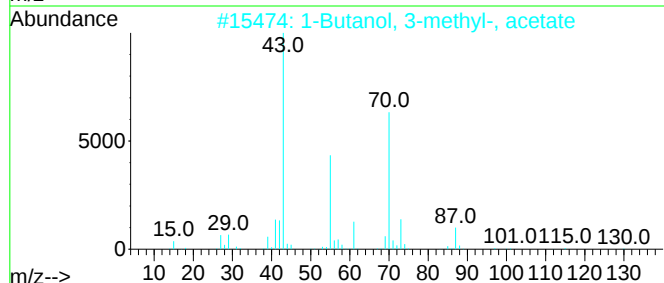
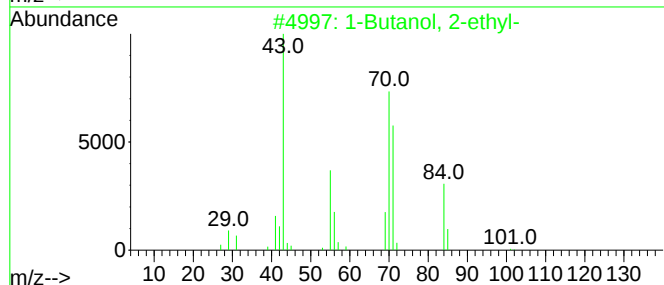
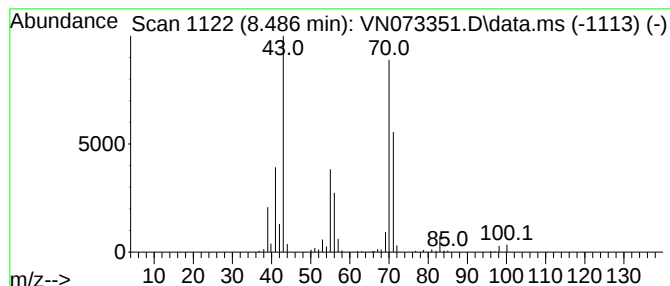
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N062822W.M  
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 1-Butanol, 2-ethyl- Concentration Rank 3

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 8.486 | 25.96 ug/l | 664462 | 1,4-Difluorobenzene | 8.904 |

| Hit# | of                            | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|-------------------------------|---|--------------|-----|---------|-------------|------|
| 1    | 1-Butanol, 2-ethyl-           |   |              | 102 | C6H14O  | 000097-95-0 | 64   |
| 2    | 1-Butanol, 3-methyl-, acetate |   |              | 130 | C7H14O2 | 000123-92-2 | 53   |
| 3    | Heptane, 3,4-dimethyl-        |   |              | 128 | C9H20   | 000922-28-1 | 50   |
| 4    | Hexane, 3-methyl-             |   |              | 100 | C7H16   | 000589-34-4 | 49   |
| 5    | Pentane, 3-ethyl-             |   |              | 100 | C7H16   | 000617-78-7 | 47   |



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ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_N  
ClientSampleId :  
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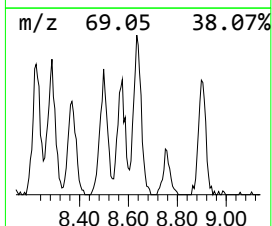
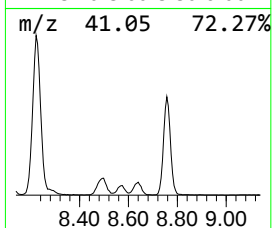
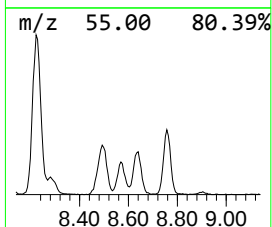
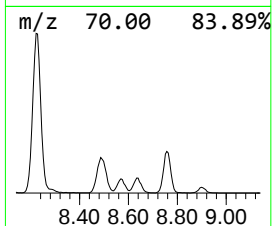
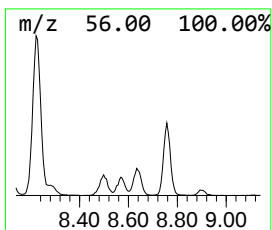
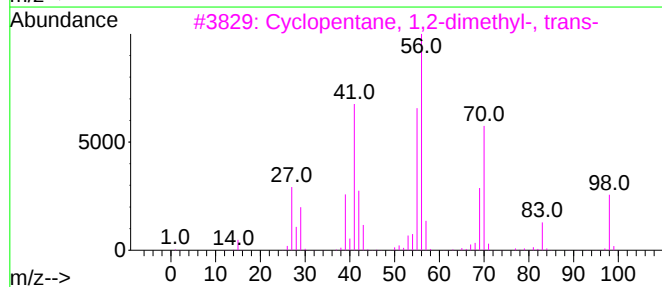
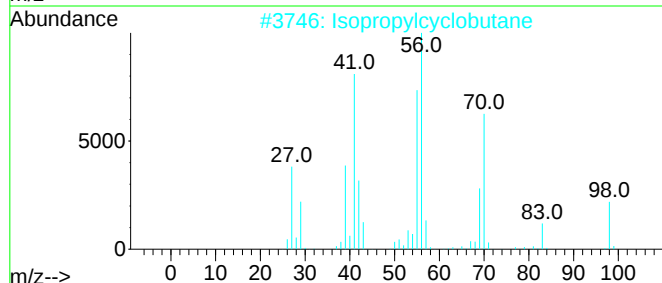
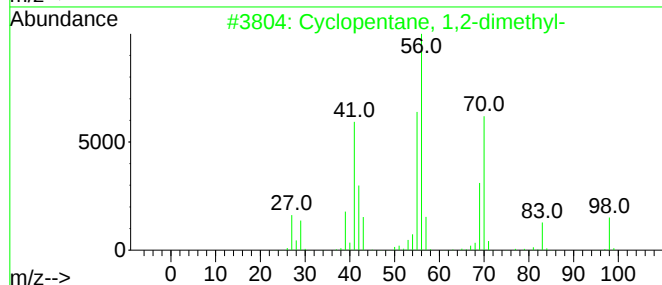
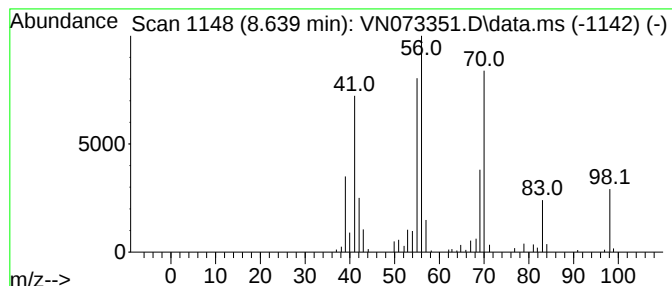
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N062822W.M  
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Cyclopentane, 1,2-dimethyl- Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 8.639 | 12.82 ug/l | 328225 | 1,4-Difluorobenzene | 8.904 |

