

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
 Data File : VU031663.D  
 Acq On : 01 May 2019 03:21  
 Operator : JC/SP  
 Sample : K2604-06 5X  
 Misc : 5.01µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 GAJ80

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 0 % 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\SOMULM042719WMA.M  
 Title : VOC Analysis

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|--------|----------|------------|----------|-----------|-------|-------------|------------|--------------|------------|
| 1      | 1.180    | 24         | 28       | 33        | rBV3  | 15145       | 13072      | 0.01%        | 0.002%     |
| 2      | 1.395    | 88         | 95       | 108       | rBV   | 436021      | 494253     | 0.26%        | 0.086%     |
| 3      | 1.678    | 176        | 183      | 195       | rBV   | 321331      | 432418     | 0.23%        | 0.075%     |
| 4      | 2.260    | 354        | 364      | 378       | rVB   | 769557      | 1178284    | 0.62%        | 0.205%     |
| 5      | 2.460    | 423        | 426      | 430       | rBV4  | 1709        | 1728       | 0.00%        | 0.000%     |
| 6      | 2.537    | 448        | 450      | 454       | rVB5  | 2749        | 1833       | 0.00%        | 0.000%     |
| 7      | 2.688    | 490        | 497      | 499       | rBV5  | 4652        | 5202       | 0.00%        | 0.001%     |
| 8      | 2.817    | 533        | 537      | 540       | rBV4  | 1384        | 1477       | 0.00%        | 0.000%     |
| 9      | 2.836    | 540        | 543      | 551       | rVV8  | 2218        | 3111       | 0.00%        | 0.001%     |
| 10     | 2.981    | 583        | 588      | 591       | rBV6  | 1385        | 1335       | 0.00%        | 0.000%     |
| 11     | 3.309    | 688        | 690      | 698       | rVB4  | 1325        | 1444       | 0.00%        | 0.000%     |
| 12     | 3.479    | 739        | 743      | 749       | rVB6  | 1547        | 1608       | 0.00%        | 0.000%     |
| 13     | 3.804    | 840        | 844      | 851       | rVV   | 2539        | 2774       | 0.00%        | 0.000%     |
| 14     | 3.842    | 851        | 856      | 860       | rVV4  | 1685        | 1801       | 0.00%        | 0.000%     |
| 15     | 4.202    | 953        | 968      | 1021      | rBV2  | 221602      | 1066738    | 0.56%        | 0.186%     |
| 16     | 4.643    | 1086       | 1105     | 1133      | rBV   | 544237      | 1342749    | 0.71%        | 0.234%     |
| 17     | 4.858    | 1166       | 1172     | 1175      | rBV7  | 1437        | 1348       | 0.00%        | 0.000%     |
| 18     | 4.887    | 1175       | 1181     | 1183      | rBV7  | 1346        | 1595       | 0.00%        | 0.000%     |
| 19     | 4.958    | 1200       | 1203     | 1206      | rBV4  | 2323        | 2080       | 0.00%        | 0.000%     |
| 20     | 5.334    | 1295       | 1320     | 1331      | rBV2  | 1199888     | 3541232    | 1.87%        | 0.616%     |
| 21     | 5.617    | 1402       | 1408     | 1411      | rBV7  | 2750        | 3452       | 0.00%        | 0.001%     |
| 22     | 5.743    | 1442       | 1447     | 1453      | rVB9  | 1319        | 1361       | 0.00%        | 0.000%     |
| 23     | 5.788    | 1454       | 1461     | 1467      | rBV7  | 2198        | 2801       | 0.00%        | 0.000%     |
| 24     | 5.881    | 1476       | 1490     | 1533      | rBV   | 1147760     | 2468393    | 1.30%        | 0.430%     |
| 25     | 6.183    | 1571       | 1584     | 1600      | rBV10 | 22736       | 55984      | 0.03%        | 0.010%     |
| 26     | 6.328    | 1614       | 1629     | 1661      | rBV   | 752165      | 1614827    | 0.85%        | 0.281%     |
| 27     | 6.562    | 1699       | 1702     | 1710      | rVB8  | 1657        | 1994       | 0.00%        | 0.000%     |
| 28     | 6.681    | 1736       | 1739     | 1745      | rVB5  | 1722        | 1679       | 0.00%        | 0.000%     |
| 29     | 6.884    | 1799       | 1802     | 1805      | rBV4  | 2590        | 2208       | 0.00%        | 0.000%     |
| 30     | 7.061    | 1850       | 1857     | 1860      | rBV6  | 1719        | 2158       | 0.00%        | 0.000%     |
| 31     | 7.225    | 1896       | 1908     | 1937      | rBV2  | 383055      | 767964     | 0.41%        | 0.134%     |
| 32     | 7.427    | 1963       | 1971     | 1979      | rVB9  | 6890        | 11511      | 0.01%        | 0.002%     |
| 33     | 7.562    | 1995       | 2013     | 2027      | rBV   | 1398775     | 2567361    | 1.36%        | 0.447%     |
| 34     | 7.636    | 2029       | 2036     | 2058      | rVB   | 262026      | 508762     | 0.27%        | 0.089%     |

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
 Data File : VU031663.D  
 Acq On : 01 May 2019 03:21  
 Operator : JC/SP  
 Sample : K2604-06 5X  
 Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 GAJ80

