

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
 Data File : VU028145.D  
 Acq On : 15 Nov 2018 01:42  
 Operator : MD/SY  
 Sample : J5943-07  
 Misc : 25.0mL/MSVOA U/WATER  
 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 H0FH6

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\SOMUTR111418WMA.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         | 1.402       | 58            | 71          | 78           | rBV2     | 1304992        | 1860869       | 8.95%           | 2.040%        |
| 2         | 1.517       | 100           | 107         | 120          | rVB      | 287683         | 330132        | 1.59%           | 0.362%        |
| 3         | 1.652       | 133           | 149         | 156          | rBV2     | 337454         | 553855        | 2.66%           | 0.607%        |
| 4         | 2.096       | 279           | 287         | 293          | rBV      | 153521         | 228245        | 1.10%           | 0.250%        |
| 5         | 2.984       | 548           | 563         | 589          | rBV      | 735412         | 1353039       | 6.51%           | 1.483%        |
| 6         | 3.447       | 685           | 707         | 733          | rBV      | 385460         | 916983        | 4.41%           | 1.005%        |
| 7         | 4.228       | 932           | 950         | 1010         | rVB      | 1158570        | 2952737       | 14.21%          | 3.236%        |
| 8         | 4.652       | 1065          | 1082        | 1104         | rBV2     | 89675          | 217765        | 1.05%           | 0.239%        |
| 9         | 5.340       | 1274          | 1296        | 1302         | rBV2     | 194147         | 502190        | 2.42%           | 0.550%        |
| 10        | 5.389       | 1302          | 1311        | 1334         | rVB      | 472149         | 1005341       | 4.84%           | 1.102%        |
| 11        | 5.887       | 1454          | 1466        | 1482         | rBV      | 163605         | 318870        | 1.53%           | 0.349%        |
| 12        | 6.215       | 1542          | 1568        | 1583         | rBV      | 629684         | 1245983       | 5.99%           | 1.366%        |
| 13        | 6.334       | 1598          | 1605        | 1619         | rVB2     | 123172         | 231384        | 1.11%           | 0.254%        |
| 14        | 7.565       | 1969          | 1988        | 1997         | rBV      | 246236         | 440542        | 2.12%           | 0.483%        |
| 15        | 7.639       | 1997          | 2011        | 2046         | rVB      | 12515762       | 20785743      | 100.00%         | 22.782%       |
| 16        | 8.318       | 2210          | 2222        | 2256         | rBV      | 364672         | 694935        | 3.34%           | 0.762%        |
| 17        | 9.089       | 2452          | 2462        | 2484         | rVB2     | 253562         | 557502        | 2.68%           | 0.611%        |
| 18        | 9.250       | 2499          | 2512        | 2535         | rBV      | 4340430        | 6698299       | 32.23%          | 7.342%        |
| 19        | 9.376       | 2535          | 2551        | 2568         | rBV      | 9772935        | 15568239      | 74.90%          | 17.063%       |
| 20        | 9.781       | 2663          | 2677        | 2705         | rVB      | 5093122        | 7923587       | 38.12%          | 8.685%        |
| 21        | 10.167      | 2782          | 2797        | 2808         | rBV      | 138737         | 238086        | 1.15%           | 0.261%        |
| 22        | 10.250      | 2808          | 2823        | 2855         | rVB      | 274020         | 497394        | 2.39%           | 0.545%        |
| 23        | 10.588      | 2914          | 2928        | 2940         | rBV      | 155280         | 248414        | 1.20%           | 0.272%        |
| 24        | 10.681      | 2945          | 2957        | 2963         | rBV      | 410630         | 719416        | 3.46%           | 0.789%        |
| 25        | 10.771      | 2975          | 2985        | 3007         | rVB      | 272762         | 429739        | 2.07%           | 0.471%        |
| 26        | 10.919      | 3018          | 3031        | 3039         | rBV      | 171302         | 264908        | 1.27%           | 0.290%        |
| 27        | 10.970      | 3039          | 3047        | 3059         | rVB      | 260882         | 407675        | 1.96%           | 0.447%        |
| 28        | 11.147      | 3091          | 3102        | 3119         | rBV      | 1074301        | 1604728       | 7.72%           | 1.759%        |
| 29        | 11.482      | 3196          | 3206        | 3222         | rVB2     | 298218         | 483292        | 2.33%           | 0.530%        |
| 30        | 11.568      | 3222          | 3233        | 3248         | rVB      | 736314         | 1126635       | 5.42%           | 1.235%        |
| 31        | 11.768      | 3284          | 3295        | 3302         | rBV      | 207780         | 359447        | 1.73%           | 0.394%        |
| 32        | 11.861      | 3315          | 3324        | 3352         | rVB4     | 344369         | 695103        | 3.34%           | 0.762%        |
| 33        | 12.250      | 3434          | 3445        | 3453         | rBV      | 204290         | 312730        | 1.50%           | 0.343%        |
| 34        | 12.353      | 3467          | 3477        | 3505         | rVB2     | 110386         | 271118        | 1.30%           | 0.297%        |

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Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 H0FH6

Integration Parameters: rteint.p  
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 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.1  
 Stop Thrs : 0  
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 Max Peaks: 100  
 Peak Location: TOP

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 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
 Title : TRACE VOA SOM01.0

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 12.700 | 3577 | 3585 | 3603 | rVB  | 139957  | 226039  | 1.09%  | 0.248% |
| 36 | 12.964 | 3658 | 3667 | 3680 | rVB  | 185307  | 281928  | 1.36%  | 0.309% |
| 37 | 13.070 | 3690 | 3700 | 3707 | rBV2 | 264392  | 412696  | 1.99%  | 0.452% |
| 38 | 13.121 | 3707 | 3716 | 3733 | rVB3 | 493605  | 800180  | 3.85%  | 0.877% |
| 39 | 13.289 | 3757 | 3768 | 3791 | rVB  | 647760  | 1030375 | 4.96%  | 1.129% |
| 40 | 13.742 | 3889 | 3909 | 3919 | rBV  | 708417  | 1132433 | 5.45%  | 1.241% |
| 41 | 13.855 | 3933 | 3944 | 3960 | rBV  | 336302  | 524688  | 2.52%  | 0.575% |
| 42 | 13.948 | 3960 | 3973 | 3985 | rBV  | 416526  | 684636  | 3.29%  | 0.750% |
| 43 | 14.154 | 4026 | 4037 | 4053 | rBV  | 131005  | 253260  | 1.22%  | 0.278% |
| 44 | 14.350 | 4082 | 4098 | 4117 | rBV2 | 998269  | 1688660 | 8.12%  | 1.851% |
| 45 | 14.539 | 4147 | 4157 | 4169 | rVB2 | 189742  | 359465  | 1.73%  | 0.394% |
| 46 | 14.674 | 4187 | 4199 | 4214 | rBV  | 794240  | 1266186 | 6.09%  | 1.388% |
| 47 | 14.874 | 4237 | 4261 | 4272 | rBV2 | 1710425 | 3191800 | 15.36% | 3.498% |
| 48 | 14.932 | 4272 | 4279 | 4293 | rVB4 | 265530  | 475931  | 2.29%  | 0.522% |
| 49 | 15.060 | 4310 | 4319 | 4331 | rVV  | 660606  | 1040471 | 5.01%  | 1.140% |
| 50 | 15.128 | 4331 | 4340 | 4346 | rVV2 | 180427  | 304943  | 1.47%  | 0.334% |
| 51 | 15.163 | 4346 | 4351 | 4364 | rVB2 | 136821  | 216859  | 1.04%  | 0.238% |
| 52 | 15.240 | 4364 | 4375 | 4388 | rBV5 | 206436  | 428144  | 2.06%  | 0.469% |
| 53 | 15.604 | 4466 | 4488 | 4499 | rBV4 | 631727  | 1682743 | 8.10%  | 1.844% |
| 54 | 15.800 | 4539 | 4549 | 4570 | rVB3 | 267151  | 558141  | 2.69%  | 0.612% |
| 55 | 15.912 | 4571 | 4584 | 4594 | rBV2 | 545697  | 995083  | 4.79%  | 1.091% |
| 56 | 16.057 | 4621 | 4629 | 4636 | rBV  | 433130  | 717103  | 3.45%  | 0.786% |
| 57 | 16.096 | 4636 | 4641 | 4660 | rVB  | 432806  | 664938  | 3.20%  | 0.729% |
| 58 | 16.282 | 4686 | 4699 | 4715 | rVB2 | 106569  | 255735  | 1.23%  | 0.280% |

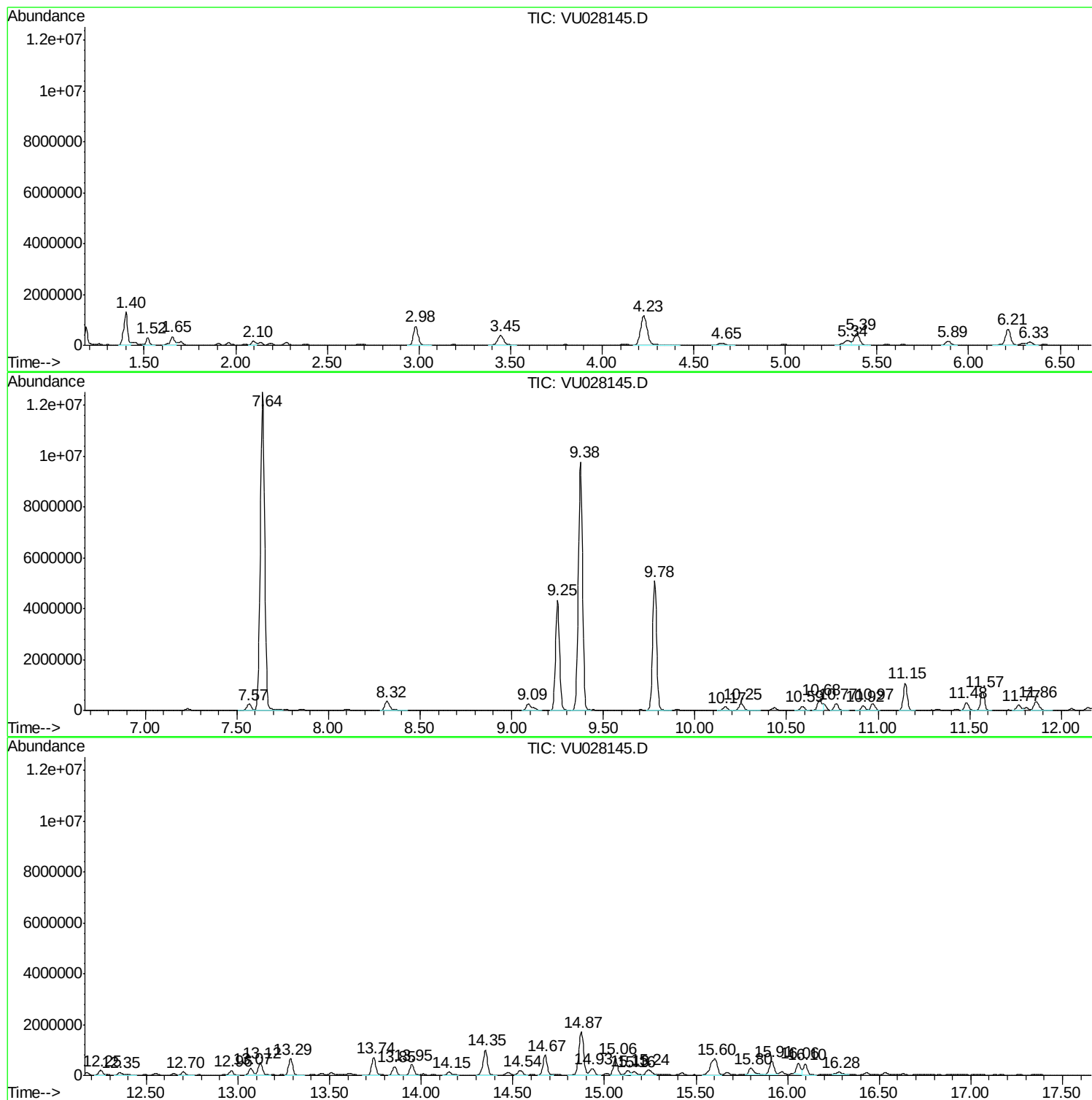
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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



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Quant Title : TRACE VOA SOM01.0

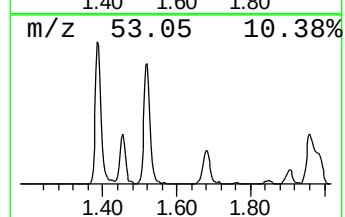
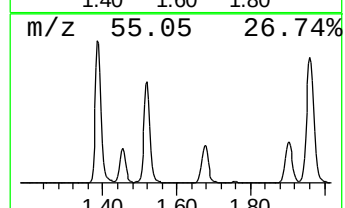
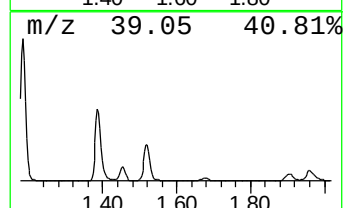
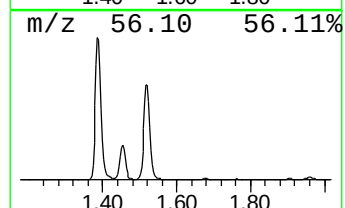
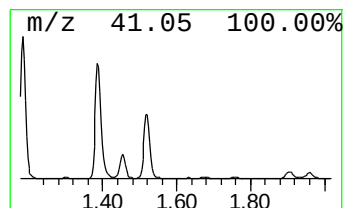
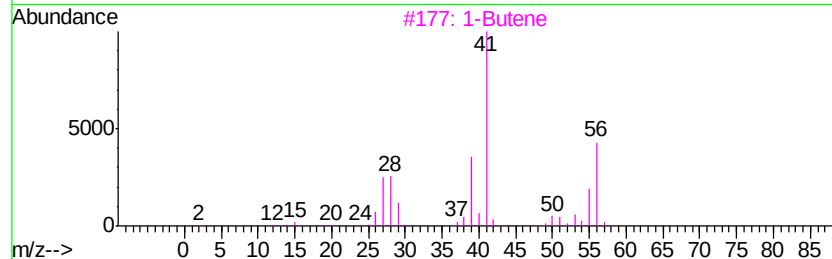
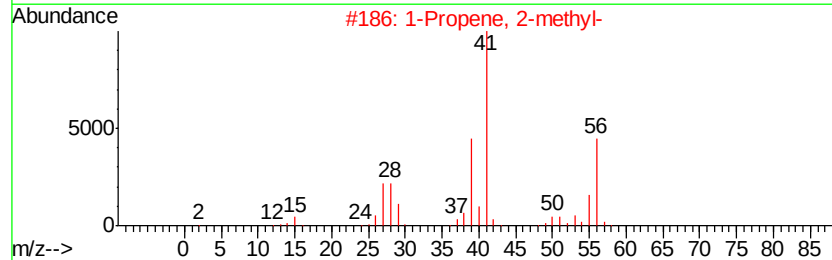
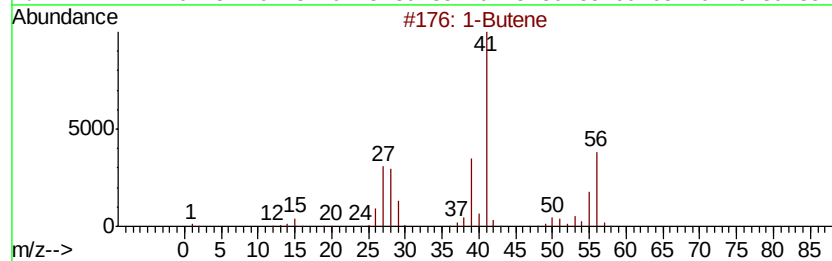
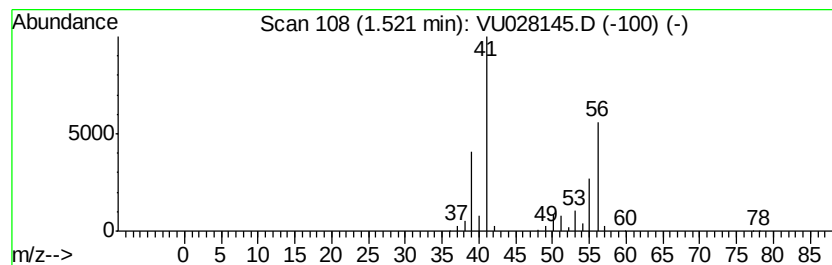
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 (DEL) Alkane: Cyclic1.52 Concentration Rank 18

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 1.52 | 5.18 ug/L | 330132 | 1,4-Difluorobenzene | 5.89 |

| Hit# | of | Tentative ID         | MW | MolForm | CAS#        | Qual |
|------|----|----------------------|----|---------|-------------|------|
| 1    | 5  | 1-Butene             | 56 | C4H8    | 000106-98-9 | 87   |
| 2    |    | 1-Propene, 2-methyl- | 56 | C4H8    | 000115-11-7 | 87   |
| 3    |    | 1-Butene             | 56 | C4H8    | 000106-98-9 | 86   |
| 4    |    | 1-Butene             | 56 | C4H8    | 000106-98-9 | 86   |
| 5    |    | 2-Butene, (Z)-       | 56 | C4H8    | 000590-18-1 | 80   |



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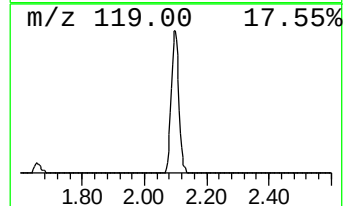
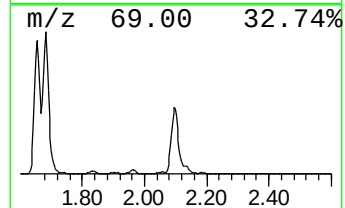
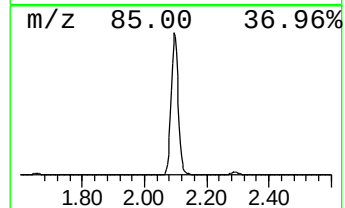
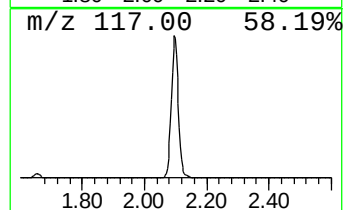
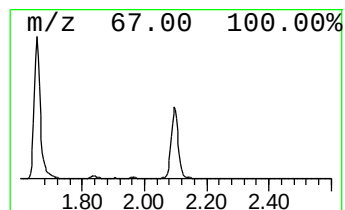
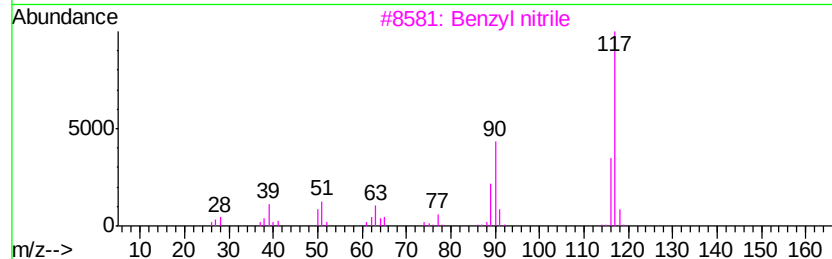
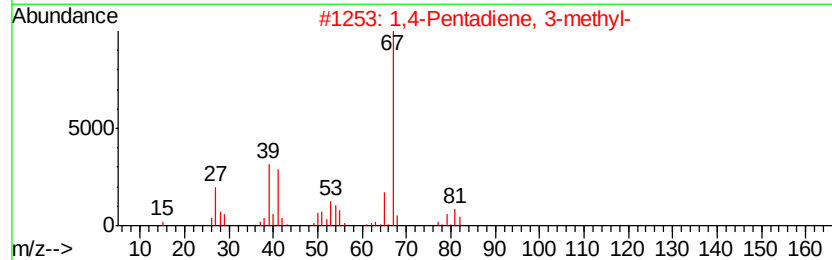
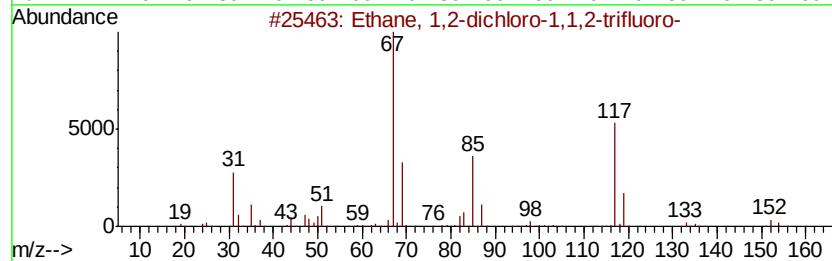
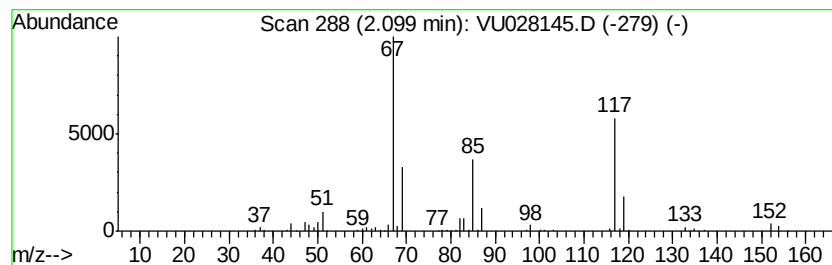
Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 2 Ethane, 1,2-dichloro-1,1,2-... Concentration Rank 27

| R.T.    | EstConc   | Area                                | Relative to ISTD    | R.T.           |
|---------|-----------|-------------------------------------|---------------------|----------------|
| 2.10    | 3.58 ug/L | 228245                              | 1,4-Difluorobenzene | 5.89           |
| Hit# of | 5         | Tentative ID                        | MW MolForm          | CAS# Qual      |
| 1       |           | Ethane, 1,2-dichloro-1,1,2-trifl... | 152 C2HCl2F3        | 000354-23-4 97 |
| 2       |           | 1,4-Pentadiene, 3-methyl-           | 82 C6H10            | 001115-08-8 9  |
| 3       |           | Benzyl nitrile                      | 117 C8H7N           | 000140-29-4 9  |
| 4       |           | 1-Butyne, 3,3-dimethyl-             | 82 C6H10            | 000917-92-0 9  |
| 5       |           | Cyclopentene, 1-methyl-             | 82 C6H10            | 000693-89-0 9  |