| Hit# | of | 5 | Tentative ID                        | MW | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|----|---------|-------------|------|
| 1    |    |   | Cyclopentane, 1,2-dimethyl-         | 98 | C7H14   | 002452-99-5 | 91   |
| 2    |    |   | Isopropylcyclobutane                | 98 | C7H14   | 000872-56-0 | 91   |
| 3    |    |   | Cyclopentane, 1,2-dimethyl-, trans- | 98 | C7H14   | 000822-50-4 | 91   |
| 4    |    |   | Cyclopentane, 1,3-dimethyl-, cis-   | 98 | C7H14   | 002532-58-3 | 90   |
| 5    |    |   | Cyclopentane, 1,2-dimethyl-, cis-   | 98 | C7H14   | 001192-18-3 | 87   |



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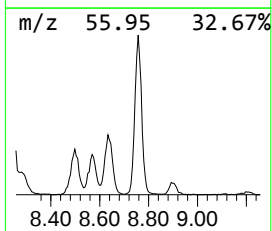
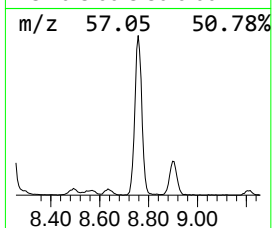
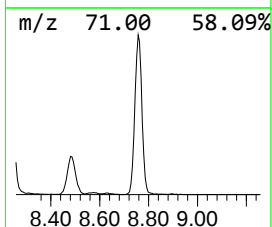
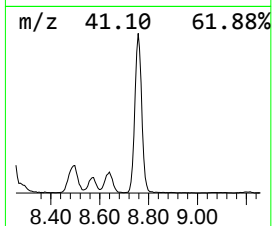
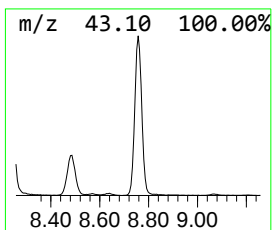
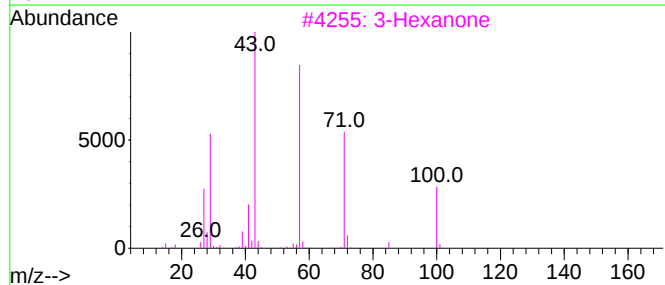
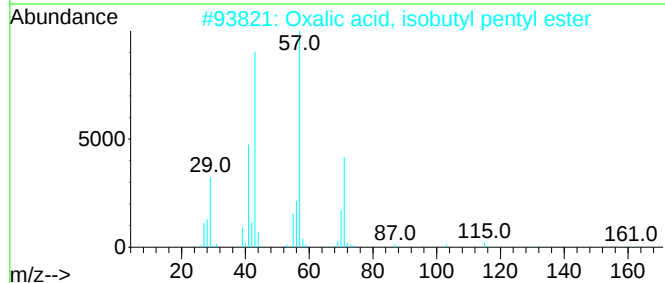
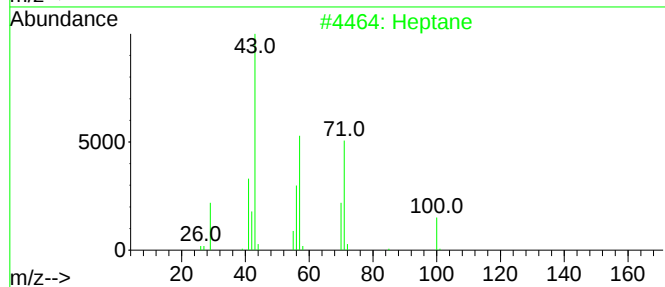
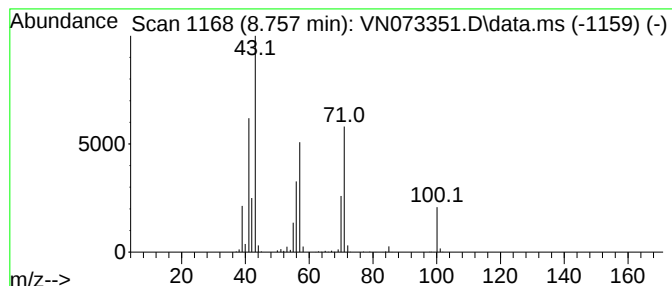
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 Heptane Concentration Rank 1