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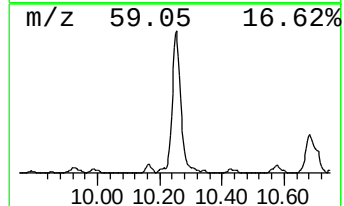
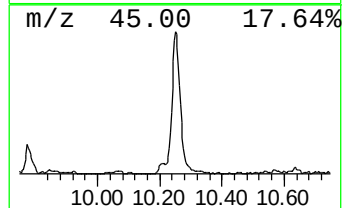
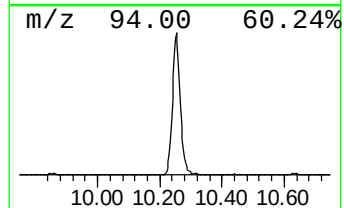
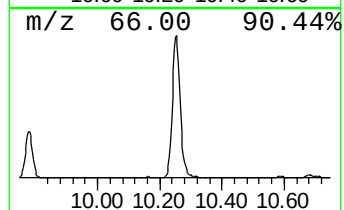
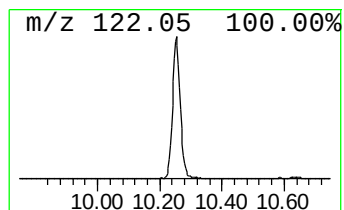
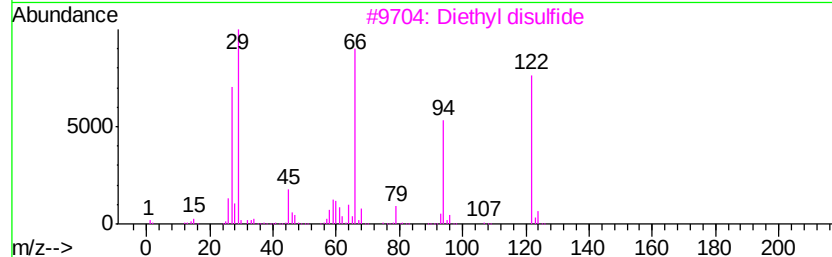
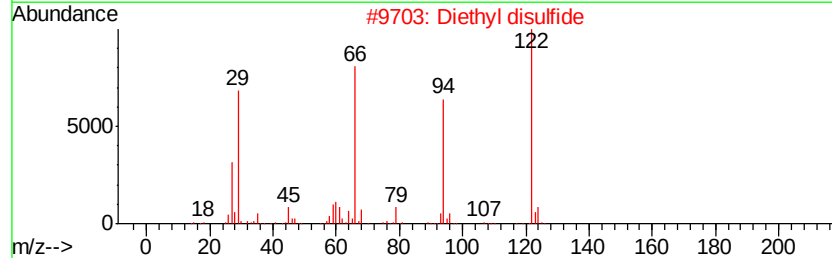
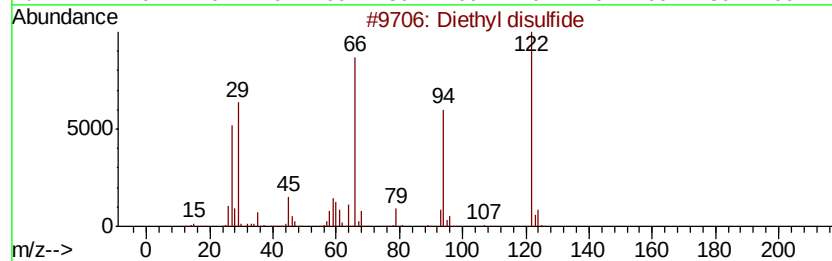
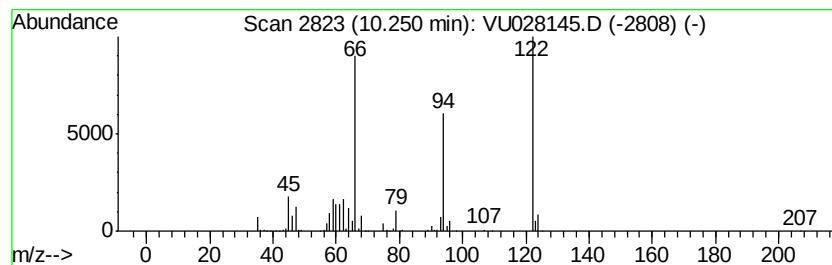
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Diethyl disulfide Concentration Rank 20

| R.T.  | EstConc   | Area   | Relative to ISTD | R.T. |
|-------|-----------|--------|------------------|------|
| 10.25 | 4.46 ug/L | 497394 | Chlorobenzene-d5 | 9.09 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Diethyl disulfide              | 122 | C4H10S2 | 000110-81-6 | 96   |
| 2    |    | Diethyl disulfide              | 122 | C4H10S2 | 000110-81-6 | 94   |
| 3    |    | Diethyl disulfide              | 122 | C4H10S2 | 000110-81-6 | 93   |
| 4    |    | 5-exo-Methyl-5-norbornenol     | 124 | C8H12O  | 003212-13-3 | 47   |
| 5    |    | 6-exo-Methyl-5-exo-norbornenol | 124 | C8H12O  | 060362-83-6 | 47   |



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Quant Title : TRACE VOA SOM01.0

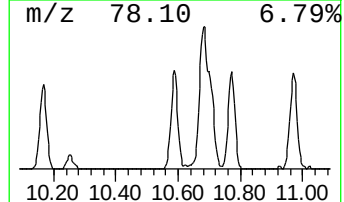
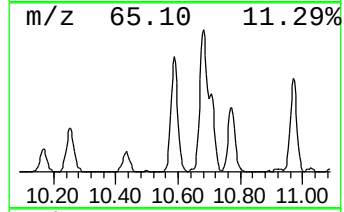
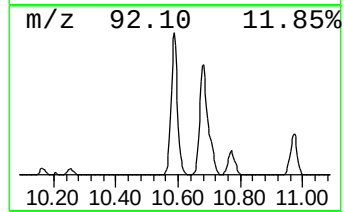
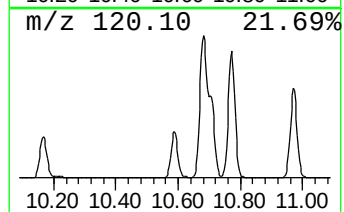
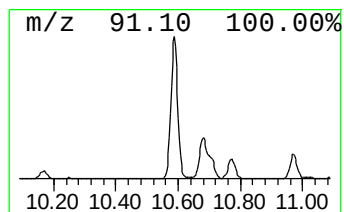
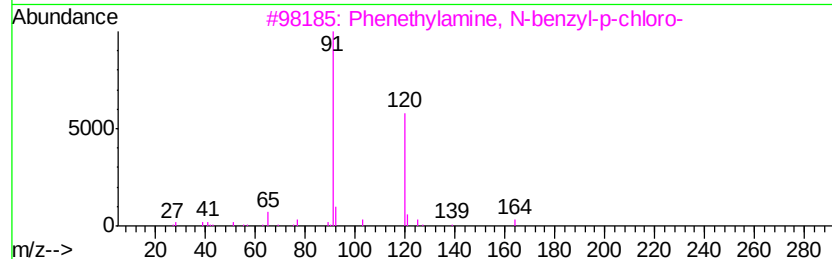
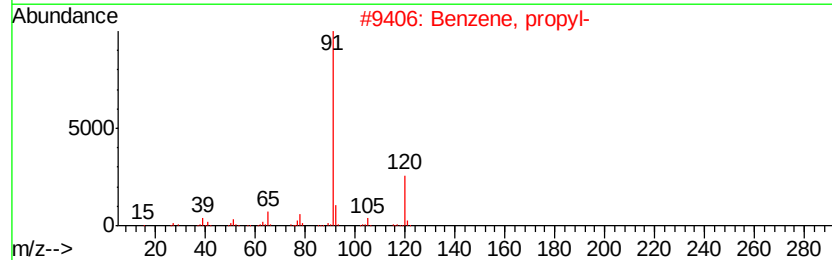
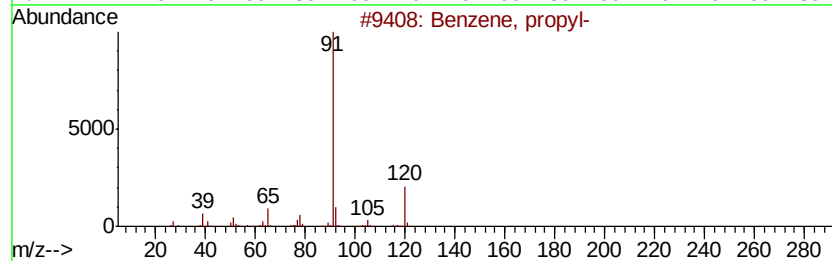
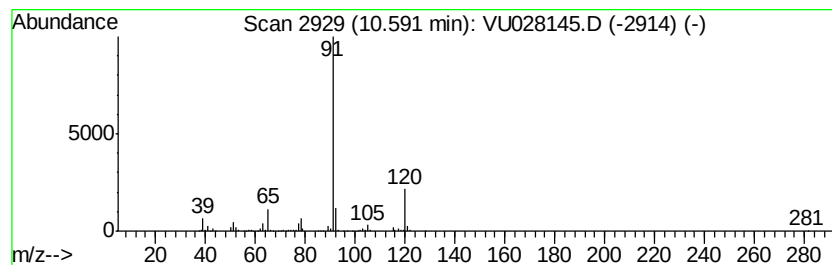
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Benzene, propyl- Concentration Rank 35

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 10.59 | 2.57 ug/L | 248414 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Benzene, propyl-                    | 120 | C9H12     | 000103-65-1 | 90   |
| 2    |    | Benzene, propyl-                    | 120 | C9H12     | 000103-65-1 | 87   |
| 3    |    | Phenethylamine, N-benzyl-p-chloro-  | 245 | C15H16ClN | 013622-43-0 | 72   |
| 4    |    | Ethanol, 2-[(phenylmethyl)amino]-   | 151 | C9H13NO   | 000104-63-2 | 64   |
| 5    |    | 1,2-Ethanediamine, N,N'-bis(phen... | 240 | C16H20N2  | 000140-28-3 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

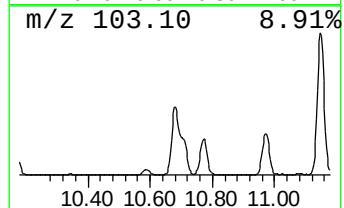
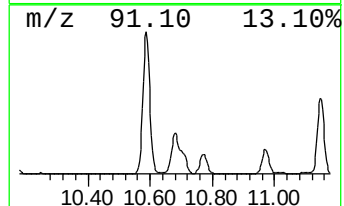
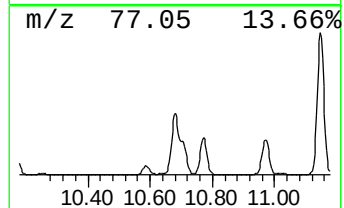
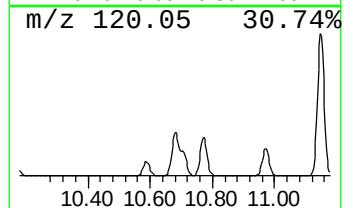
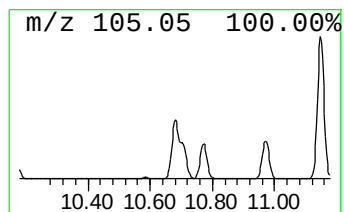
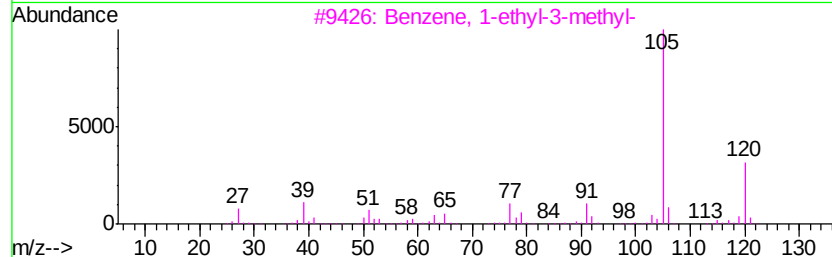
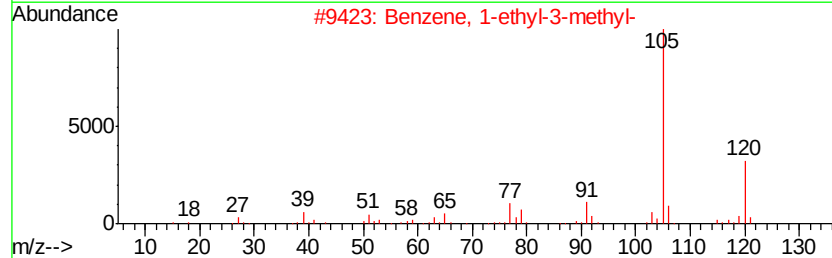
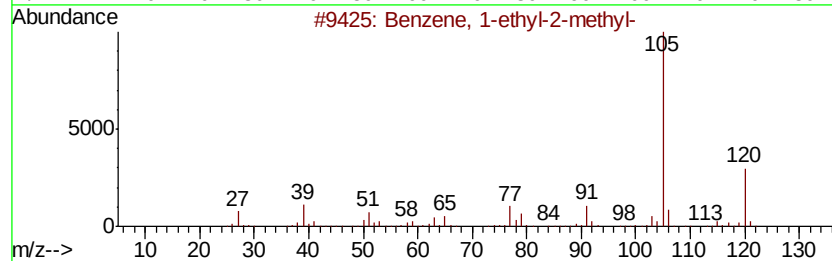
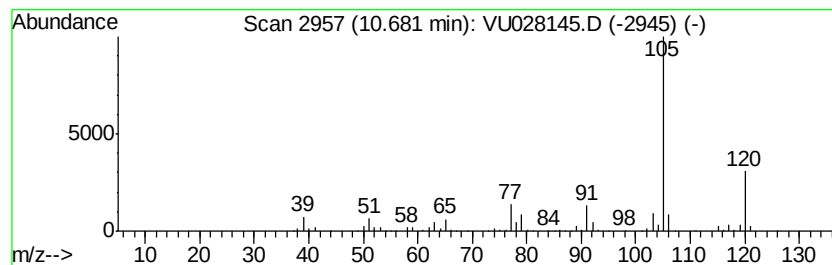
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Quant Title : TRACE VOA SOM01.0

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

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Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 12

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 10.68 | 7.44 ug/L | 719416 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 2    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 95   |
| 3    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 94   |
| 4    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |
| 5    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 91   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

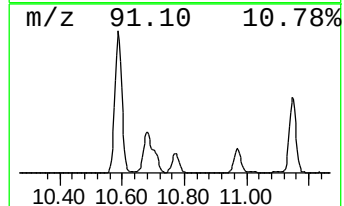
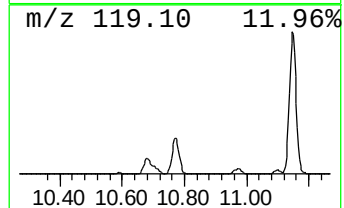
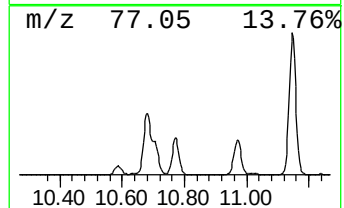
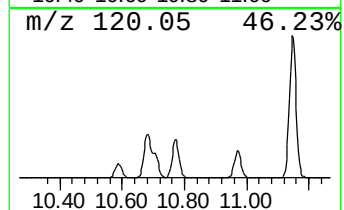
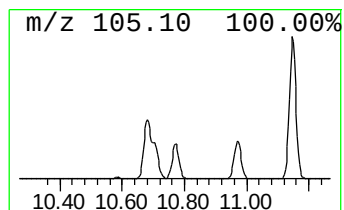
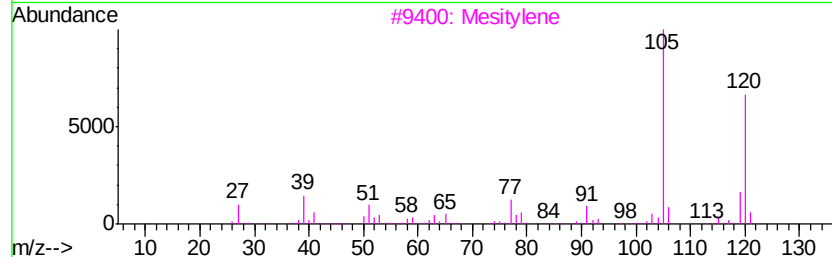
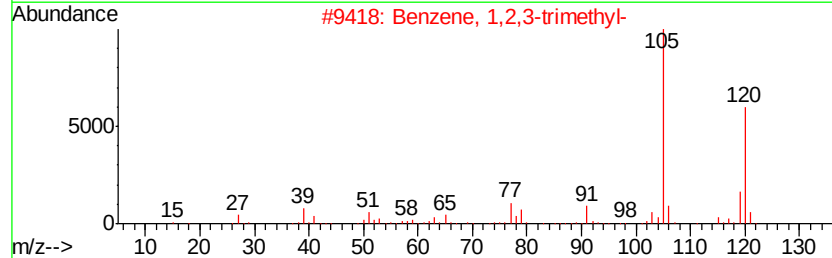
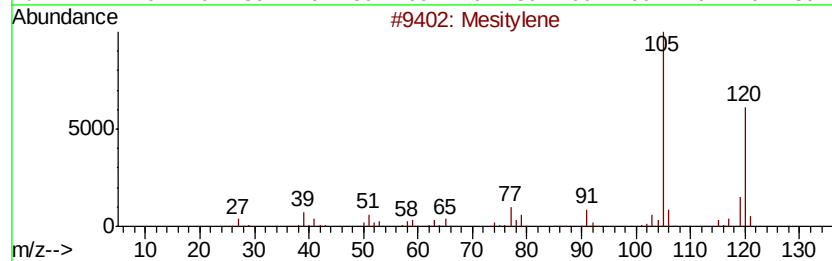
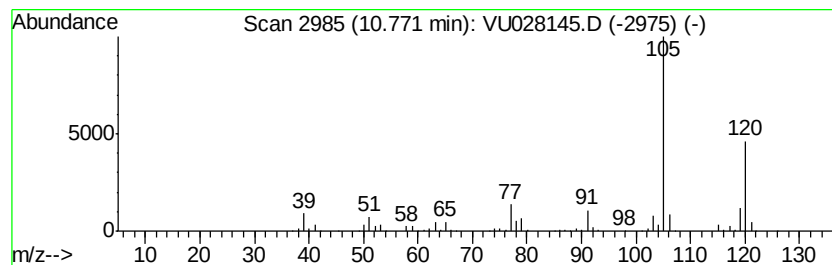
Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 6 Mesitylene Concentration Rank 21

| R.T.      | EstConc                   | Area   | Relative to ISTD       | R.T.        |      |
|-----------|---------------------------|--------|------------------------|-------------|------|
| 10.77     | 4.45 ug/L                 | 429739 | 1,4-Dichlorobenzene-d4 | 11.48       |      |
| Hit# of 5 | Tentative ID              | MW     | MolForm                | CAS#        | Qual |
| 1         | Mesitylene                | 120    | C9H12                  | 000108-67-8 | 97   |
| 2         | Benzene, 1,2,3-trimethyl- | 120    | C9H12                  | 000526-73-8 | 97   |
| 3         | Mesitylene                | 120    | C9H12                  | 000108-67-8 | 95   |
| 4         | Mesitylene                | 120    | C9H12                  | 000108-67-8 | 95   |
| 5         | Benzene, 1,2,4-trimethyl- | 120    | C9H12                  | 000095-63-6 | 95   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