| R.T.  | EstConc    | Area    | Relative to ISTD    | R.T.  |
|-------|------------|---------|---------------------|-------|
| 8.757 | 76.80 ug/l | 1965660 | 1,4-Difluorobenzene | 8.904 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|----|---|------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Heptane                            | 100 | C7H16    | 000142-82-5  | 91   |
| 2    |    |   | Oxalic acid, isobutyl pentyl ester | 216 | C11H20O4 | 1000309-37-0 | 53   |
| 3    |    |   | 3-Hexanone                         | 100 | C6H12O   | 000589-38-8  | 46   |
| 4    |    |   | Acetaldehyde, propylhydrazone      | 100 | C5H12N2  | 007422-88-0  | 38   |
| 5    |    |   | Butane, 2,2-dimethyl-              | 86  | C6H14    | 000075-83-2  | 37   |



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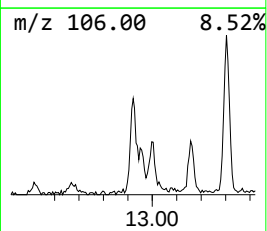
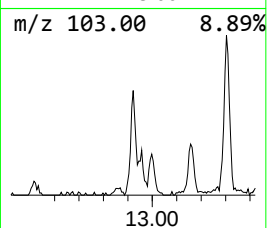
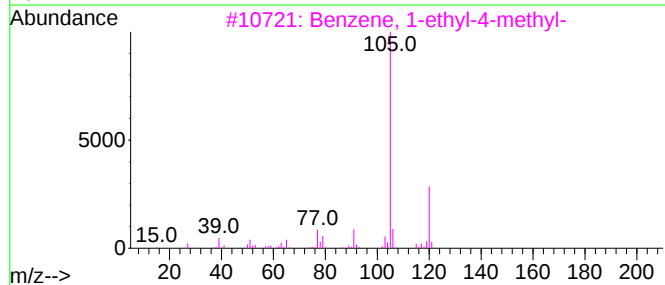
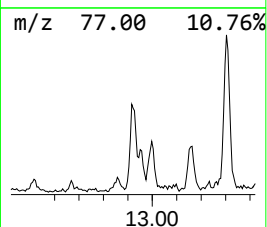
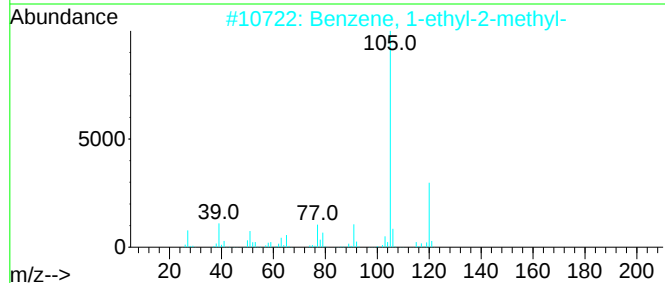
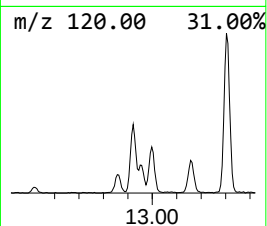
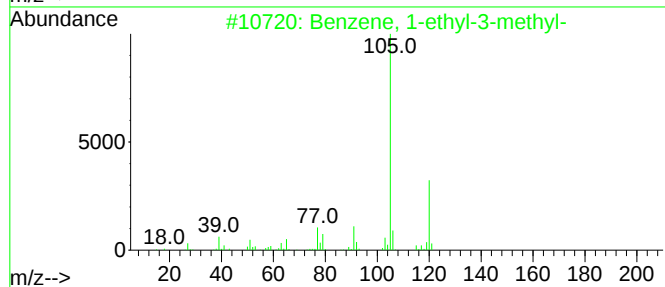
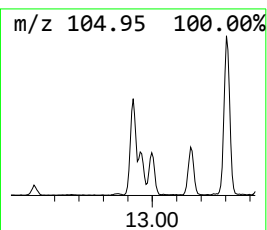
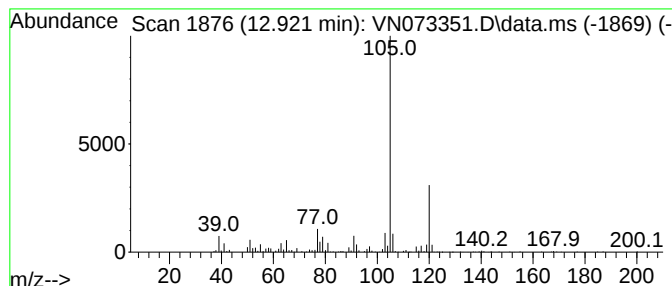
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 10

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 12.922 | 12.39 ug/l | 444153 | 1,4-Dichlorobenzene-d4 | 13.610 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 94   |
| 2    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 3    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |
| 4    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 90   |
| 5    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 90   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN070622\  
Data File : VN073351.D  
Acq On : 06 Jul 2022 17:38  
Operator : JC\MD  
Sample : N3586-04 10X  
Misc : 1.11g/10mL/100uL/5.00mL/MSVOA\_N/MEOH  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_N  
ClientSampleId :  
31457

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N062822W.M  
Quant Title : SW846 8260

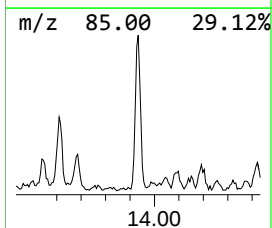
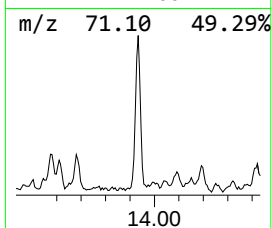
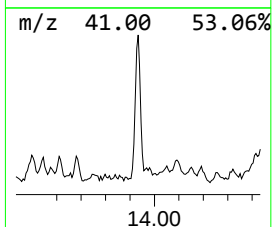
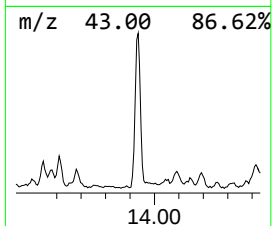
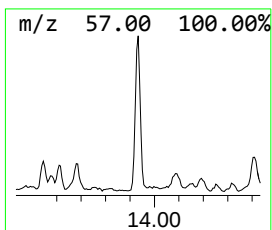
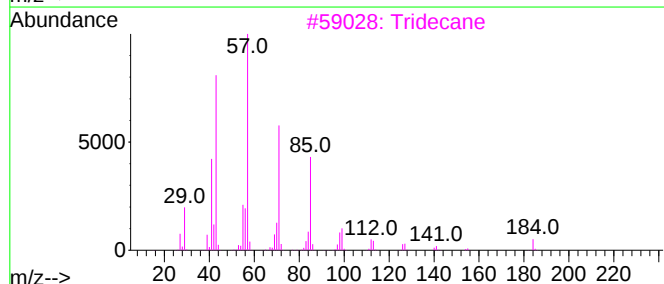
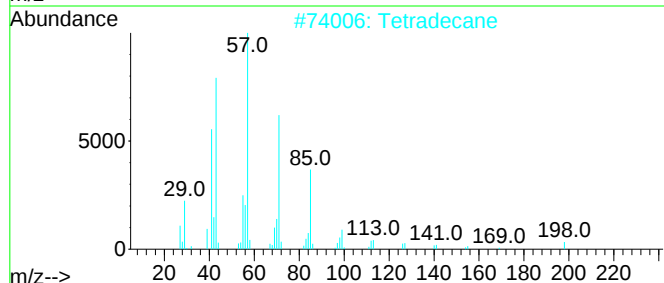
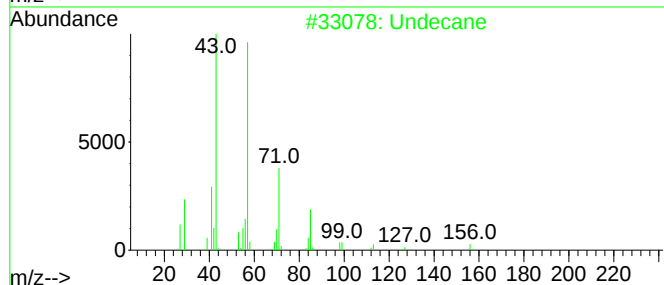
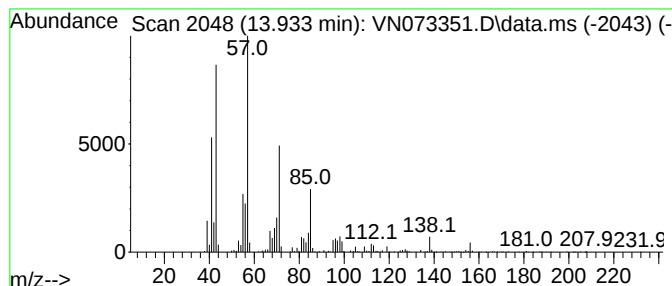
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Undecane Concentration Rank 4

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 13.933 | 19.92 ug/l | 713816 | 1,4-Dichlorobenzene-d4 | 13.610 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Undecane                  | 156 | C11H24  | 001120-21-4 | 90   |
| 2    |    |   | Tetradecane               | 198 | C14H30  | 000629-59-4 | 80   |
| 3    |    |   | Tridecane                 | 184 | C13H28  | 000629-50-5 | 80   |
| 4    |    |   | Decane, 6-ethyl-2-methyl- | 184 | C13H28  | 062108-21-8 | 59   |
| 5    |    |   | Dodecane, 1-iodo-         | 296 | C12H25I | 004292-19-7 | 58   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN070622\  
Data File : VN073351.D  
Acq On : 06 Jul 2022 17:38  
Operator : JC\MD  
Sample : N3586-04 10X  
Misc : 1.11g/10mL/100uL/5.00mL/MSVOA\_N/MEOH  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_N  
ClientSampleId :  
31457