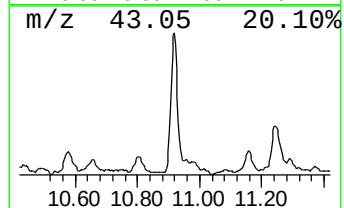
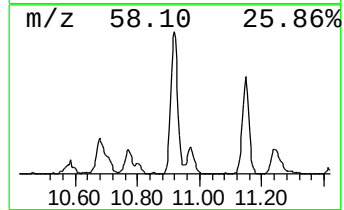
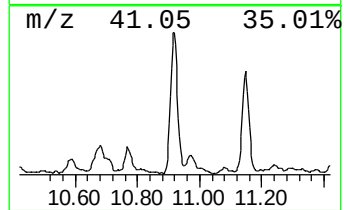
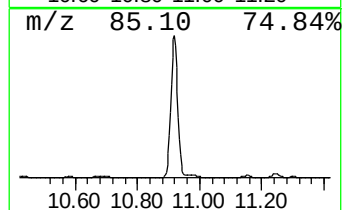
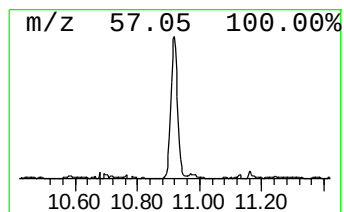
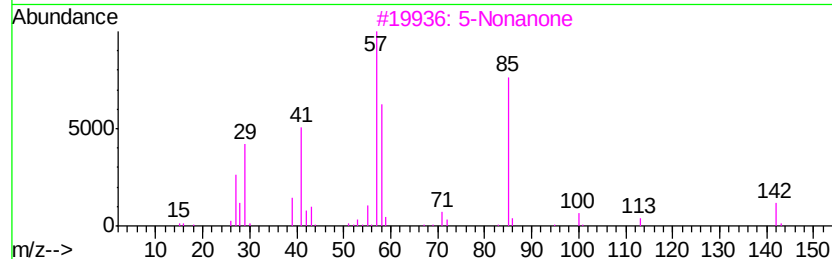
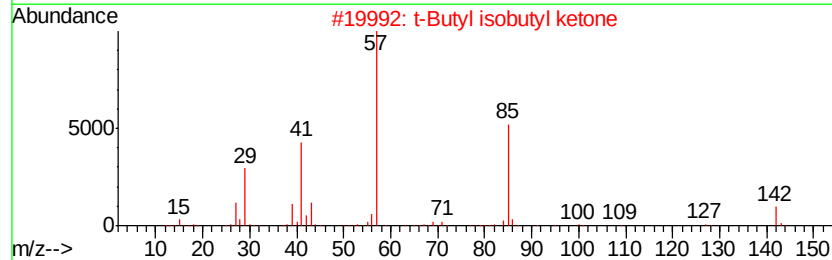
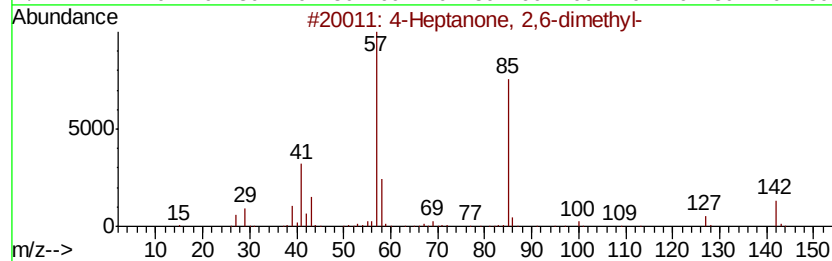
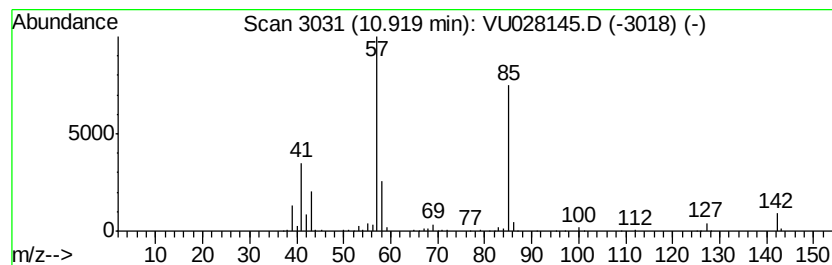
Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

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Peak Number 7 4-Heptanone, 2,6-dimethyl- Concentration Rank 32

| R.T.      | EstConc                        | Area   | Relative to ISTD       | R.T.        |      |
|-----------|--------------------------------|--------|------------------------|-------------|------|
| 10.92     | 2.74 ug/L                      | 264908 | 1,4-Dichlorobenzene-d4 | 11.48       |      |
| Hit# of 5 | Tentative ID                   | MW     | MolForm                | CAS#        | Qual |
| 1         | 4-Heptanone, 2,6-dimethyl-     | 142    | C9H18O                 | 000108-83-8 | 91   |
| 2         | t-Butyl isobutyl ketone        | 142    | C9H18O                 | 014705-50-1 | 64   |
| 3         | 5-Nonanone                     | 142    | C9H18O                 | 000502-56-7 | 59   |
| 4         | 2,4-Hexanedione, 5,5-dimethyl- | 142    | C8H14O2                | 007307-04-2 | 59   |
| 5         | 4-Octanone, 2-methyl-          | 142    | C9H18O                 | 007492-38-8 | 59   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

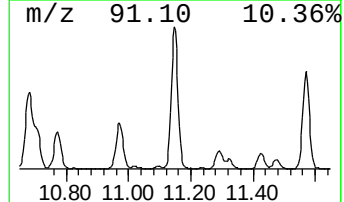
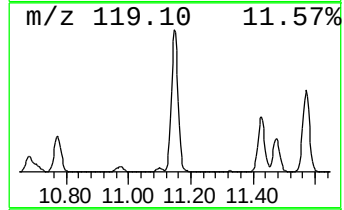
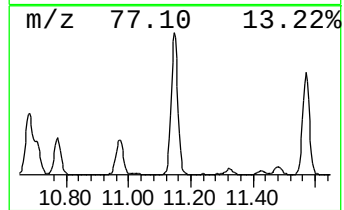
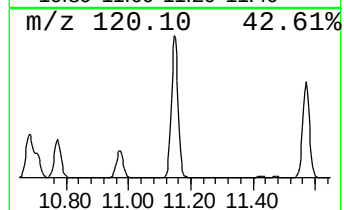
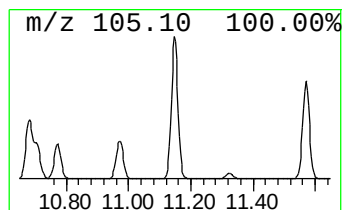
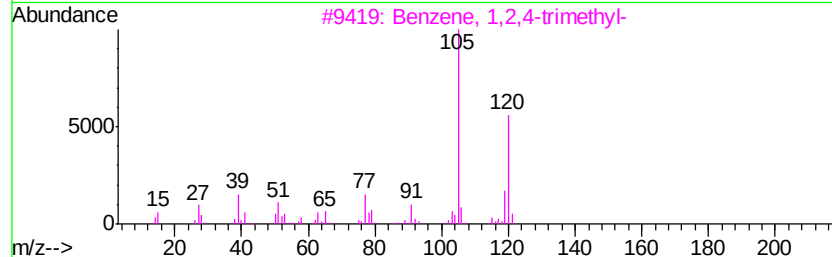
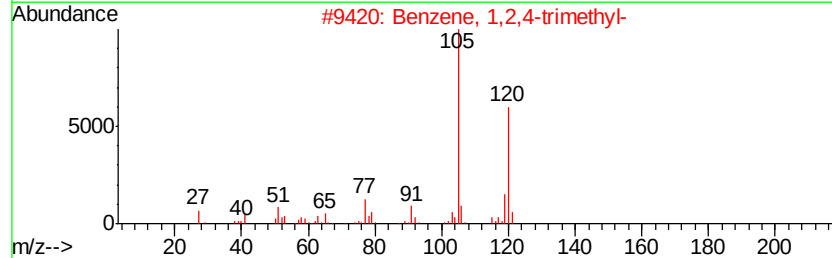
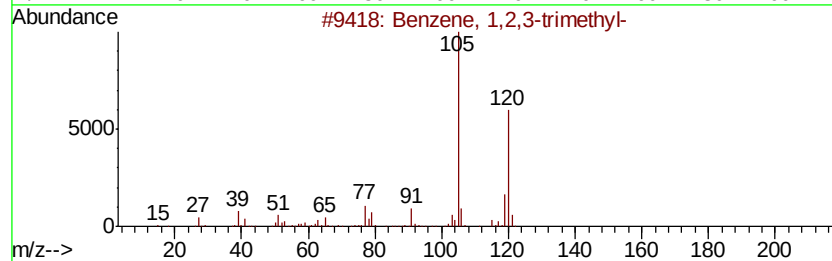
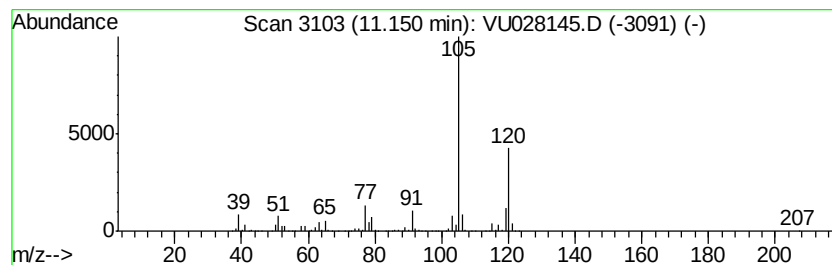
Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

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Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 4

| R.T.    | EstConc    | Area                       | Relative to ISTD       | R.T.           |
|---------|------------|----------------------------|------------------------|----------------|
| 11.15   | 16.60 ug/L | 1604730                    | 1,4-Dichlorobenzene-d4 | 11.48          |
| 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 95 |
| 4       |            | Benzene, 1,2,4-trimethyl-  | 120 C9H12              | 000095-63-6 94 |
| 5       |            | Benzene, 1-ethyl-4-methyl- | 120 C9H12              | 000622-96-8 93 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

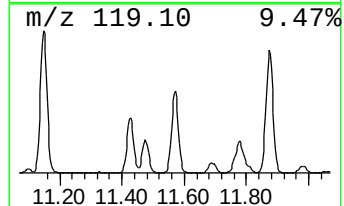
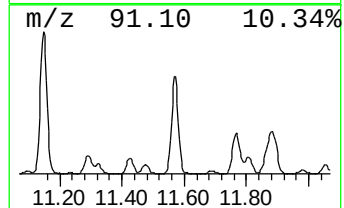
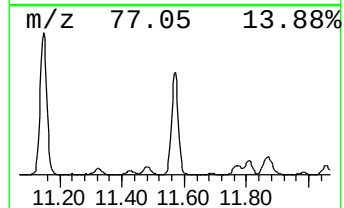
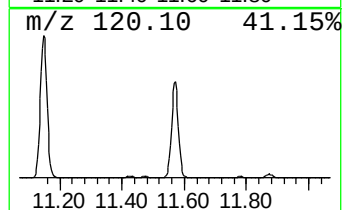
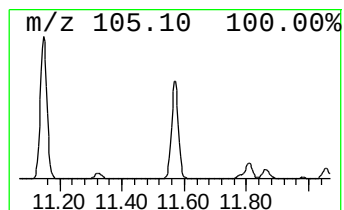
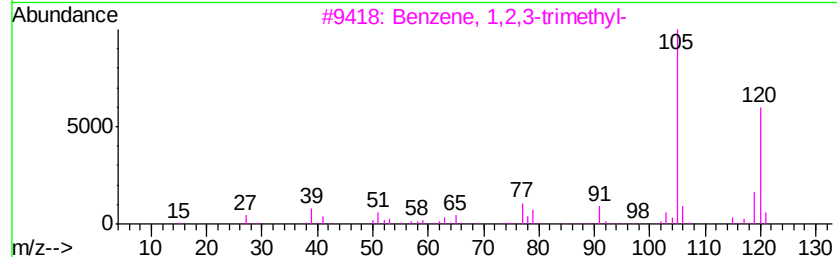
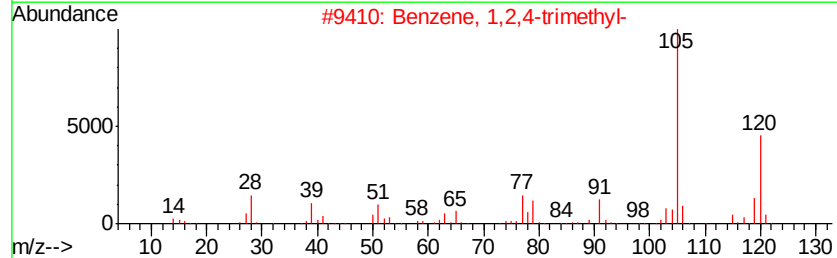
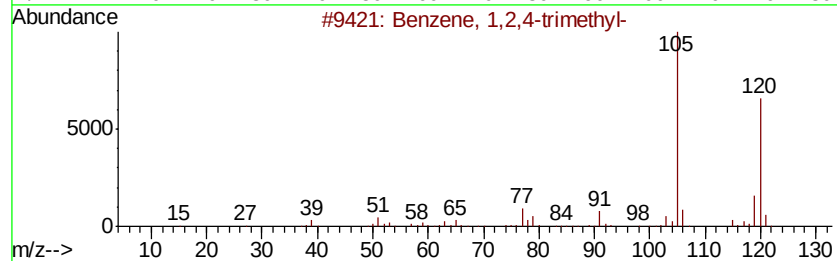
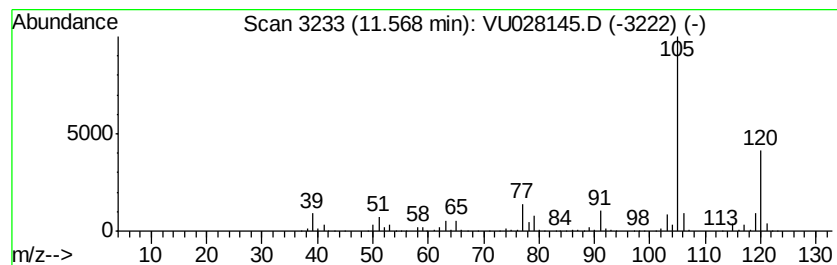
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Benzene, 1,2,4-trimethyl- Concentration Rank 7

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 11.57 | 11.66 ug/L | 1126640 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 2    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 3    |    | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 94   |
| 4    |    | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 94   |
| 5    |    | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

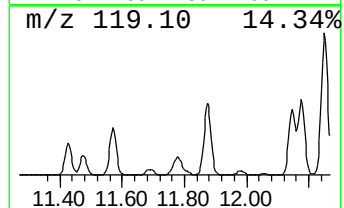
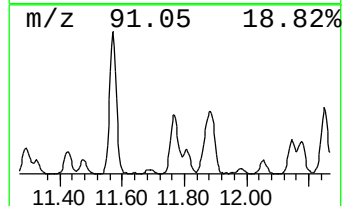
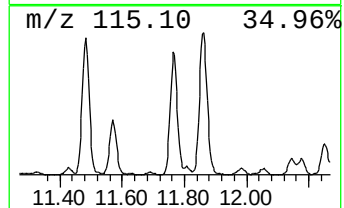
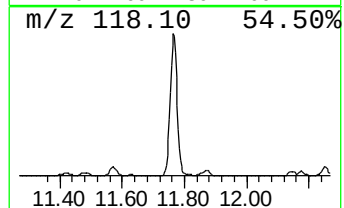
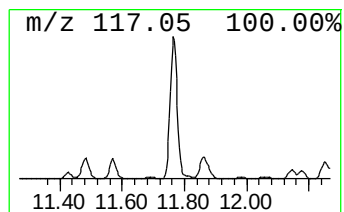
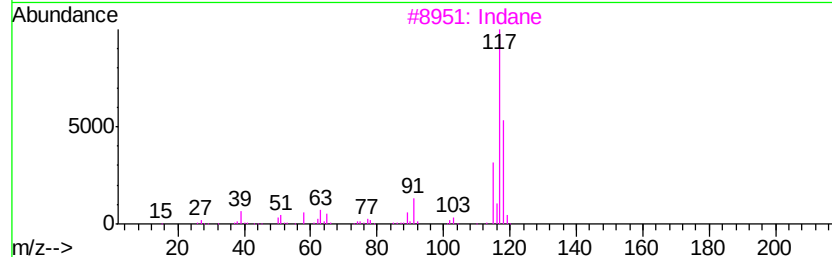
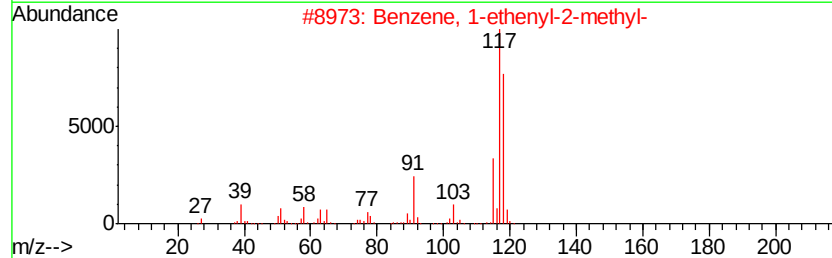
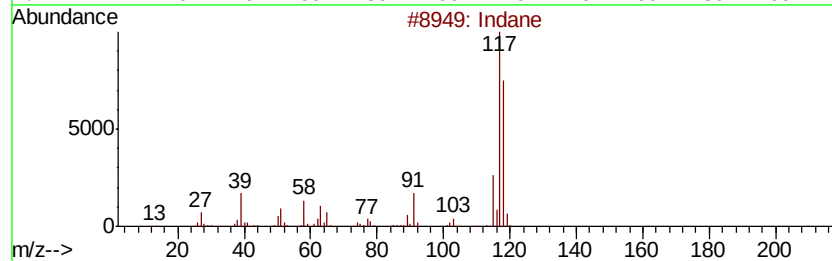
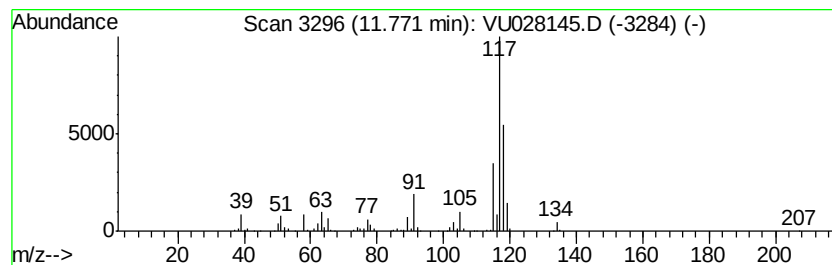
Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 11 Indane Concentration Rank 26

| R.T.    | EstConc                      | Area         | Relative to ISTD       | R.T.           |
|---------|------------------------------|--------------|------------------------|----------------|
| 11.77   | 3.72 ug/L                    | 359447       | 1,4-Dichlorobenzene-d4 | 11.48          |
| Hit# of | 5                            | Tentative ID | MW MolForm             | CAS# Qual      |
| 1       | Indane                       |              | 118 C9H10              | 000496-11-7 95 |
| 2       | Benzene, 1-ethenyl-2-methyl- |              | 118 C9H10              | 000611-15-4 89 |
| 3       | Indane                       |              | 118 C9H10              | 000496-11-7 87 |
| 4       | Indane                       |              | 118 C9H10              | 000496-11-7 81 |
| 5       | Benzene, cyclopropyl-        |              | 118 C9H10              | 000873-49-4 81 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

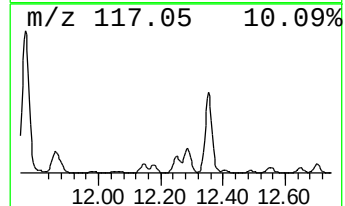
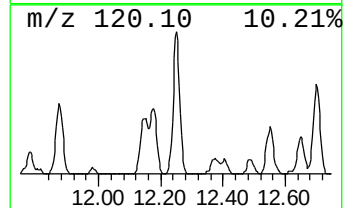
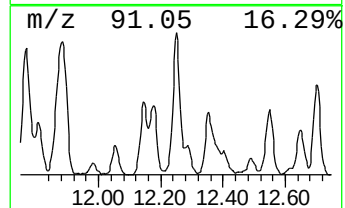
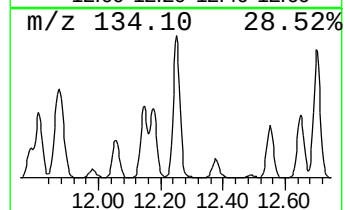
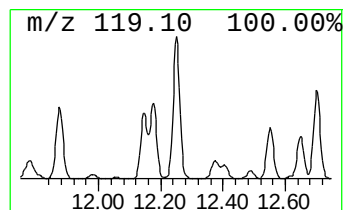
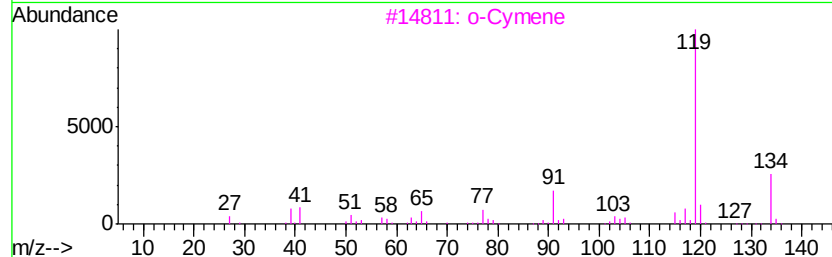
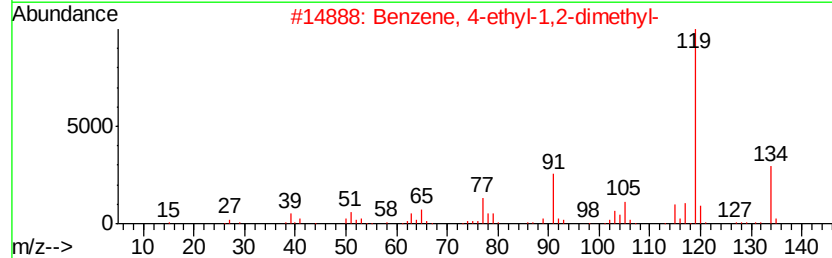
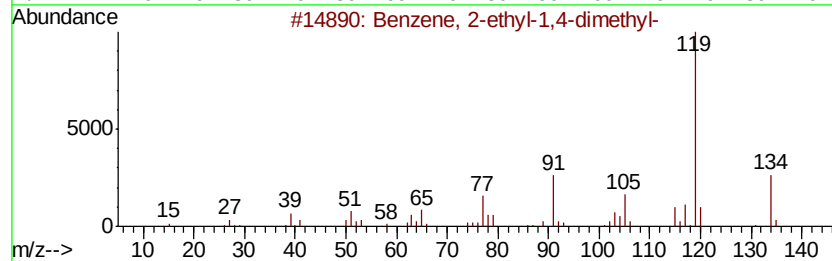
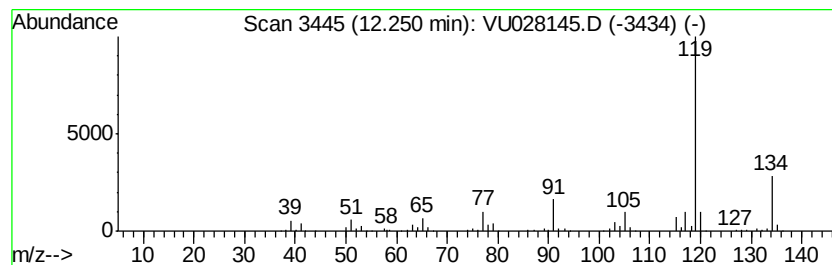
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 28

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.25 | 3.24 ug/L | 312730 | 1,4-Dichlorobenzene-d4 | 11.48 |

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

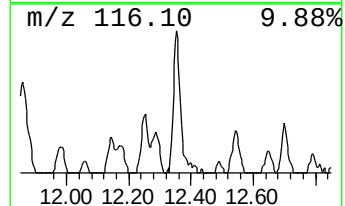
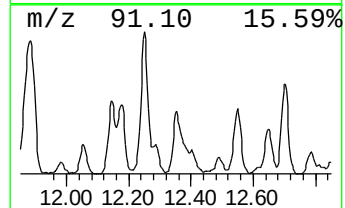
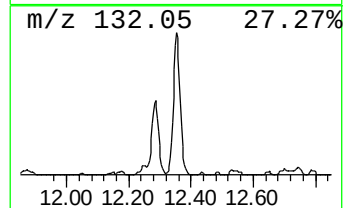
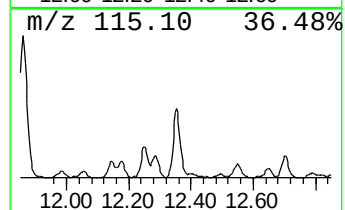
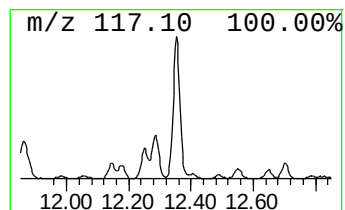
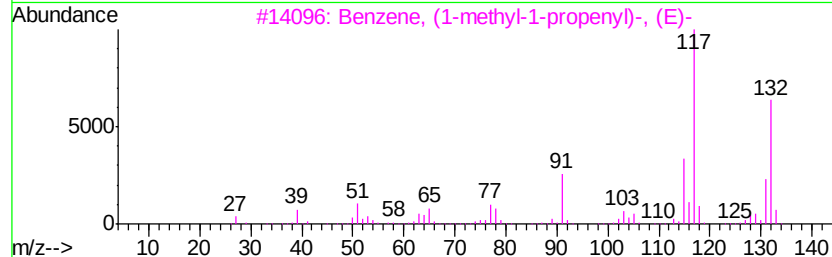
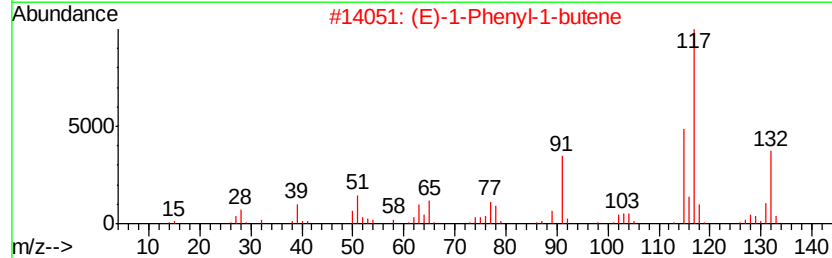
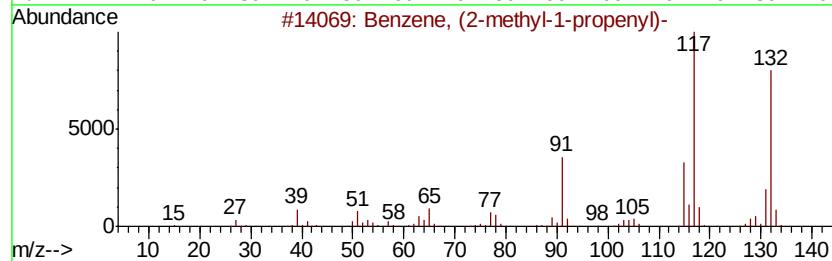
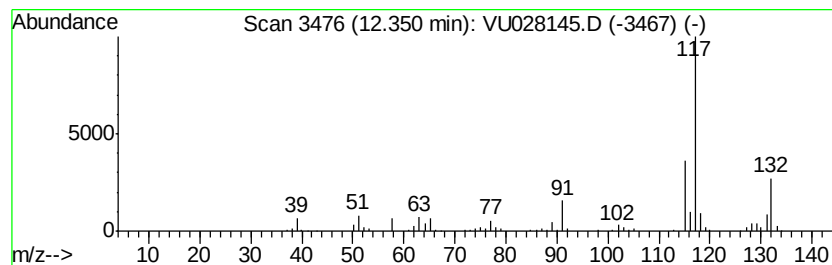
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Benzene, (2-methyl-1-propen... Concentration Rank 31

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.35 | 2.80 ug/L | 271118 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (2-methyl-1-propenyl)-     | 132 | C10H12  | 000768-49-0 | 90   |
| 2    |    |   | (E)-1-Phenyl-1-butene               | 132 | C10H12  | 001005-64-7 | 87   |
| 3    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 87   |
| 4    |    |   | Indan, 1-methyl-                    | 132 | C10H12  | 000767-58-8 | 87   |
| 5    |    |   | Benzene, 1-ethenyl-4-ethyl-         | 132 | C10H12  | 003454-07-7 | 86   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

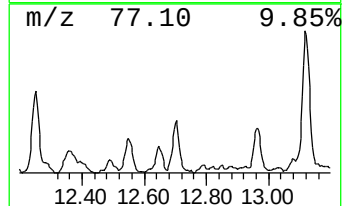
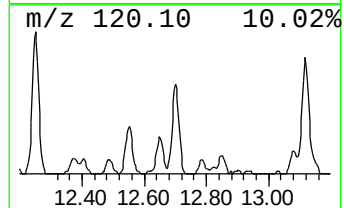
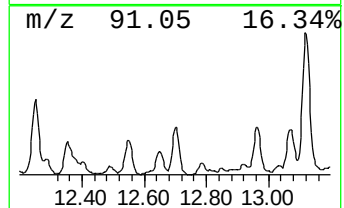
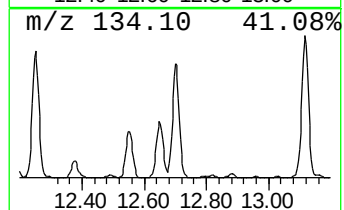
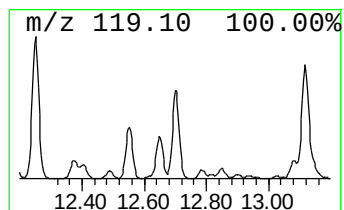
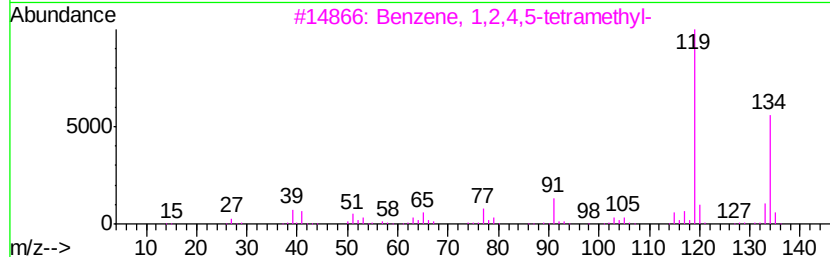
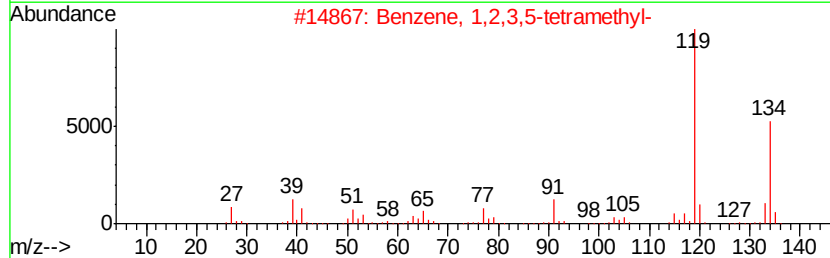
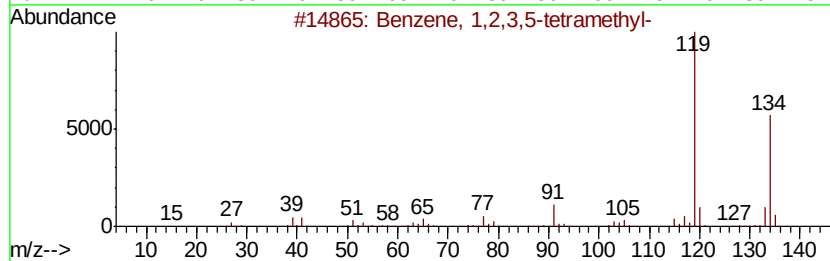
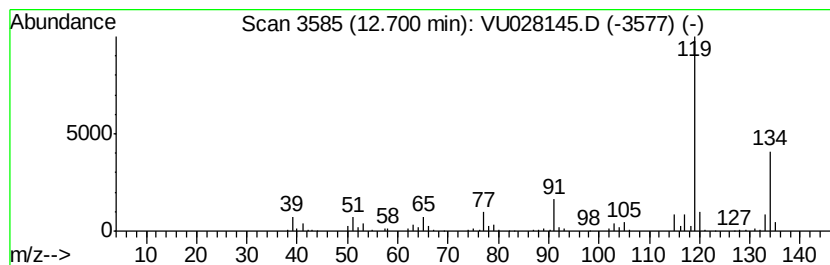
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 36

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.70 | 2.34 ug/L | 226039 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 96   |
| 2    |    | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 95   |
| 3    |    | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 95   |
| 4    |    | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 95   |
| 5    |    | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 95   |





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

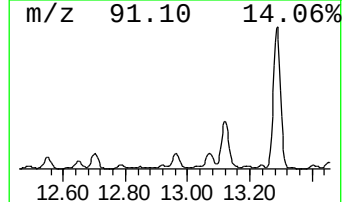
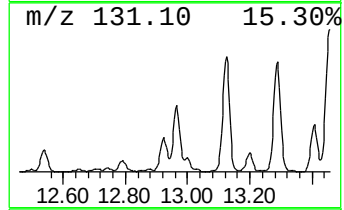
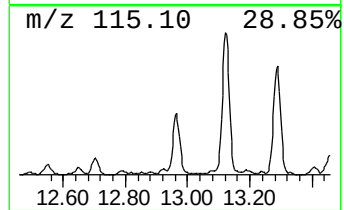
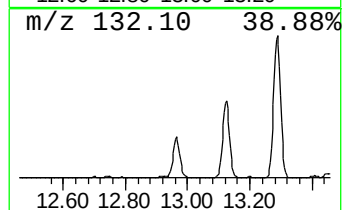
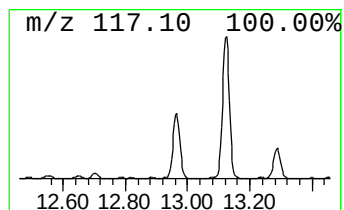
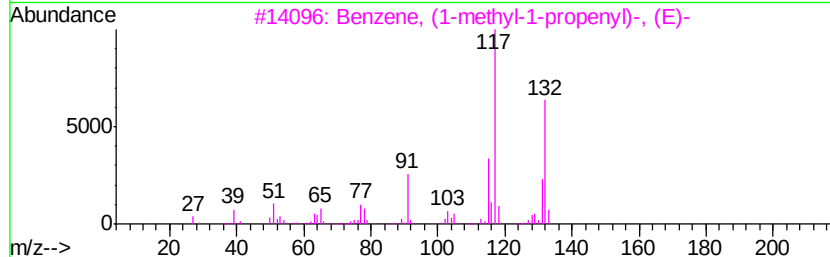
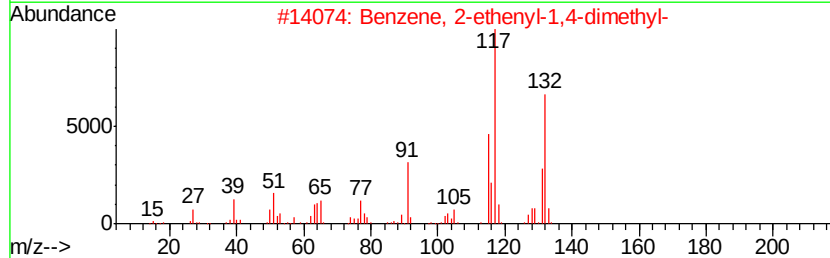
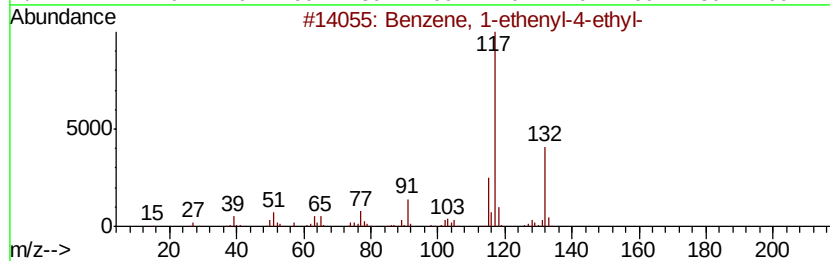
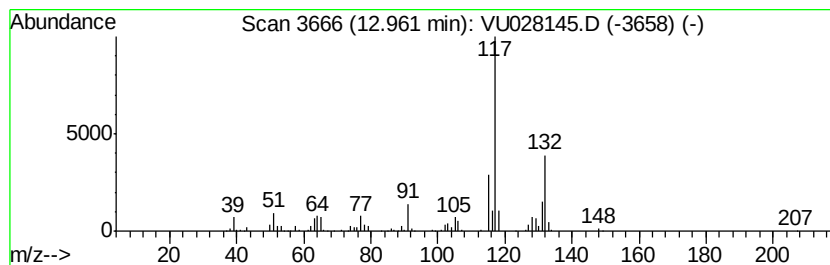
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 30

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.96 | 2.92 ug/L | 281928 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethenyl-4-ethyl-         | 132 | C10H12  | 003454-07-7 | 94   |
| 2    |    | Benzene, 2-ethenyl-1,4-dimethyl-    | 132 | C10H12  | 002039-89-6 | 92   |
| 3    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 91   |
| 4    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000767-99-7 | 91   |
| 5    |    | Benzene, (2-methyl-2-propenyl)-     | 132 | C10H12  | 003290-53-7 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

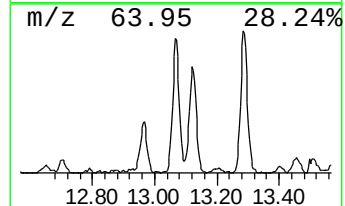
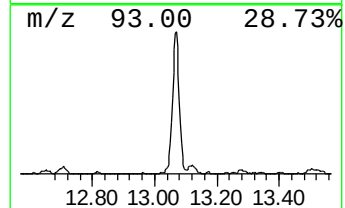
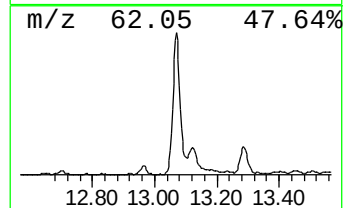
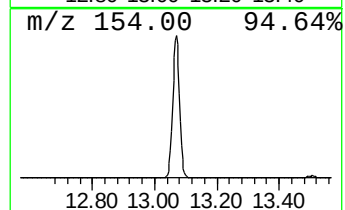
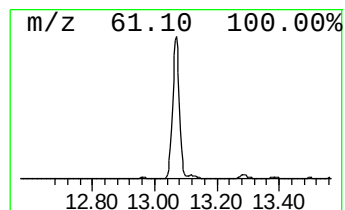
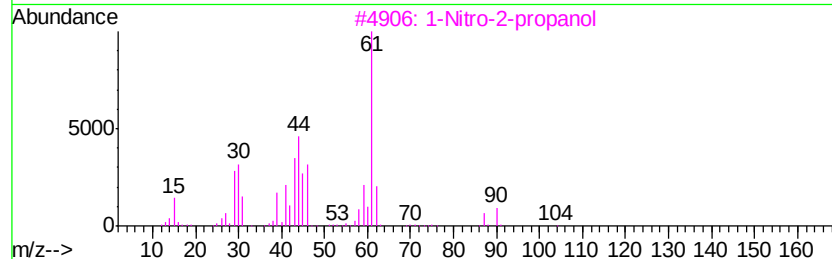
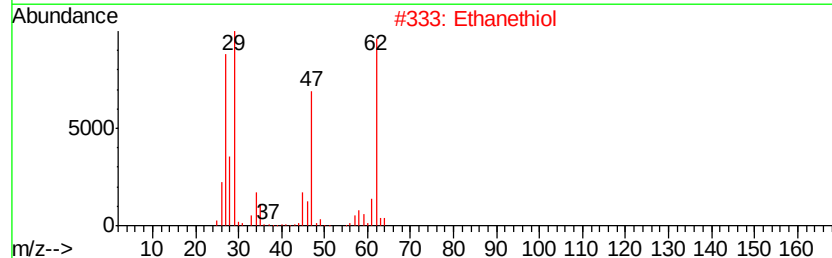
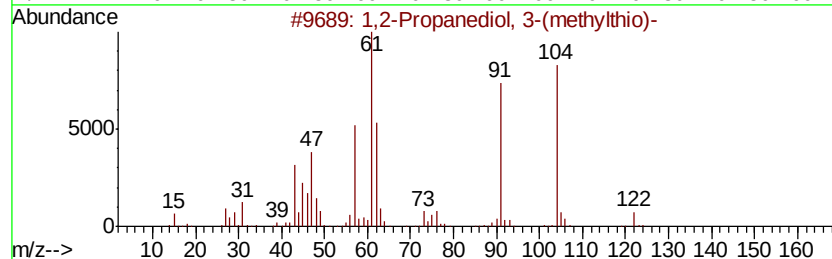
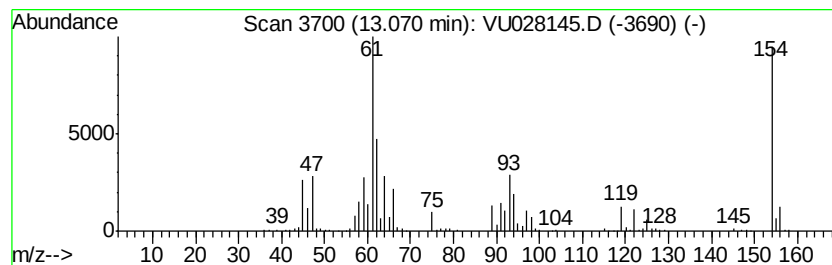
Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 16 unknown-01 Concentration Rank 23

| R.T.    | EstConc                          | Area         | Relative to ISTD       | R.T.      |
|---------|----------------------------------|--------------|------------------------|-----------|
| 13.07   | 4.27 ug/L                        | 412696       | 1,4-Dichlorobenzene-d4 | 11.48     |
| Hit# of | 5                                | Tentative ID | MW MolForm             | CAS# Qual |
| 1       | 1,2-Propanediol, 3-(methylthio)- | 122 C4H10O2S | 022551-26-4            | 37        |
| 2       | Ethanethiol                      | 62 C2H6S     | 000075-08-1            | 16        |
| 3       | 1-Nitro-2-propanol               | 105 C3H7NO3  | 003156-73-8            | 10        |
| 4       | Trisulfide, diethyl              | 154 C4H10S3  | 003600-24-6            | 10        |
| 5       | 2,4-Dithiapentane                | 108 C3H8S2   | 001618-26-4            | 10        |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