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N062822W.M  
Quant Title : SW846 8260

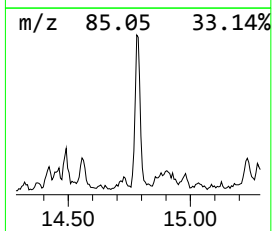
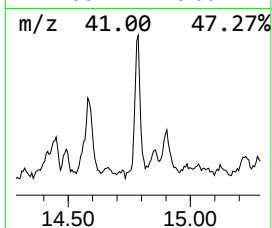
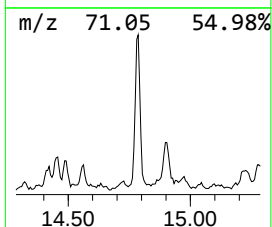
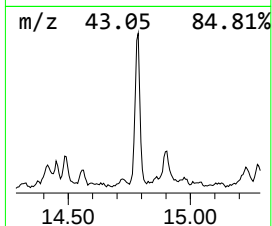
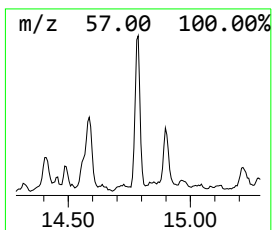
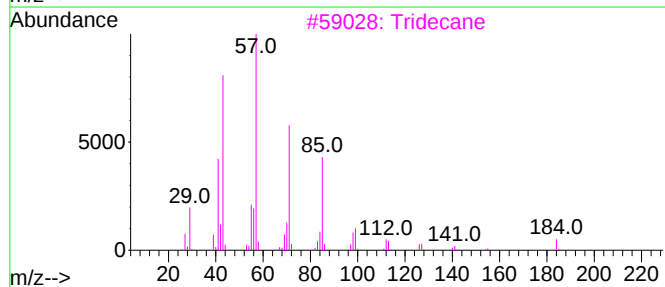
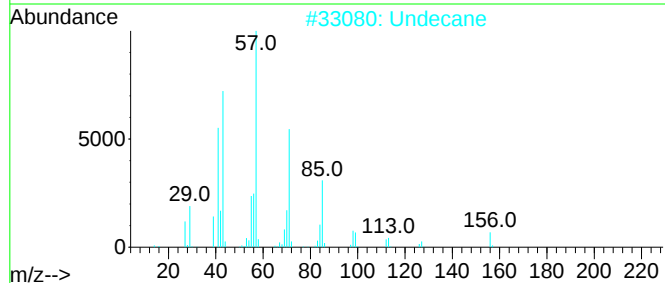
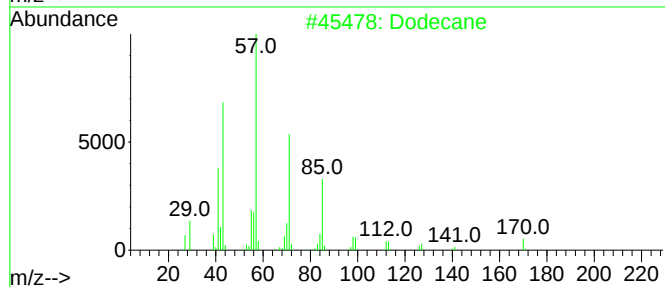
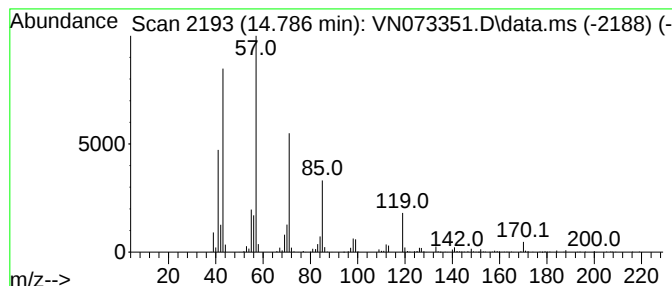
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Dodecane Concentration Rank 5

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 14.786 | 16.65 ug/l | 596764 | 1,4-Dichlorobenzene-d4 | 13.610 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane     | 170 | C12H26  | 000112-40-3 | 93   |
| 2    |    |   | Undecane     | 156 | C11H24  | 001120-21-4 | 80   |
| 3    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 80   |
| 4    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 72   |
| 5    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 72   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN070622\  
Data File : VN073351.D  
Acq On : 06 Jul 2022 17:38  
Operator : JC\MD  
Sample : N3586-04 10X  
Misc : 1.11g/10mL/100uL/5.00mL/MSVOA\_N/MEOH  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_N  
ClientSampleId :  
31457