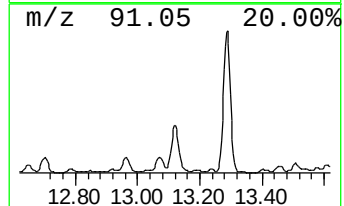
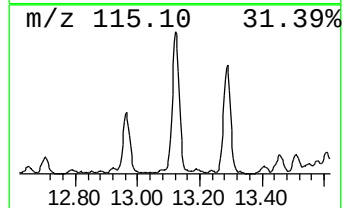
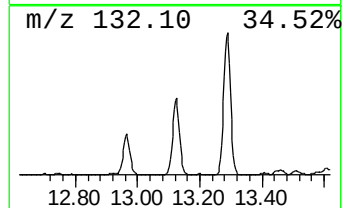
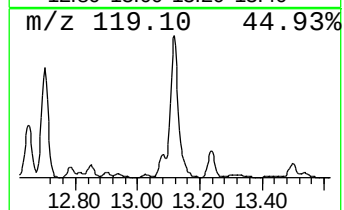
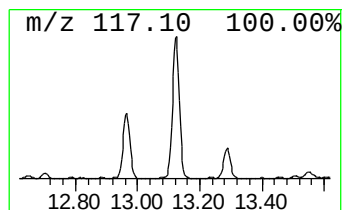
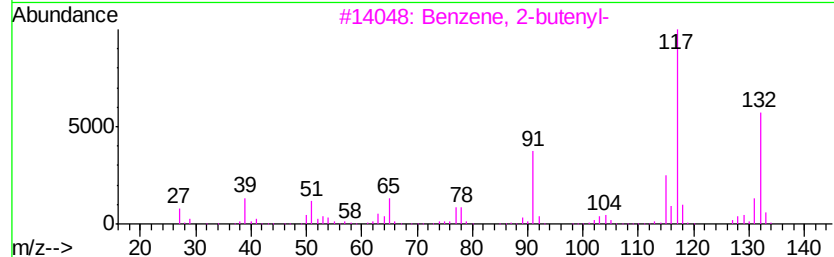
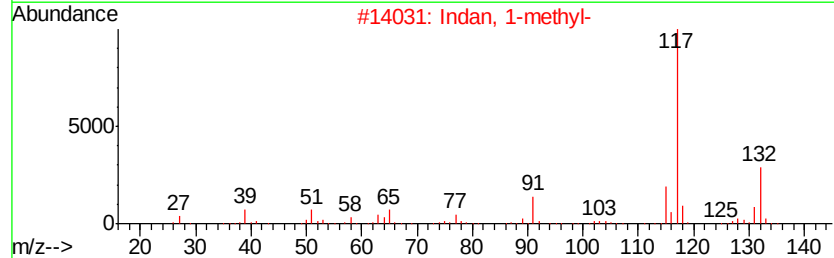
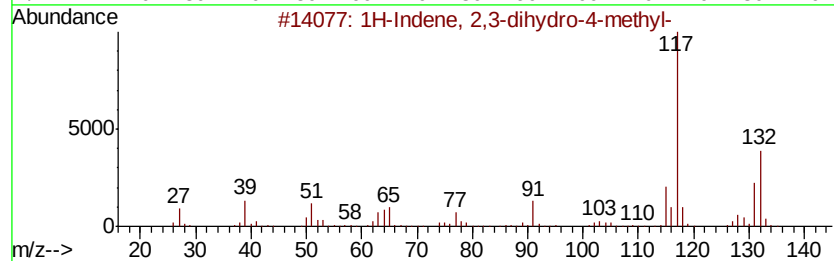
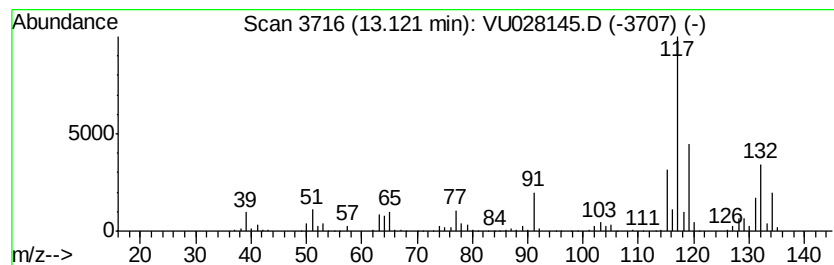
Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 17 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 11

| R.T.    | EstConc   | Area                             | Relative to ISTD       | R.T.    |             |      |
|---------|-----------|----------------------------------|------------------------|---------|-------------|------|
| 13.12   | 8.28 ug/L | 800180                           | 1,4-Dichlorobenzene-d4 | 11.48   |             |      |
| Hit# of | 5         | Tentative ID                     | MW                     | MolForm | CAS#        | Qual |
| 1       |           | 1H-Indene, 2,3-dihydro-4-methyl- | 132                    | C10H12  | 000824-22-6 | 91   |
| 2       |           | Indan, 1-methyl-                 | 132                    | C10H12  | 000767-58-8 | 70   |
| 3       |           | Benzene, 2-butenyl-              | 132                    | C10H12  | 001560-06-1 | 70   |
| 4       |           | 3-Phenylbut-1-ene                | 132                    | C10H12  | 000934-10-1 | 70   |
| 5       |           | Benzene, 1-ethenyl-4-ethyl-      | 132                    | C10H12  | 003454-07-7 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

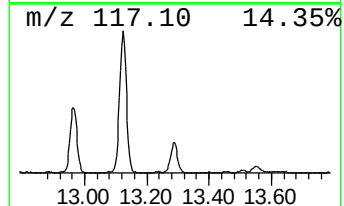
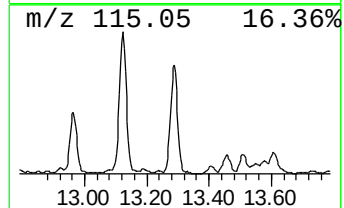
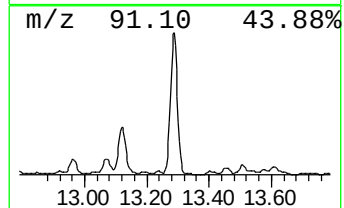
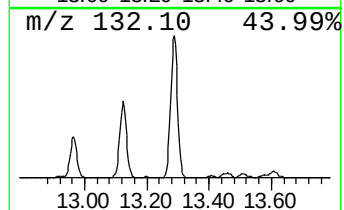
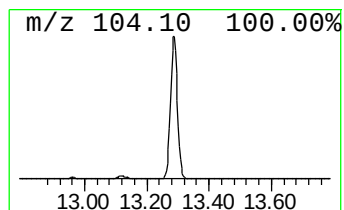
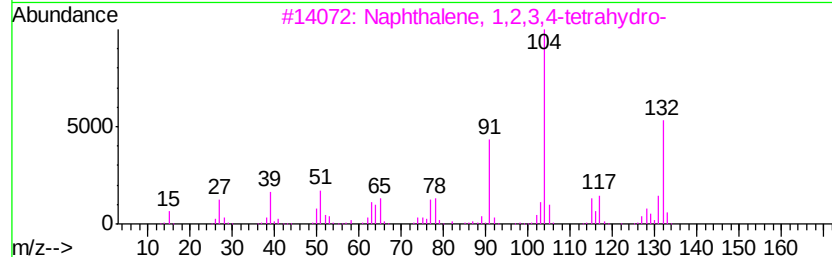
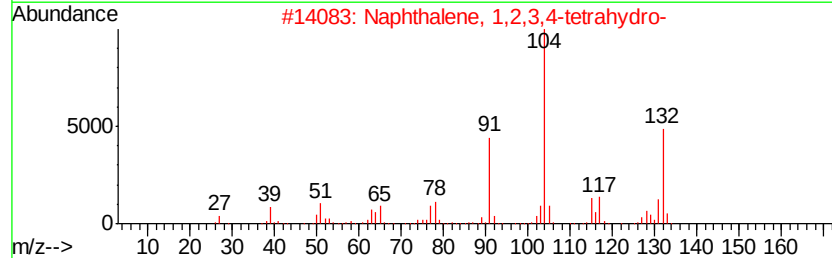
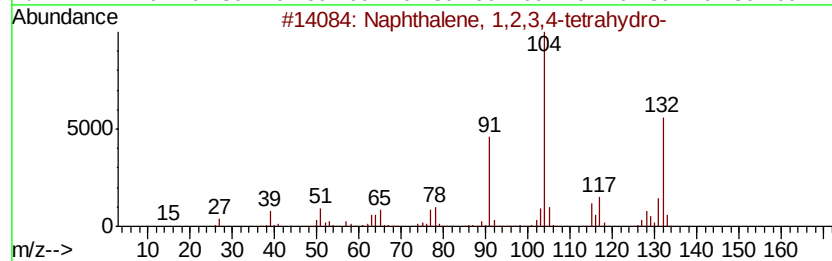
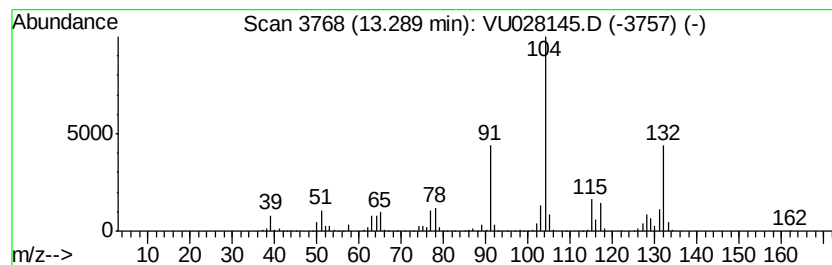
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 13.29 | 10.66 ug/L | 1030380 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 96   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 96   |
| 3    |    | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 96   |
| 4    |    | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 95   |
| 5    |    | 2H-Inden-2-one, 1,3-dihydro-     | 132 | C9H8O   | 000615-13-4 | 62   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

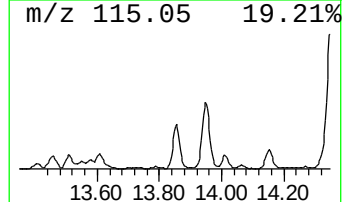
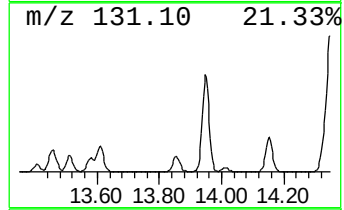
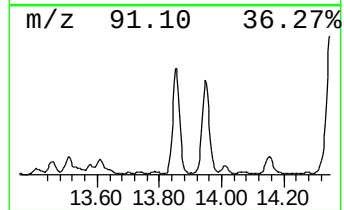
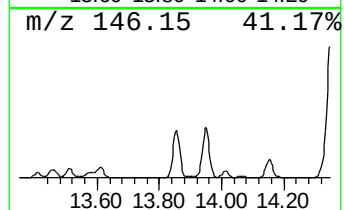
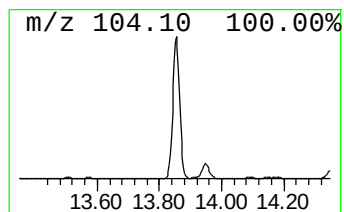
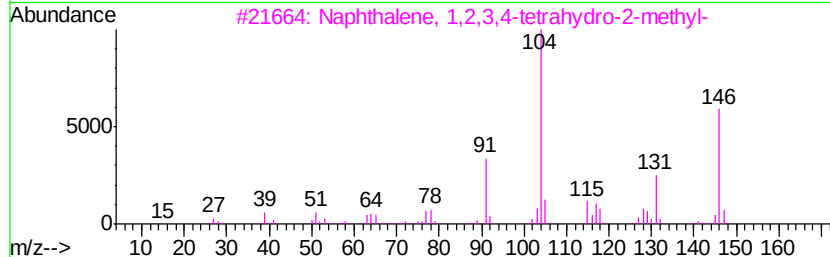
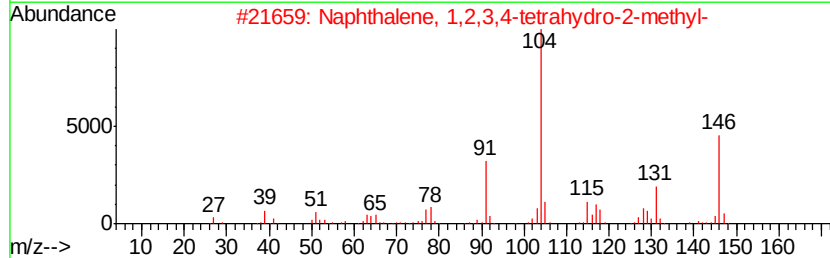
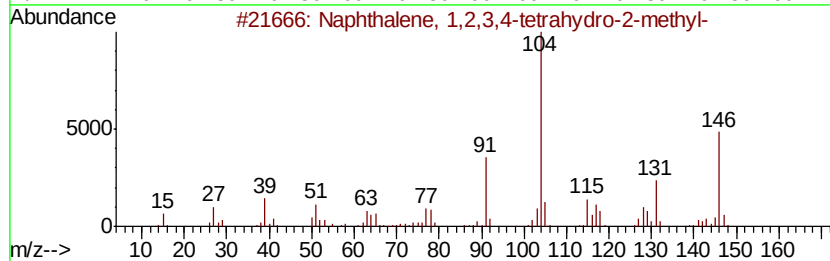
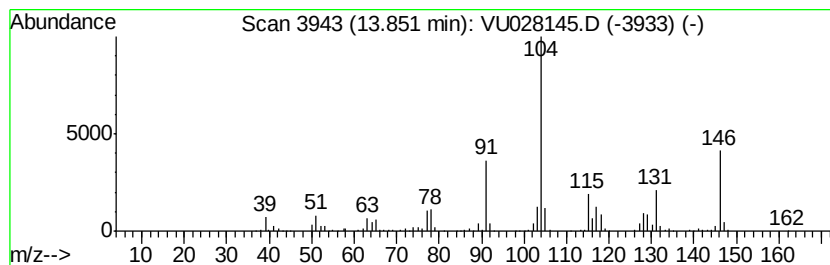
Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 20 Naphthalene, 1,2,3,4-tetra... Concentration Rank 17

| R.T.    | EstConc   | Area                                | Relative to ISTD       | R.T.           |
|---------|-----------|-------------------------------------|------------------------|----------------|
| 13.85   | 5.43 ug/L | 524688                              | 1,4-Dichlorobenzene-d4 | 11.48          |
| Hit# of | 5         | Tentative ID                        | MW MolForm             | CAS# Qual      |
| 1       |           | Naphthalene, 1,2,3,4-tetrahydro-... | 146 C11H14             | 003877-19-8 93 |
| 2       |           | Naphthalene, 1,2,3,4-tetrahydro-... | 146 C11H14             | 003877-19-8 91 |
| 3       |           | Naphthalene, 1,2,3,4-tetrahydro-... | 146 C11H14             | 003877-19-8 87 |
| 4       |           | Naphthalene, 1,2,3,4-tetrahydro-... | 146 C11H14             | 003877-19-8 81 |
| 5       |           | Naphth[1,2-b]oxirene, 1a,2,3,7b-... | 146 C10H10O            | 002461-34-9 70 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

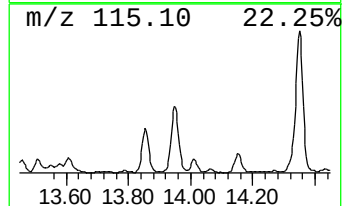
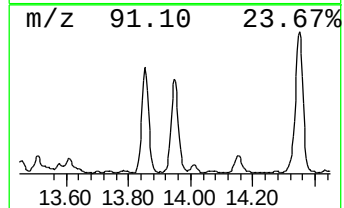
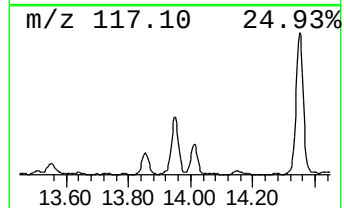
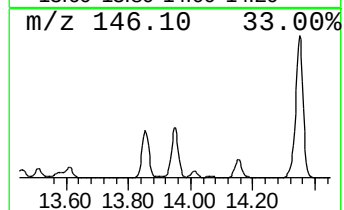
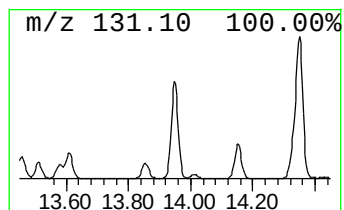
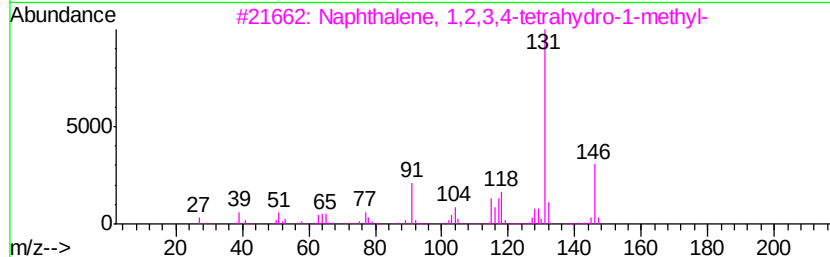
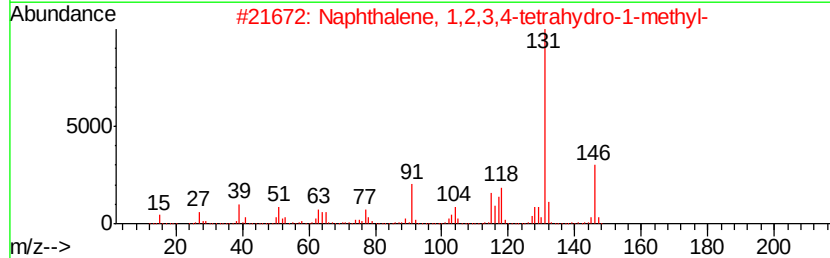
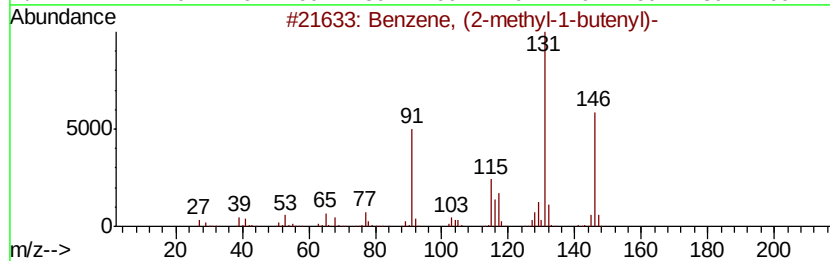
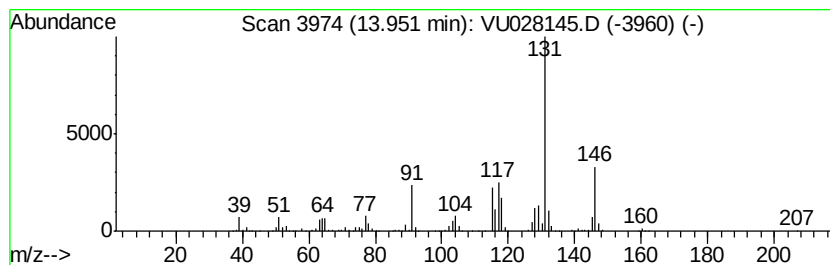
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 Benzene, (2-methyl-1-butenyl)- Concentration Rank 14

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 13.95 | 7.08 ug/L | 684636 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (2-methyl-1-butenyl)-      | 146 | C11H14  | 056253-64-6 | 95   |
| 2    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001559-81-5 | 94   |
| 3    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001559-81-5 | 93   |
| 4    |    |   | 2,2-Dimethylindene, 2,3-dihydro-    | 146 | C11H14  | 020836-11-7 | 91   |
| 5    |    |   | 1H-Inden-1-one, 2,3-dihydro-2-me... | 146 | C10H10O | 017496-14-9 | 87   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

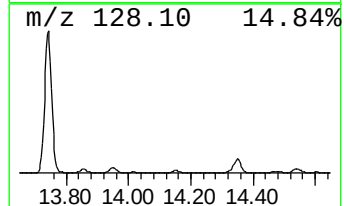
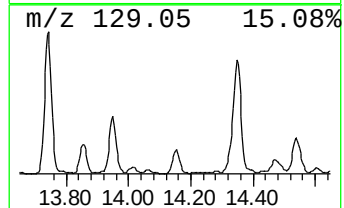
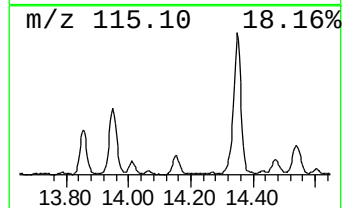
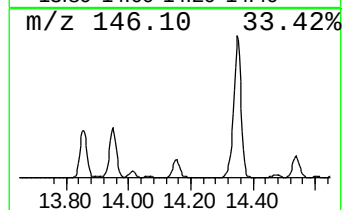
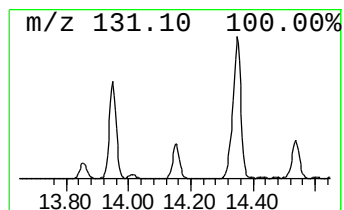
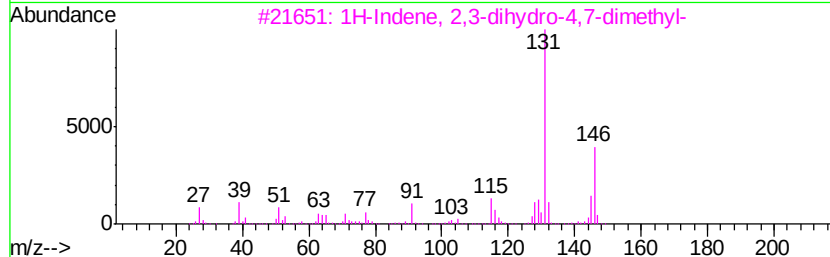
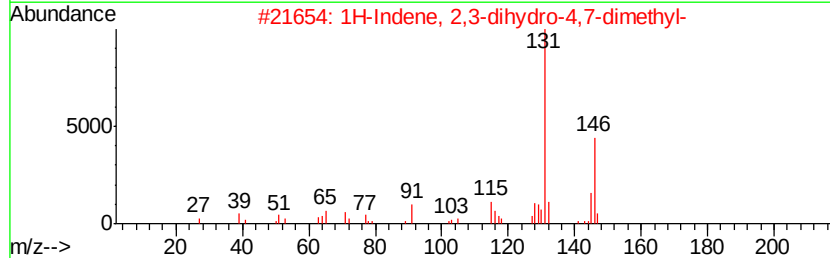
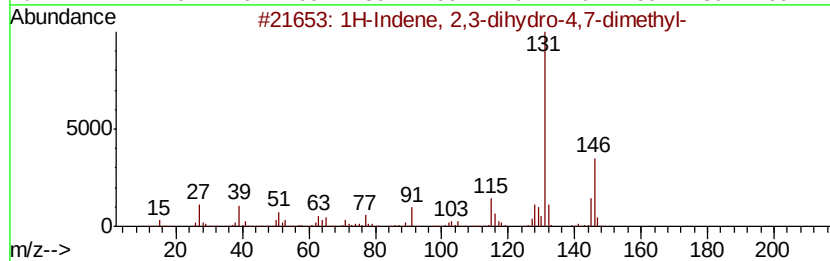
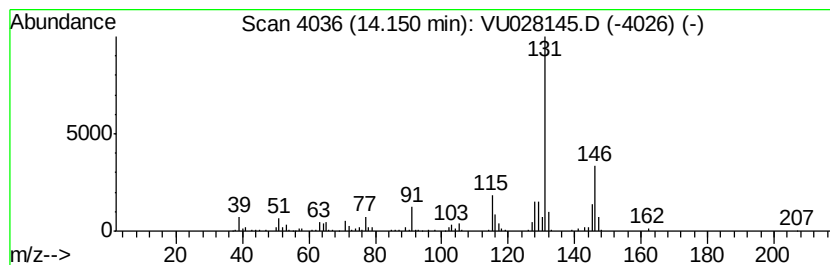
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 34

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 14.15 | 2.62 ug/L | 253260 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 97   |
| 2    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 96   |
| 3    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 96   |
| 4    |    | 1H-Indene, 2,3-dihydro-5,6-dimet... | 146 | C11H14  | 001075-22-5 | 94   |
| 5    |    | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

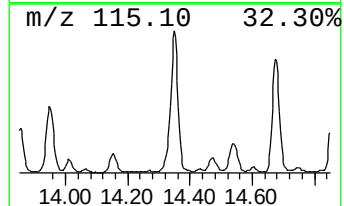
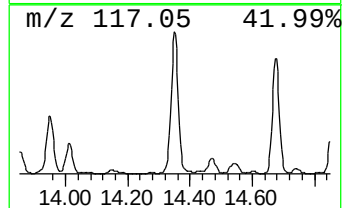
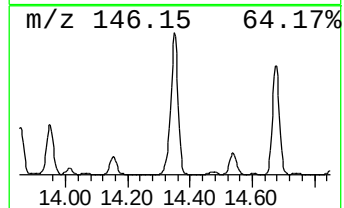
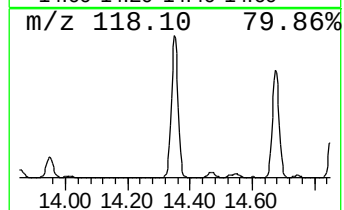
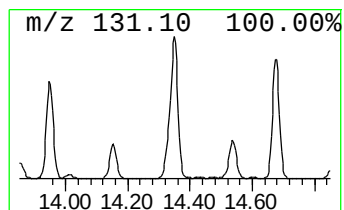
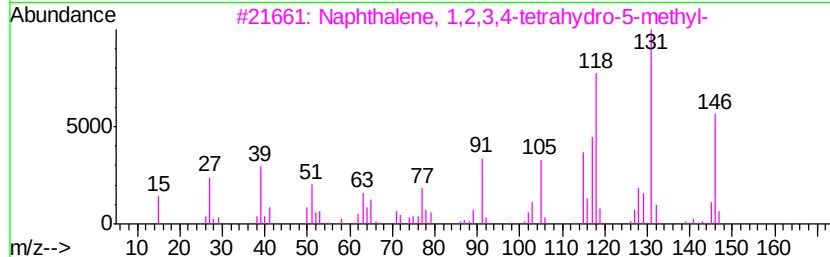
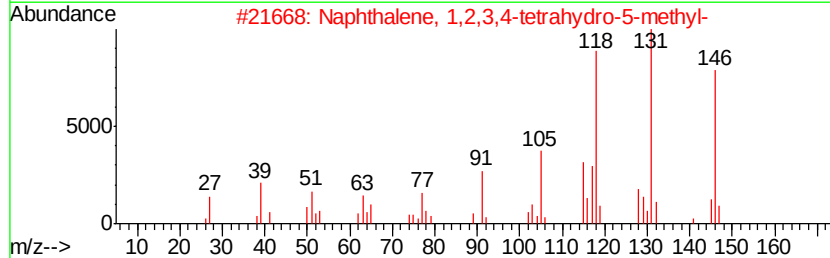
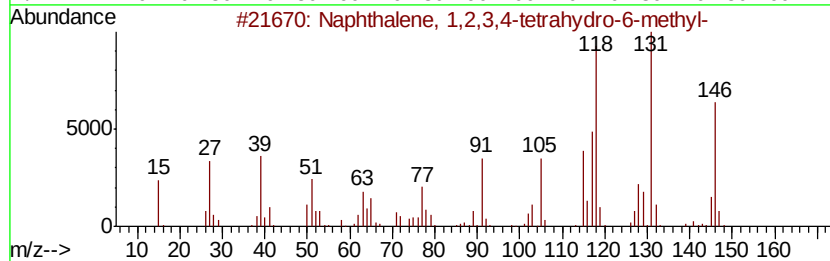
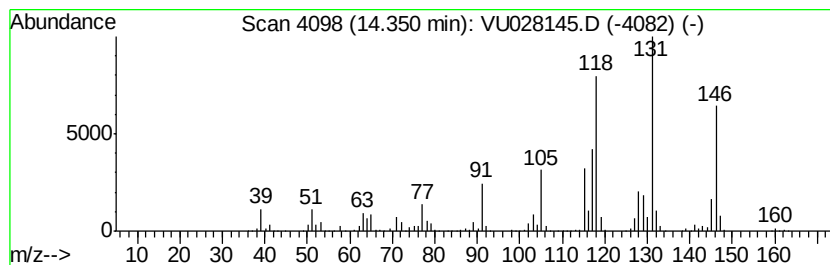
Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 23 Naphthalene, 1,2,3,4-tetra... Concentration Rank 2

| R.T.    | EstConc    | Area                                | Relative to ISTD       | R.T.    |             |      |
|---------|------------|-------------------------------------|------------------------|---------|-------------|------|
| 14.35   | 17.47 ug/L | 1688660                             | 1,4-Dichlorobenzene-d4 | 11.48   |             |      |
| Hit# of | 5          | Tentative ID                        | MW                     | MolForm | CAS#        | Qual |
| 1       |            | Naphthalene, 1,2,3,4-tetrahydro-... | 146                    | C11H14  | 001680-51-9 | 96   |
| 2       |            | Naphthalene, 1,2,3,4-tetrahydro-... | 146                    | C11H14  | 002809-64-5 | 96   |
| 3       |            | Naphthalene, 1,2,3,4-tetrahydro-... | 146                    | C11H14  | 002809-64-5 | 95   |
| 4       |            | Naphthalene, 1,2,3,4-tetrahydro-... | 146                    | C11H14  | 002809-64-5 | 95   |
| 5       |            | Naphthalene, 1,2,3,4-tetrahydro-... | 146                    | C11H14  | 001680-51-9 | 94   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

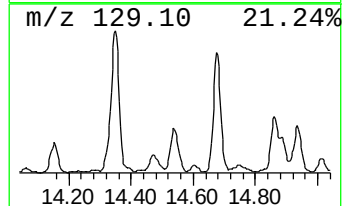
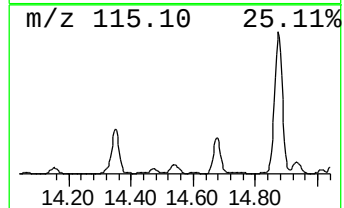
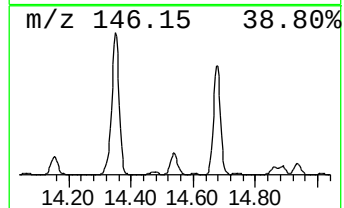
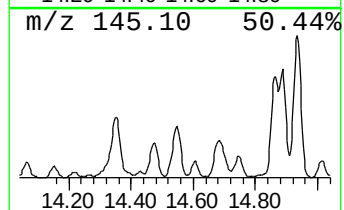
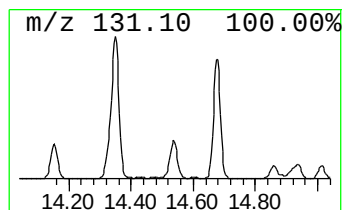
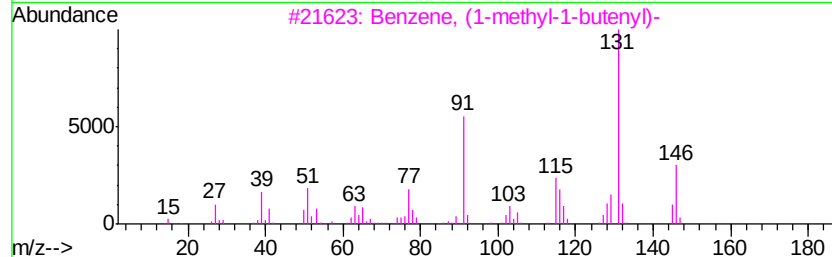
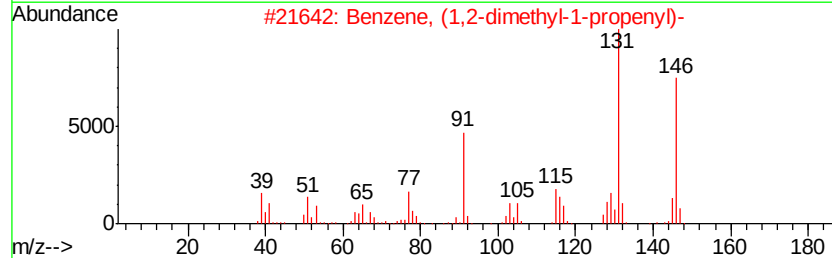
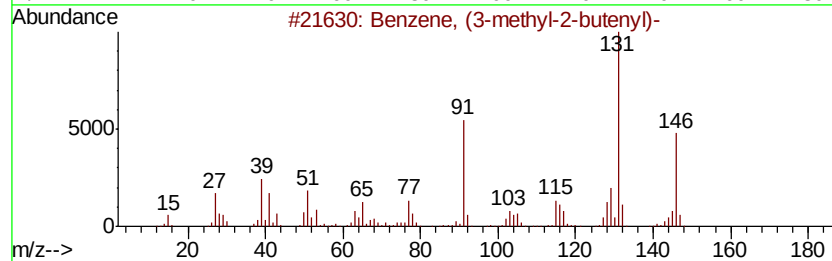
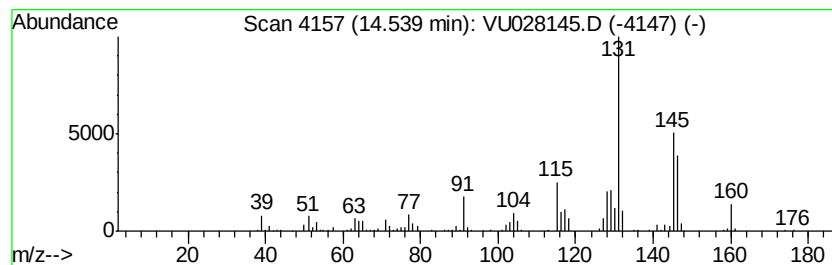
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 Benzene, (3-methyl-2-butenyl)- Concentration Rank 25

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 14.54 | 3.72 ug/L | 359465 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 76   |
| 2    |    | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 70   |
| 3    |    | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 70   |
| 4    |    | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 58   |
| 5    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 58   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

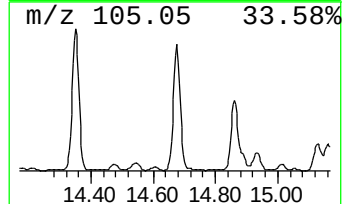
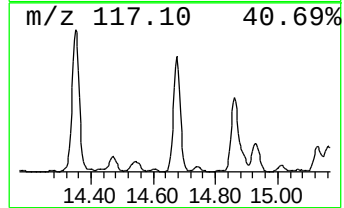
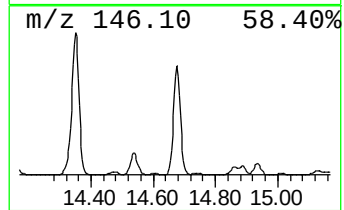
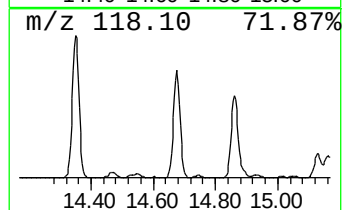
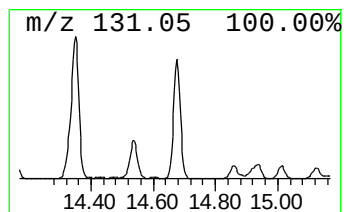
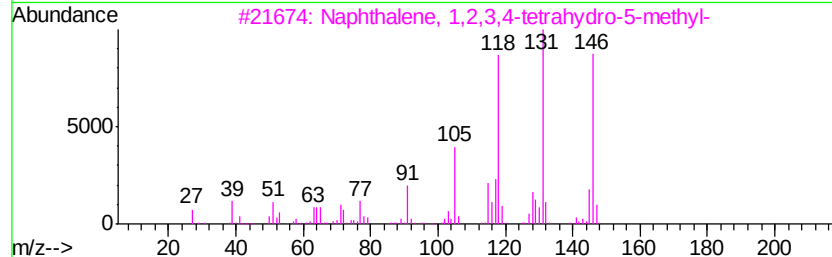
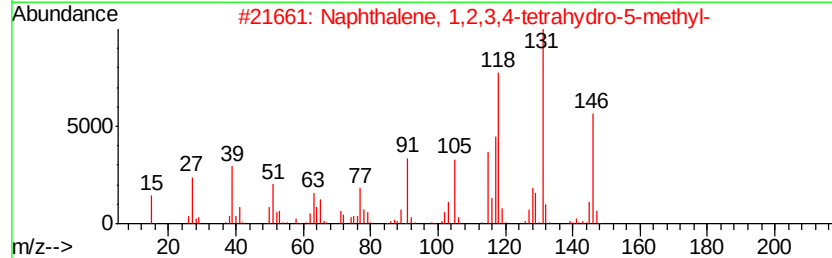
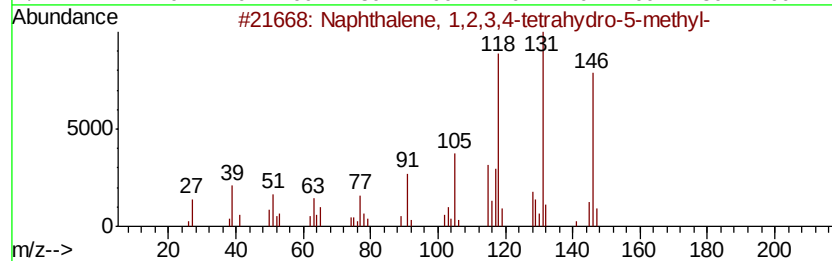
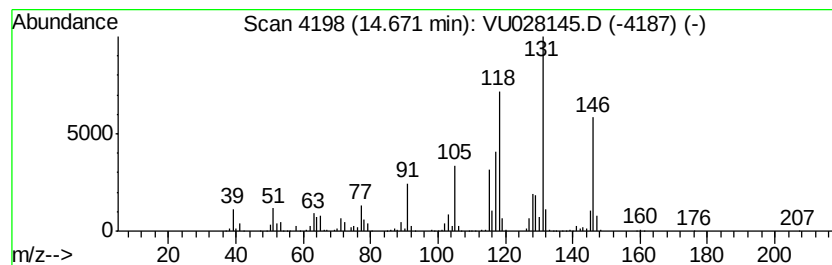
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MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 25 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 5

| R.T.    | EstConc                             | Area         | Relative to ISTD       | R.T.      |
|---------|-------------------------------------|--------------|------------------------|-----------|
| 14.67   | 13.10 ug/L                          | 1266190      | 1,4-Dichlorobenzene-d4 | 11.48     |
| Hit# of | 5                                   | Tentative ID | MW MolForm             | CAS# Qual |
| 1       | Naphthalene, 1,2,3,4-tetrahydro-... | 146 C11H14   | 002809-64-5            | 97        |
| 2       | Naphthalene, 1,2,3,4-tetrahydro-... | 146 C11H14   | 002809-64-5            | 96        |
| 3       | Naphthalene, 1,2,3,4-tetrahydro-... | 146 C11H14   | 002809-64-5            | 95        |
| 4       | Naphthalene, 1,2,3,4-tetrahydro-... | 146 C11H14   | 001680-51-9            | 95        |
| 5       | Naphthalene, 1,2,3,4-tetrahydro-... | 146 C11H14   | 002809-64-5            | 94        |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