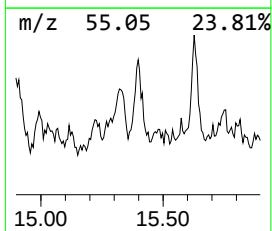
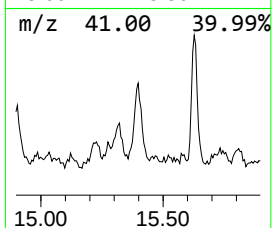
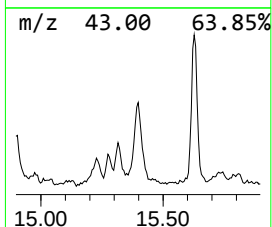
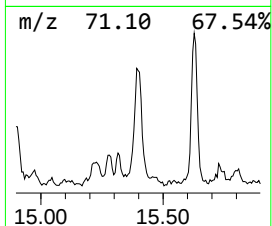
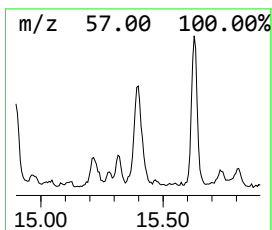
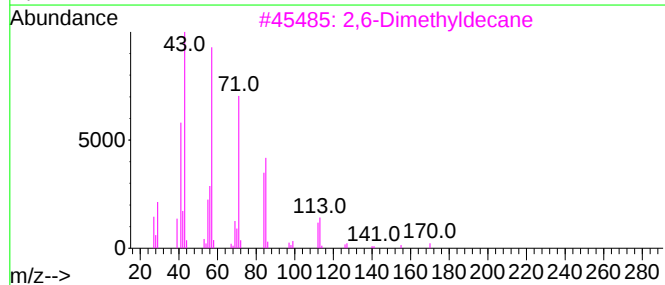
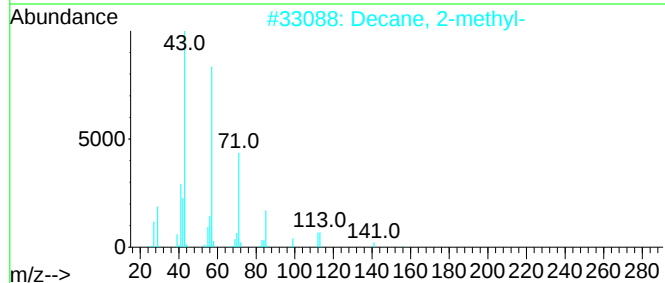
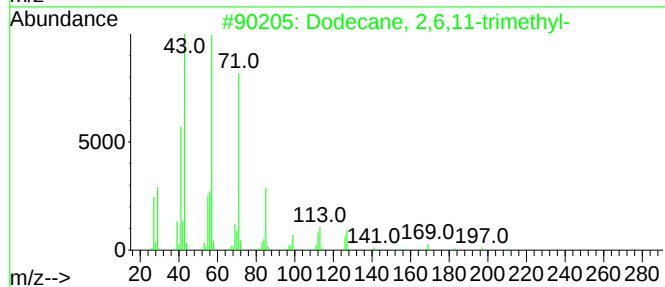
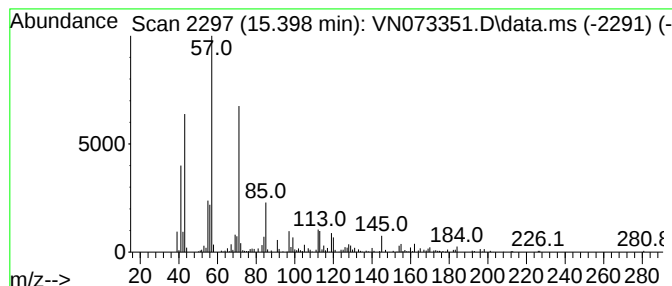
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Dodecane, 2,6,11-trimethyl- Concentration Rank 7

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 15.398 | 13.33 ug/l | 477615 | 1,4-Dichlorobenzene-d4 | 13.610 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 90   |
| 2    |    |   | Decane, 2-methyl-           | 156 | C11H24  | 006975-98-0 | 87   |
| 3    |    |   | 2,6-Dimethyldecane          | 170 | C12H26  | 013150-81-7 | 81   |
| 4    |    |   | Pentadecane, 7-methyl-      | 226 | C16H34  | 006165-40-8 | 76   |
| 5    |    |   | 2,6,10-Trimethyltridecane   | 226 | C16H34  | 003891-99-4 | 62   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN070622\  
Data File : VN073351.D  
Acq On : 06 Jul 2022 17:38  
Operator : JC\MD  
Sample : N3586-04 10X  
Misc : 1.11g/10mL/100uL/5.00mL/MSVOA\_N/MEOH  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_N  
ClientSampleId :  
31457

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N062822W.M  
Quant Title : SW846 8260