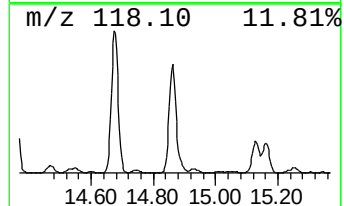
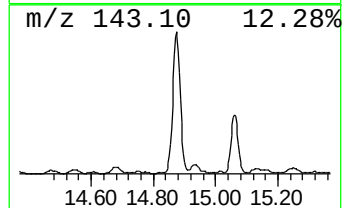
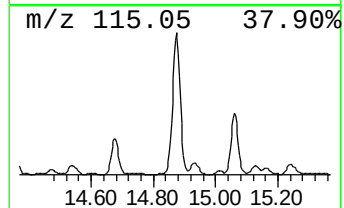
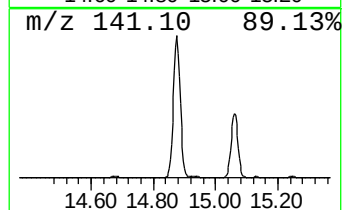
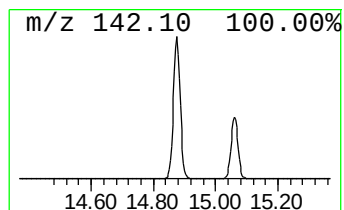
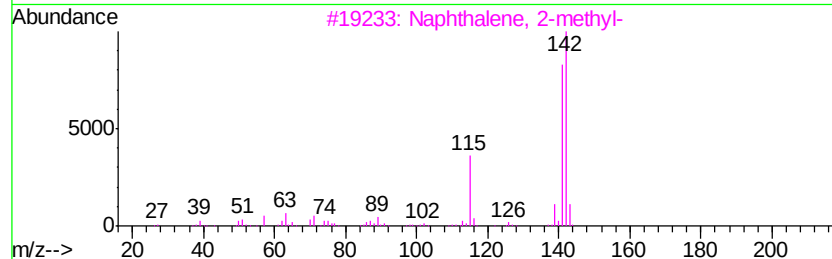
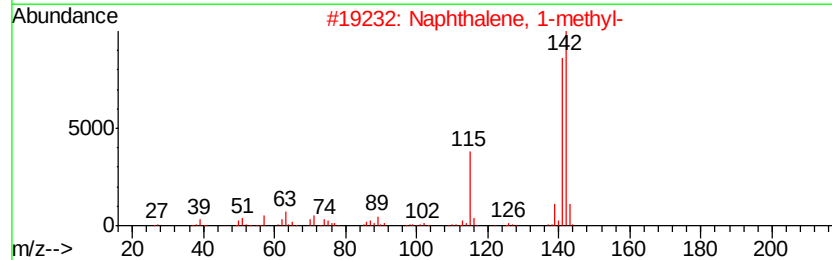
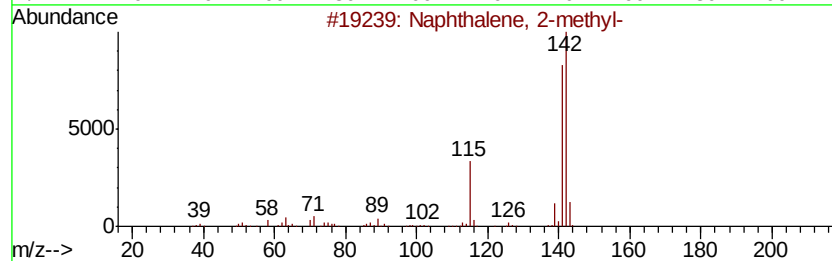
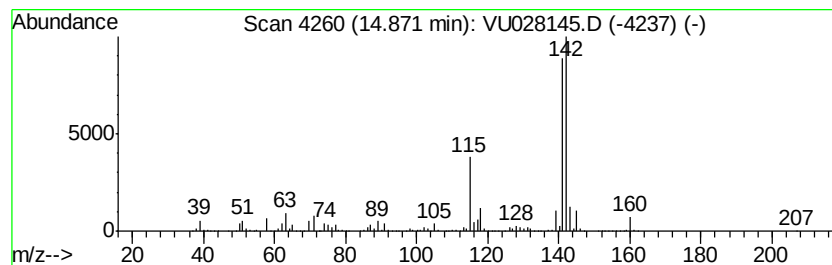
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 Naphthalene, 1-methyl- Concentration Rank 1

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 14.87 | 33.02 ug/L | 3191800 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 96   |
| 2    |    | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 96   |
| 3    |    | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 96   |
| 4    |    | Benzocycloheptatriene  | 142 | C11H10  | 000264-09-5 | 93   |
| 5    |    | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

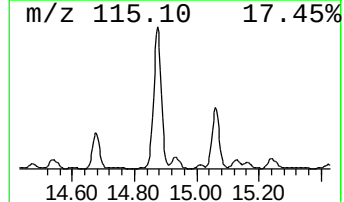
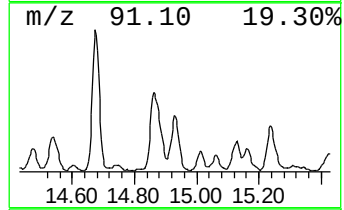
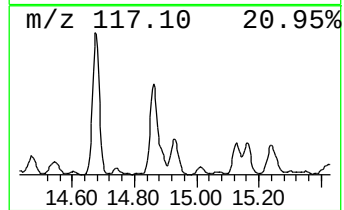
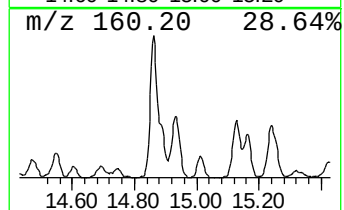
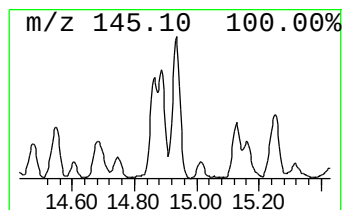
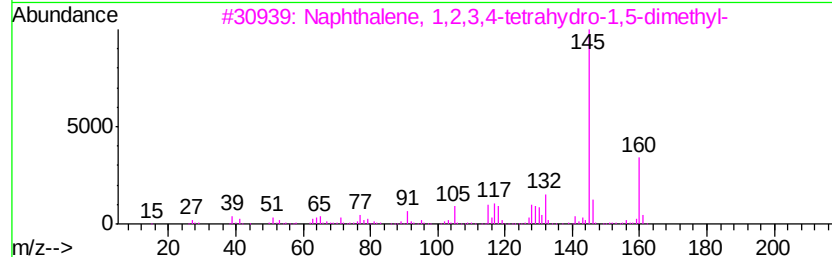
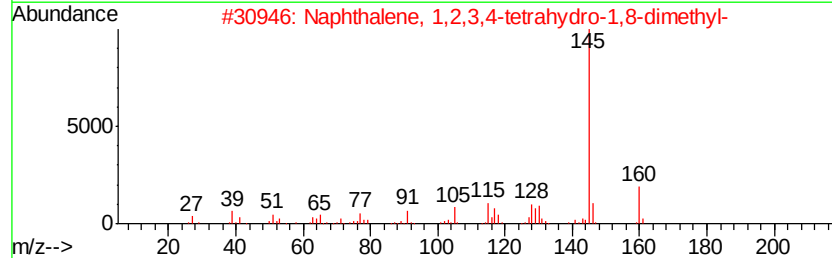
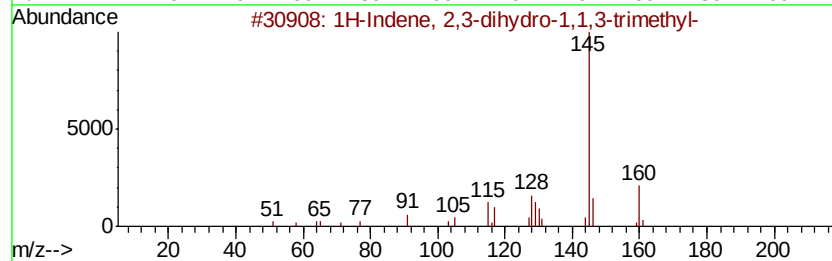
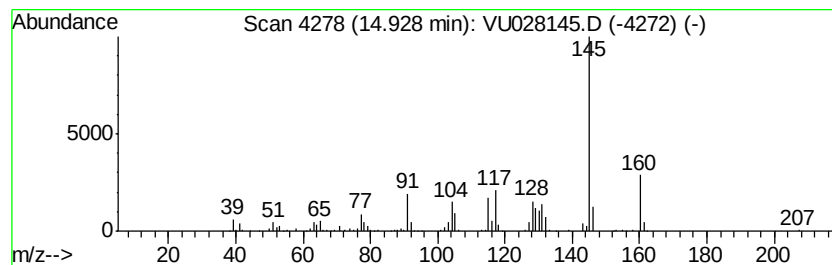
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 19

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 14.93 | 4.92 ug/L | 475931 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 2,3-dihydro-1,1,3-tri... | 160 | C12H16  | 002613-76-5 | 90   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 025419-33-4 | 90   |
| 3    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 021564-91-0 | 90   |
| 4    |    | 1H-Indene, 2,3-dihydro-1,1,5-tri... | 160 | C12H16  | 040650-41-7 | 87   |
| 5    |    | Benzene, 1-(1-methylethenyl)-2-(... | 160 | C12H16  | 005557-93-7 | 87   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

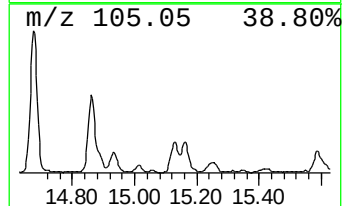
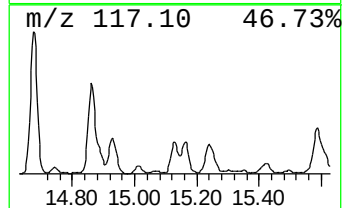
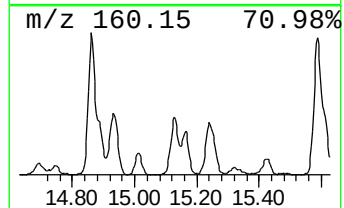
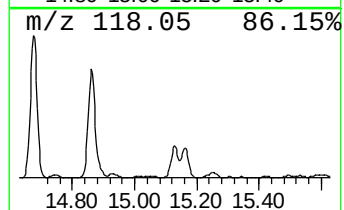
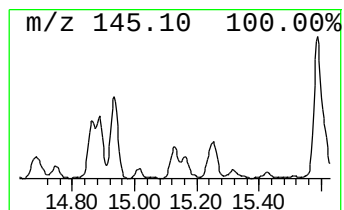
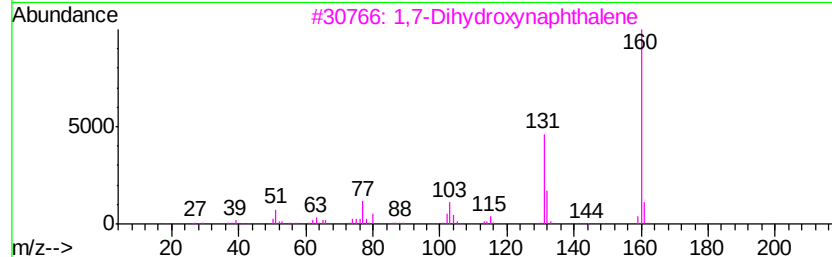
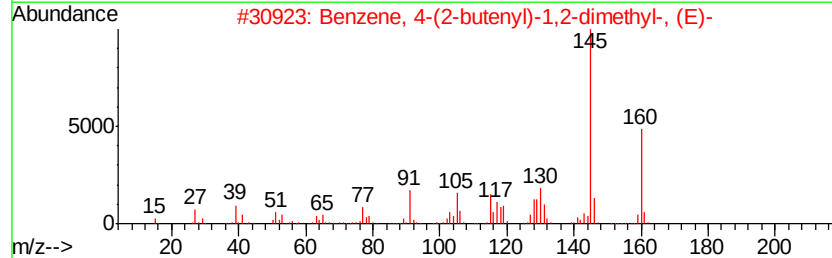
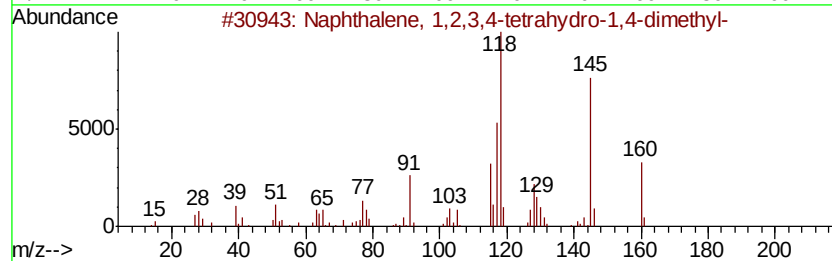
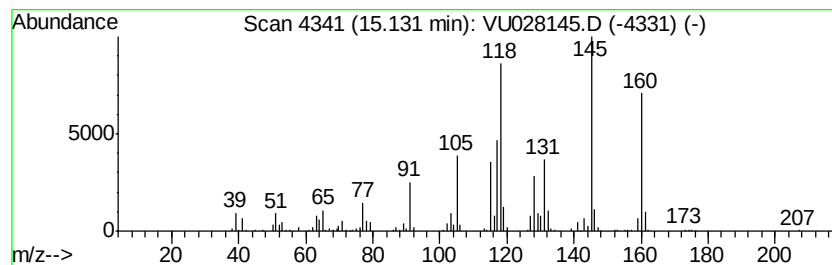
Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 29 Naphthalene, 1,2,3,4-tetra... Concentration Rank 29

| R.T.    | EstConc   | Area                                | Relative to ISTD       | R.T.    |             |      |
|---------|-----------|-------------------------------------|------------------------|---------|-------------|------|
| 15.13   | 3.15 ug/L | 304943                              | 1,4-Dichlorobenzene-d4 | 11.48   |             |      |
| Hit# of | 5         | Tentative ID                        | MW                     | MolForm | CAS#        | Qual |
| 1       |           | Naphthalene, 1,2,3,4-tetrahydro-... | 160                    | C12H16  | 004175-54-6 | 95   |
| 2       |           | Benzene, 4-(2-butenyl)-1,2-dimet... | 160                    | C12H16  | 054340-86-2 | 83   |
| 3       |           | 1,7-Dihydroxynaphthalene            | 160                    | C10H8O2 | 000575-38-2 | 64   |
| 4       |           | Benzene, 1,3,5-trimethyl-2-(1-me... | 160                    | C12H16  | 014679-13-1 | 64   |
| 5       |           | Benzene, 1-(1-methylethenyl)-3-(... | 160                    | C12H16  | 001129-29-9 | 60   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

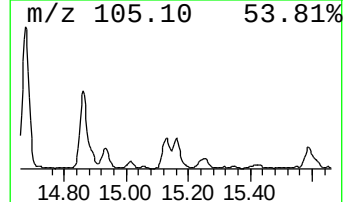
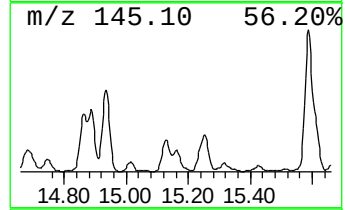
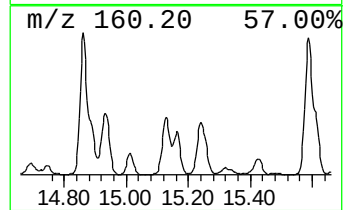
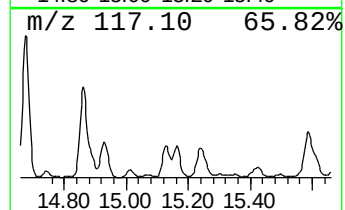
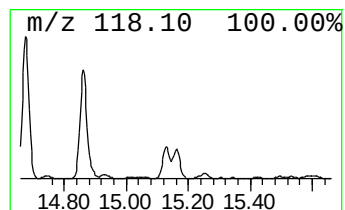
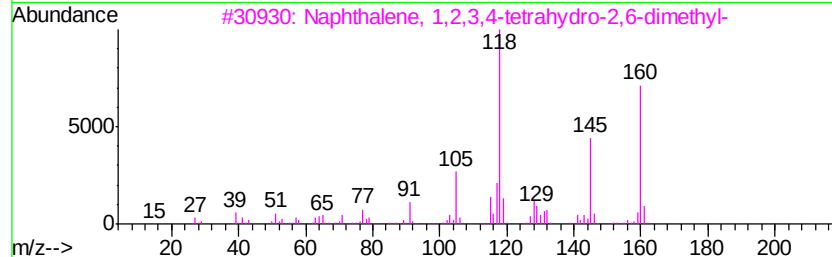
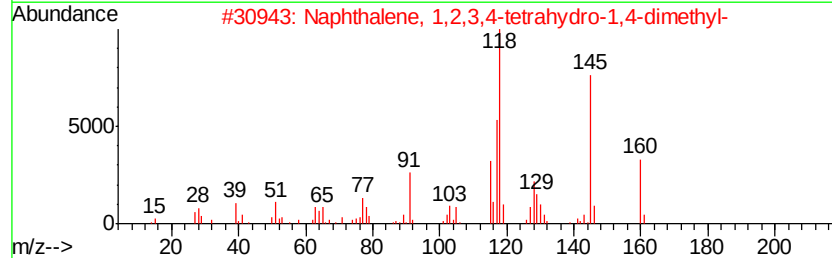
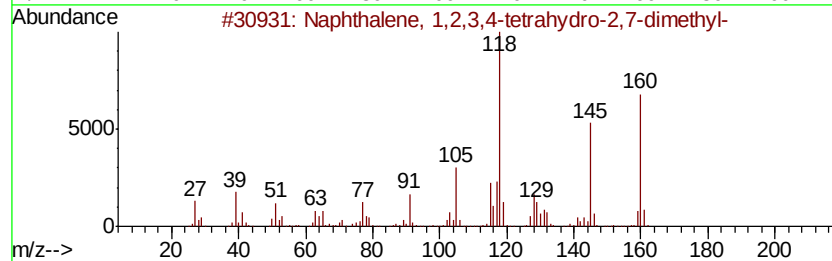
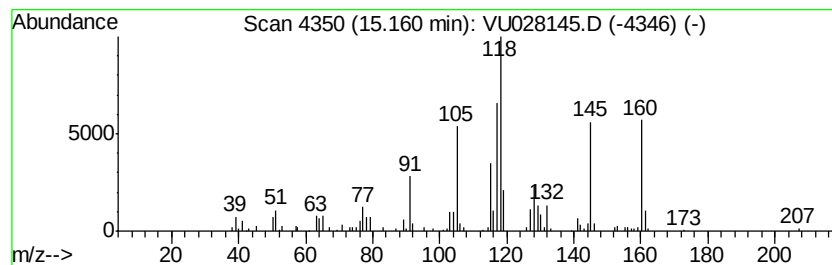
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 Naphthalene, 1,2,3,4-tetra... Concentration Rank 37

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 15.16 | 2.24 ug/L | 216859 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 013065-07-1 | 76   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 004175-54-6 | 74   |
| 3    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 007524-63-2 | 68   |
| 4    |    | Benzene, 1,3,5-trimethyl-2-(1-me... | 160 | C12H16  | 014679-13-1 | 55   |
| 5    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 007524-63-2 | 49   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

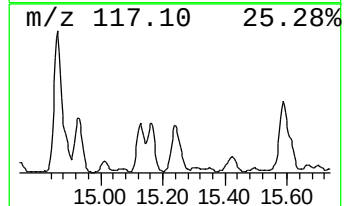
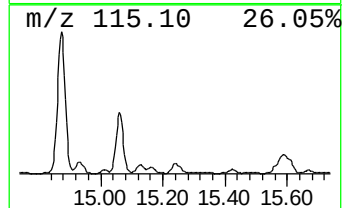
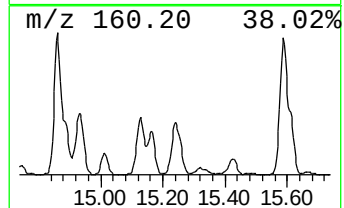
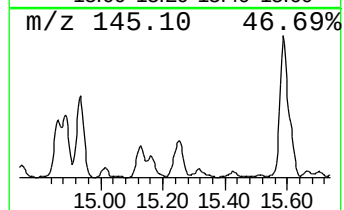
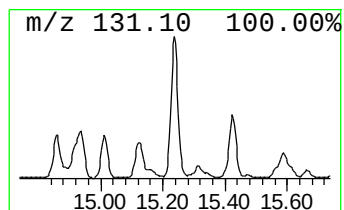
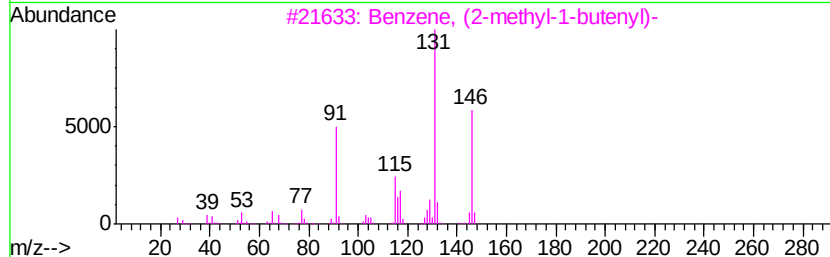
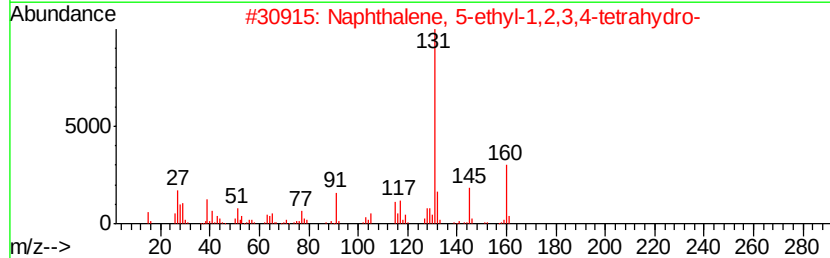
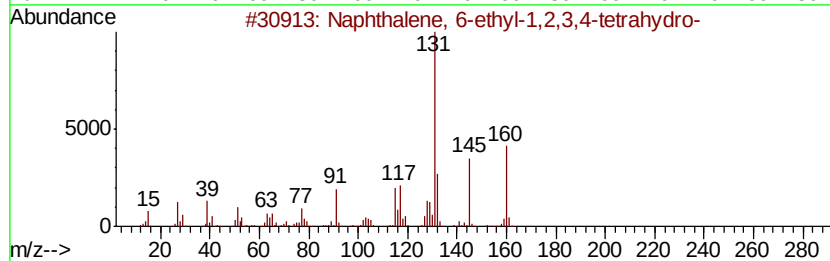
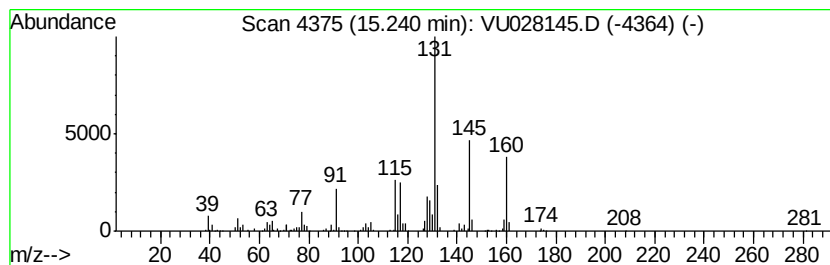
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 31 Naphthalene, 6-ethyl-1,2,3,... Concentration Rank 22