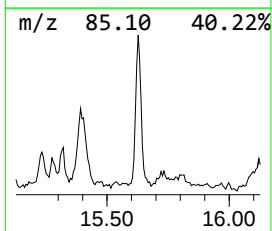
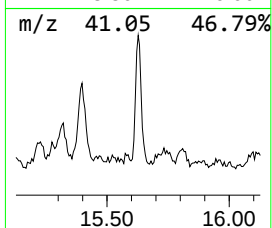
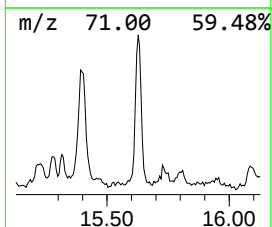
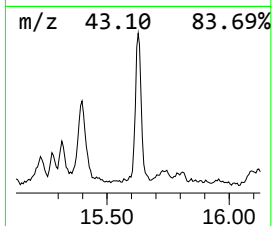
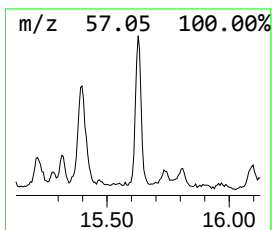
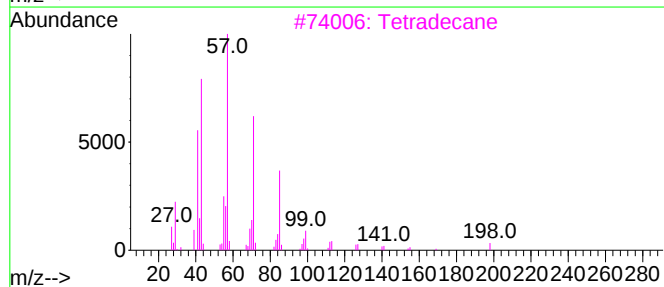
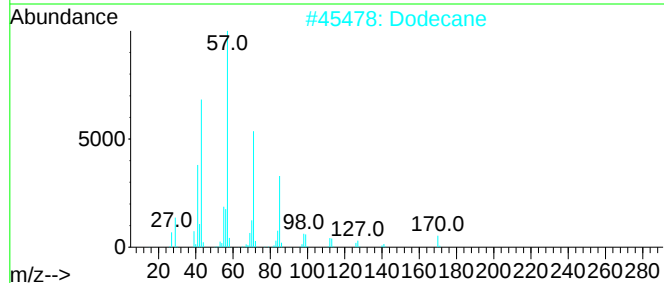
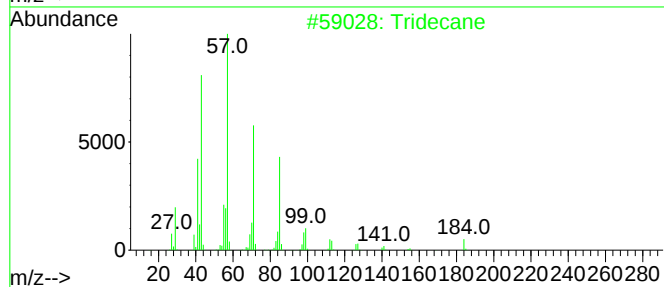
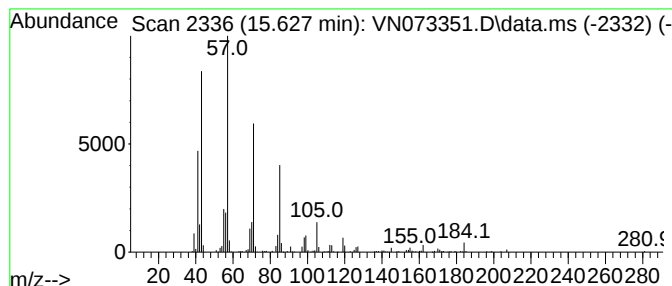
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Tridecane Concentration Rank 6

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 15.627 | 15.64 ug/l | 560632 | 1,4-Dichlorobenzene-d4 | 13.610 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane                | 184 | C13H28  | 000629-50-5 | 96   |
| 2    |    |   | Dodecane                 | 170 | C12H26  | 000112-40-3 | 74   |
| 3    |    |   | Tetradecane              | 198 | C14H30  | 000629-59-4 | 74   |
| 4    |    |   | Octane, 2,4,6-trimethyl- | 156 | C11H24  | 062016-37-9 | 72   |
| 5    |    |   | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 72   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN070622\  
Data File : VN073351.D  
Acq On : 06 Jul 2022 17:38  
Operator : JC\MD  
Sample : N3586-04 10X  
Misc : 1.11g/10mL/100uL/5.00mL/MSVOA\_N/MEOH  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_N  
ClientSampleId :  
31457