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 15.24 | 4.43 ug/L | 428144 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 6-ethyl-1,2,3,4-tet... | 160 | C12H16  | 022531-20-0 | 93   |
| 2    |    | Naphthalene, 5-ethyl-1,2,3,4-tet... | 160 | C12H16  | 042775-75-7 | 58   |
| 3    |    | Benzene, (2-methyl-1-butenyl)-      | 146 | C11H14  | 056253-64-6 | 53   |
| 4    |    | 1H-Indene, 2,3-dihydro-4-propyl-    | 160 | C12H16  | 092013-16-6 | 50   |
| 5    |    | 1,4-Methanonaphthalen-9-ol, 1,2,... | 160 | C11H12O | 055255-94-2 | 49   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

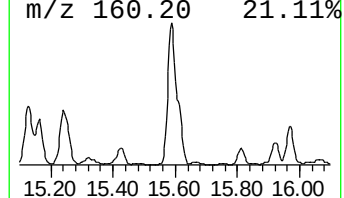
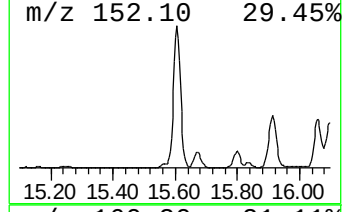
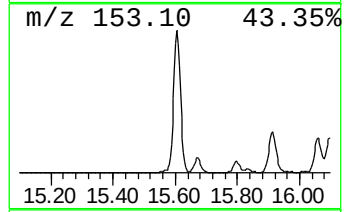
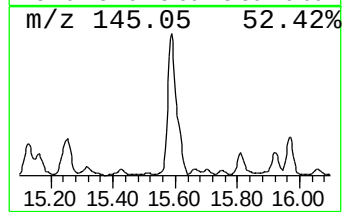
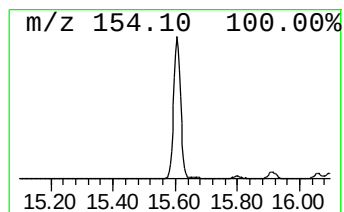
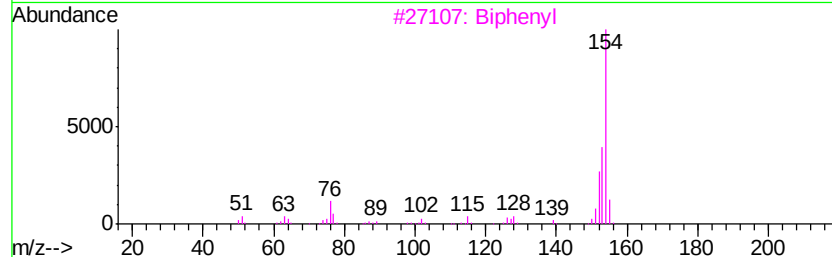
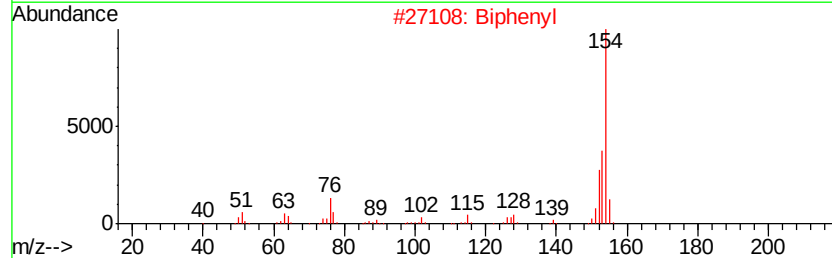
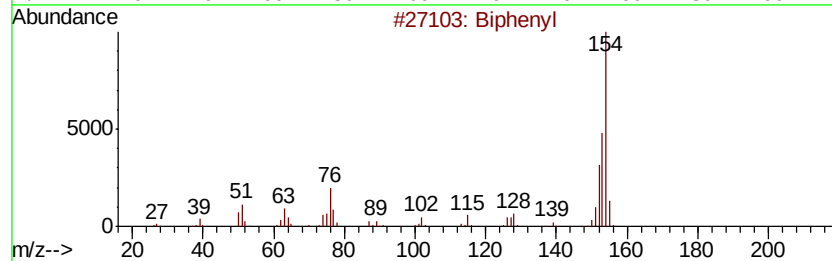
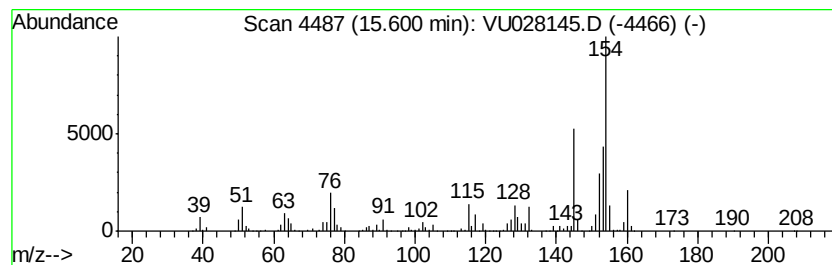
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 32 Biphenyl Concentration Rank 3

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 15.60 | 17.41 ug/L | 1682740 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Biphenyl                | 154 | C12H10  | 000092-52-4 | 95   |
| 2    |    | Biphenyl                | 154 | C12H10  | 000092-52-4 | 90   |
| 3    |    | Biphenyl                | 154 | C12H10  | 000092-52-4 | 60   |
| 4    |    | Naphthalene, 2-ethenyl- | 154 | C12H10  | 000827-54-3 | 60   |
| 5    |    | Naphthalene, 2-ethenyl- | 154 | C12H10  | 000827-54-3 | 60   |





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

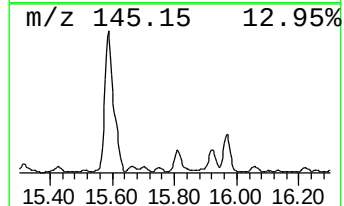
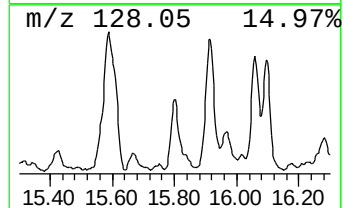
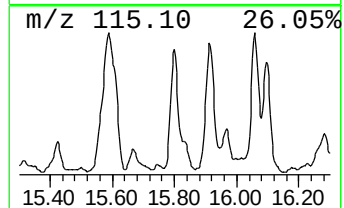
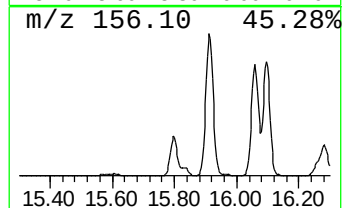
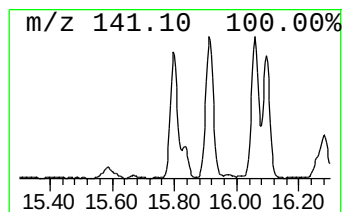
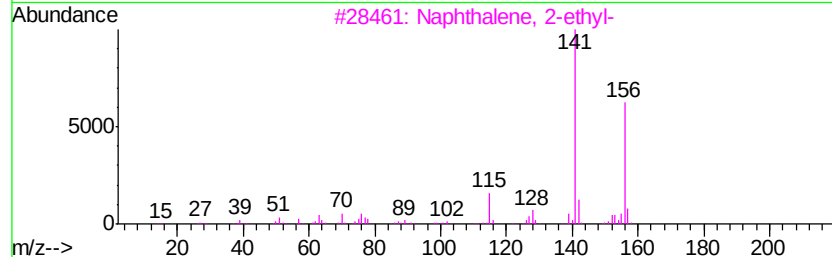
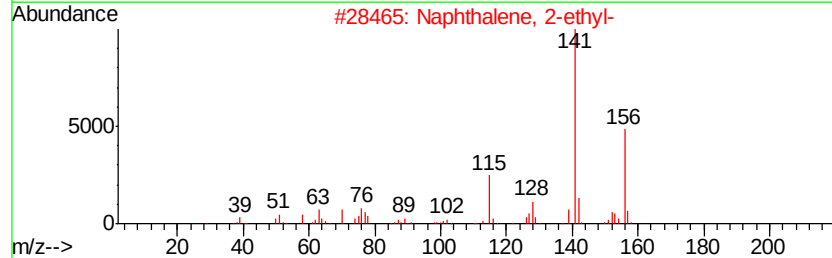
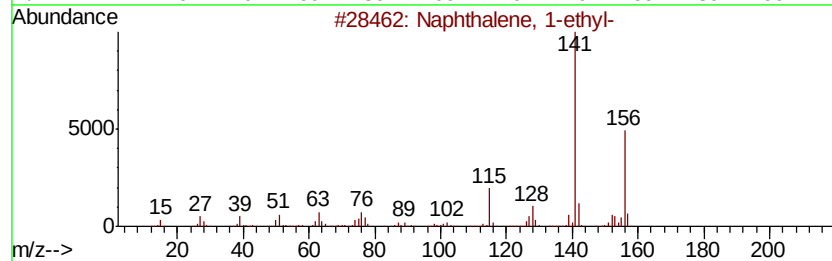
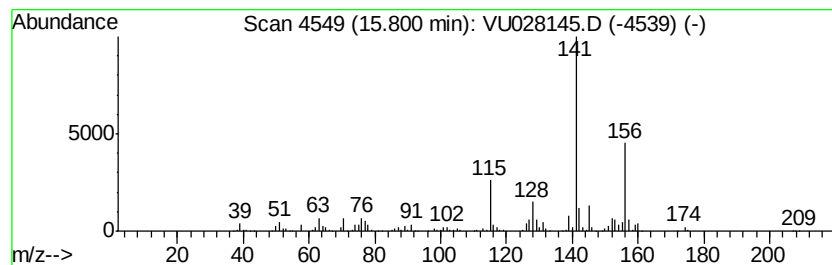
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 33 Naphthalene, 1-ethyl- Concentration Rank 16

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 15.80 | 5.77 ug/L | 558141 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1-ethyl-      | 156 | C12H12  | 001127-76-0 | 94   |
| 2    |    | Naphthalene, 2-ethyl-      | 156 | C12H12  | 000939-27-5 | 94   |
| 3    |    | Naphthalene, 2-ethyl-      | 156 | C12H12  | 000939-27-5 | 93   |
| 4    |    | Naphthalene, 1-ethyl-      | 156 | C12H12  | 001127-76-0 | 93   |
| 5    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU111418\  
Data File : VU028145.D  
Acq On : 15 Nov 2018 01:42  
Operator : MD/SY  
Sample : J5943-07  
Misc : 25.0mL/MSVOA U/WATER  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
H0FH6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.M  
Quant Title : TRACE VOA SOM01.0

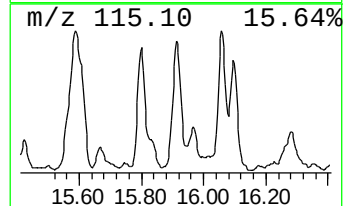
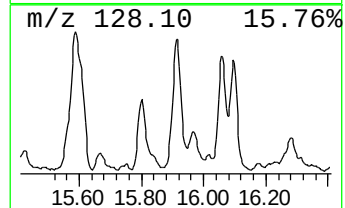
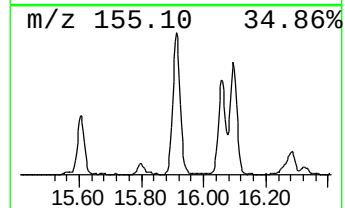
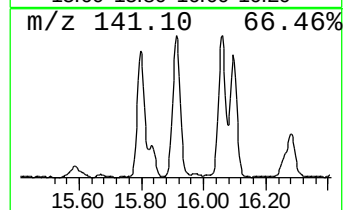
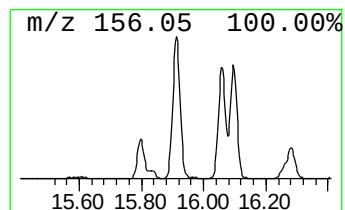
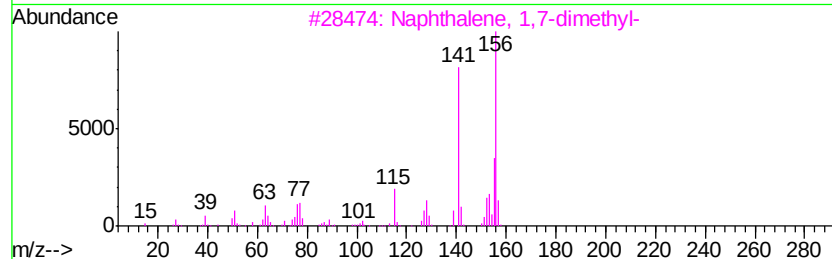
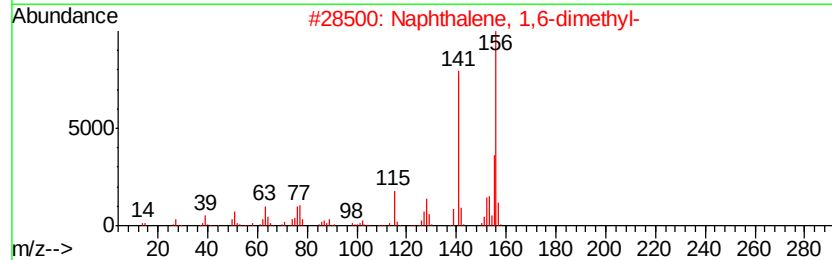
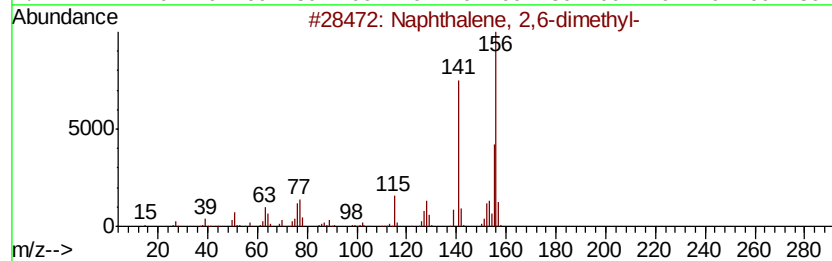
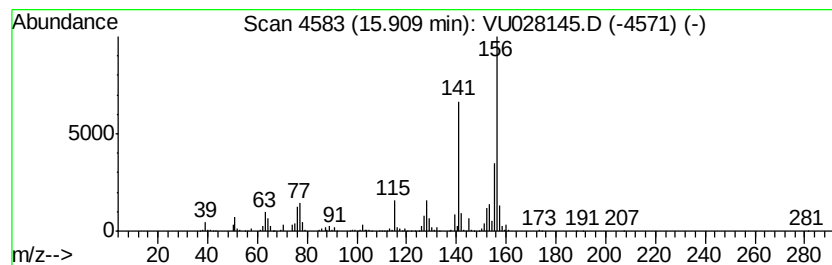
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 34 Naphthalene, 2,6-dimethyl- Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 15.91 | 10.29 ug/L | 995083 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 2    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 3    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 98   |
| 4    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 5    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU111418\  
 Data File : VU028145.D  
 Acq On : 15 Nov 2018 01:42  
 Operator : MD/SY  
 Sample : J5943-07  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 H0FH6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR111418WMA.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: Cyc... | 1.52  | 5.2     | ug/L  | 330132   | 1                     | 5.89  | 318870 | 5.0  |
| Ethane, 1,2-dichl... | 2.10  | 3.6     | ug/L  | 228245   | 1                     | 5.89  | 318870 | 5.0  |
| Diethyl disulfide    | 10.25 | 4.5     | ug/L  | 497394   | 2                     | 9.09  | 557502 | 5.0  |
| Benzene, propyl-     | 10.59 | 2.6     | ug/L  | 248414   | 3                     | 11.48 | 483292 | 5.0  |
| Benzene, 1-ethyl-... | 10.68 | 7.4     | ug/L  | 719416   | 3                     | 11.48 | 483292 | 5.0  |
| Mesitylene           | 10.77 | 4.5     | ug/L  | 429739   | 3                     | 11.48 | 483292 | 5.0  |
| 4-Heptanone, 2,6-... | 10.92 | 2.7     | ug/L  | 264908   | 3                     | 11.48 | 483292 | 5.0  |
| Benzene, 1,2,3-tr... | 11.15 | 16.6    | ug/L  | 1604730  | 3                     | 11.48 | 483292 | 5.0  |
| Benzene, 1,2,4-tr... | 11.57 | 11.7    | ug/L  | 1126640  | 3                     | 11.48 | 483292 | 5.0  |
| Indane               | 11.77 | 3.7     | ug/L  | 359447   | 3                     | 11.48 | 483292 | 5.0  |
| Benzene, 2-ethyl-... | 12.25 | 3.2     | ug/L  | 312730   | 3                     | 11.48 | 483292 | 5.0  |
| Benzene, (2-methy... | 12.35 | 2.8     | ug/L  | 271118   | 3                     | 11.48 | 483292 | 5.0  |
| Benzene, 1,2,3,5-... | 12.70 | 2.3     | ug/L  | 226039   | 3                     | 11.48 | 483292 | 5.0  |
| Benzene, 1-etheny... | 12.96 | 2.9     | ug/L  | 281928   | 3                     | 11.48 | 483292 | 5.0  |
| unknown-01           | 13.07 | 4.3     | ug/L  | 412696   | 3                     | 11.48 | 483292 | 5.0  |
| 1H-Indene, 2,3-di... | 13.12 | 8.3     | ug/L  | 800180   | 3                     | 11.48 | 483292 | 5.0  |
| Naphthalene, 1,2,... | 13.29 | 10.7    | ug/L  | 1030380  | 3                     | 11.48 | 483292 | 5.0  |
| Naphthalene, 1,2,... | 13.85 | 5.4     | ug/L  | 524688   | 3                     | 11.48 | 483292 | 5.0  |
| Benzene, (2-methy... | 13.95 | 7.1     | ug/L  | 684636   | 3                     | 11.48 | 483292 | 5.0  |
| 1H-Indene, 2,3-di... | 14.15 | 2.6     | ug/L  | 253260   | 3                     | 11.48 | 483292 | 5.0  |
| Naphthalene, 1,2,... | 14.35 | 17.5    | ug/L  | 1688660  | 3                     | 11.48 | 483292 | 5.0  |
| Benzene, (3-methy... | 14.54 | 3.7     | ug/L  | 359465   | 3                     | 11.48 | 483292 | 5.0  |
| Naphthalene, 1,2,... | 14.67 | 13.1    | ug/L  | 1266190  | 3                     | 11.48 | 483292 | 5.0  |
| Naphthalene, 1-me... | 14.87 | 33.0    | ug/L  | 3191800  | 3                     | 11.48 | 483292 | 5.0  |
| 1H-Indene, 2,3-di... | 14.93 | 4.9     | ug/L  | 475931   | 3                     | 11.48 | 483292 | 5.0  |
| Naphthalene, 1,2,... | 15.13 | 3.1     | ug/L  | 304943   | 3                     | 11.48 | 483292 | 5.0  |
| Naphthalene, 1,2,... | 15.16 | 2.2     | ug/L  | 216859   | 3                     | 11.48 | 483292 | 5.0  |
| Naphthalene, 6-et... | 15.24 | 4.4     | ug/L  | 428144   | 3                     | 11.48 | 483292 | 5.0  |
| Biphenyl             | 15.60 | 17.4    | ug/L  | 1682740  | 3                     | 11.48 | 483292 | 5.0  |
| Naphthalene, 1-et... | 15.80 | 5.8     | ug/L  | 558141   | 3                     | 11.48 | 483292 | 5.0  |
| Naphthalene, 2,6-... | 15.91 | 10.3    | ug/L  | 995083   | 3                     | 11.48 | 483292 | 5.0  |