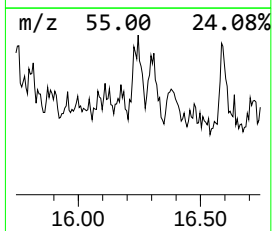
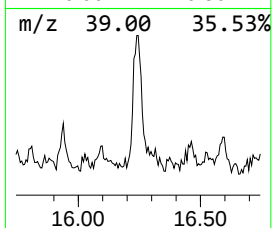
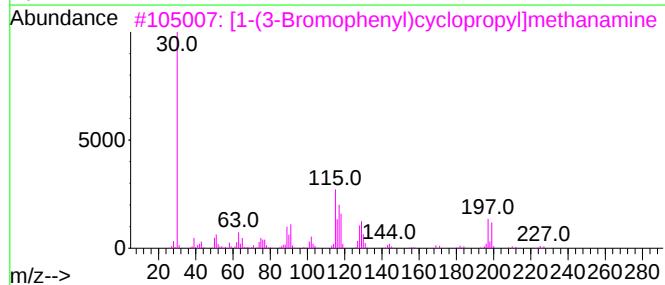
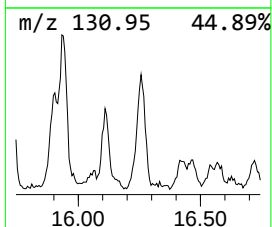
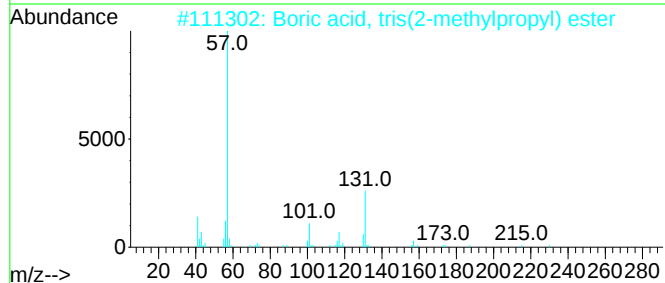
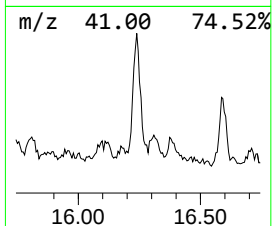
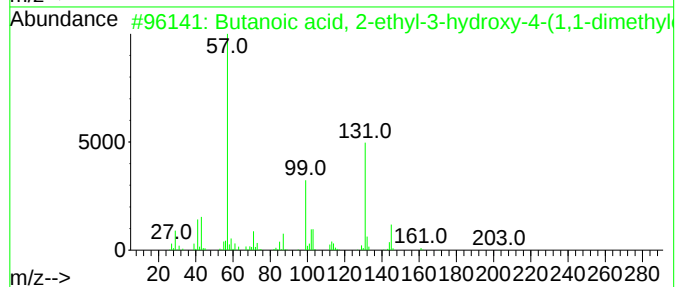
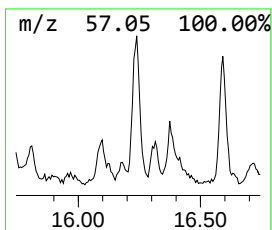
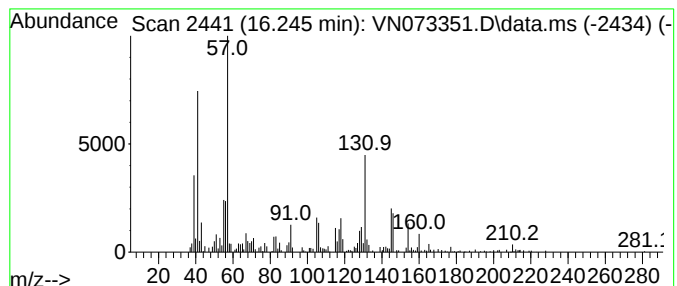
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 unknown16.245 Concentration Rank 8

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 16.245 | 13.17 ug/l | 471862 | 1,4-Dichlorobenzene-d4 | 13.610 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Butanoic acid, 2-ethyl-3-hydroxy... | 218 | C11H22O4  | 106059-00-1  | 25   |
| 2    |    |   | Boric acid, tris(2-methylpropyl)... | 230 | C12H27BO3 | 013195-76-1  | 12   |
| 3    |    |   | [1-(3-Bromophenyl)cyclopropyl]me... | 225 | C10H12BrN | 886365-87-3  | 10   |
| 4    |    |   | 1,1,1,2,3,3,3-Heptafluoro-2-meth... | 200 | C4H3F7O   | 1000322-37-6 | 9    |
| 5    |    |   | Iodothioacetic acid, S-t-butyl e... | 258 | C6H11IOS  | 1010192-69-7 | 9    |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_N\Data\VN070622\  
Data File : VN073351.D  
Acq On : 06 Jul 2022 17:38  
Operator : JC\MD  
Sample : N3586-04 10X  
Misc : 1.11g/10mL/100uL/5.00mL/MSVOA\_N/MEOH  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_N  
ClientSampleId :  
31457

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_N\methods\82N062822W.M  
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| Hexane, 3-methyl-  | 8.222  | 45.3    | ug/l  | 4429960  | 1                     | 8.016  | 4891110 | 50.0 |
| 1-Butanol, 2-et... | 8.486  | 26.0    | ug/l  | 664462   | 2                     | 8.904  | 1279750 | 50.0 |
| Cyclopentane, 1... | 8.639  | 12.8    | ug/l  | 328225   | 2                     | 8.904  | 1279750 | 50.0 |
| Heptane            | 8.757  | 76.8    | ug/l  | 1965660  | 2                     | 8.904  | 1279750 | 50.0 |
| Benzene, 1-ethy... | 12.922 | 12.4    | ug/l  | 444153   | 4                     | 13.610 | 1791760 | 50.0 |
| Undecane           | 13.933 | 19.9    | ug/l  | 713816   | 4                     | 13.610 | 1791760 | 50.0 |
| Dodecane           | 14.786 | 16.6    | ug/l  | 596764   | 4                     | 13.610 | 1791760 | 50.0 |
| Dodecane, 2,6,1... | 15.398 | 13.3    | ug/l  | 477615   | 4                     | 13.610 | 1791760 | 50.0 |
| Tridecane          | 15.627 | 15.6    | ug/l  | 560632   | 4                     | 13.610 | 1791760 | 50.0 |
| unknown16.245      | 16.245 | 13.2    | ug/l  | 471862   | 4                     | 13.610 | 1791760 | 50.0 |