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 0 % 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\SOMULM042719WMA.M  
 Title : VOC Analysis

|    |        |      |      |      |      |          |          |        |        |
|----|--------|------|------|------|------|----------|----------|--------|--------|
| 35 | 7.852  | 2086 | 2103 | 2134 | rVB2 | 244974   | 520612   | 0.27%  | 0.091% |
| 36 | 8.225  | 2212 | 2219 | 2227 | rBV9 | 5138     | 8709     | 0.00%  | 0.002% |
| 37 | 8.324  | 2237 | 2250 | 2288 | rBV  | 1076867  | 2362772  | 1.25%  | 0.411% |
| 38 | 8.595  | 2330 | 2334 | 2335 | rBV4 | 6758     | 4909     | 0.00%  | 0.001% |
| 39 | 8.746  | 2375 | 2381 | 2394 | rVB9 | 14532    | 26791    | 0.01%  | 0.005% |
| 40 | 8.813  | 2394 | 2402 | 2405 | rBV6 | 7010     | 9887     | 0.01%  | 0.002% |
| 41 | 9.086  | 2475 | 2487 | 2525 | rBV  | 1699427  | 3026932  | 1.60%  | 0.527% |
| 42 | 9.247  | 2526 | 2537 | 2562 | rBV  | 1093089  | 1871466  | 0.99%  | 0.326% |
| 43 | 9.369  | 2562 | 2575 | 2592 | rBV  | 8170241  | 13838756 | 7.30%  | 2.409% |
| 44 | 9.643  | 2656 | 2660 | 2661 | rBV4 | 6420     | 4690     | 0.00%  | 0.001% |
| 45 | 9.787  | 2690 | 2705 | 2740 | rBV3 | 13915223 | 25661030 | 13.54% | 4.466% |
| 46 | 10.164 | 2815 | 2822 | 2830 | rBV  | 46475    | 75332    | 0.04%  | 0.013% |
| 47 | 10.430 | 2895 | 2905 | 2917 | rBV  | 1027763  | 1688947  | 0.89%  | 0.294% |
| 48 | 10.508 | 2922 | 2929 | 2942 | rVB2 | 221863   | 364763   | 0.19%  | 0.063% |
| 49 | 10.585 | 2944 | 2953 | 2970 | rBV  | 286250   | 450038   | 0.24%  | 0.078% |
| 50 | 10.678 | 2972 | 2982 | 2998 | rBV  | 1143194  | 2479615  | 1.31%  | 0.432% |
| 51 | 10.768 | 2999 | 3010 | 3019 | rBV  | 1503297  | 2486919  | 1.31%  | 0.433% |
| 52 | 10.993 | 3063 | 3080 | 3094 | rBV2 | 982461   | 2089505  | 1.10%  | 0.364% |
| 53 | 11.144 | 3110 | 3127 | 3138 | rBV  | 5052383  | 8252228  | 4.36%  | 1.436% |
| 54 | 11.212 | 3138 | 3148 | 3157 | rVV  | 8559381  | 12903877 | 6.81%  | 2.246% |
| 55 | 11.263 | 3157 | 3164 | 3177 | rVB  | 3107123  | 4684745  | 2.47%  | 0.815% |
| 56 | 11.398 | 3200 | 3206 | 3211 | rBV3 | 108656   | 162423   | 0.09%  | 0.028% |
| 57 | 11.482 | 3222 | 3232 | 3243 | rBV  | 1883573  | 2921186  | 1.54%  | 0.508% |
| 58 | 11.565 | 3250 | 3258 | 3268 | rBV  | 1603926  | 2392406  | 1.26%  | 0.416% |
| 59 | 11.623 | 3268 | 3276 | 3293 | rVB  | 1179945  | 1816023  | 0.96%  | 0.316% |
| 60 | 11.762 | 3309 | 3319 | 3328 | rBV  | 1136529  | 1904003  | 1.01%  | 0.331% |
| 61 | 11.858 | 3340 | 3349 | 3365 | rVB3 | 1709108  | 3099343  | 1.64%  | 0.539% |
| 62 | 11.977 | 3374 | 3386 | 3398 | rBV2 | 30496905 | 47923470 | 25.30% | 8.341% |
| 63 | 12.144 | 3430 | 3438 | 3441 | rBV2 | 233935   | 327161   | 0.17%  | 0.057% |
| 64 | 12.218 | 3454 | 3461 | 3466 | rBV2 | 470708   | 728979   | 0.38%  | 0.127% |
| 65 | 12.421 | 3515 | 3524 | 3535 | rVB  | 2022180  | 3134361  | 1.65%  | 0.546% |
| 66 | 12.482 | 3535 | 3543 | 3556 | rVB2 | 535765   | 829554   | 0.44%  | 0.144% |
| 67 | 12.646 | 3588 | 3594 | 3601 | rVB  | 379933   | 522070   | 0.28%  | 0.091% |
| 68 | 12.704 | 3602 | 3612 | 3623 | rBV4 | 1394693  | 2816394  | 1.49%  | 0.490% |
| 69 | 12.819 | 3639 | 3648 | 3658 | rBV  | 1481061  | 2433955  | 1.28%  | 0.424% |
| 70 | 12.961 | 3685 | 3692 | 3705 | rVB  | 362408   | 528213   | 0.28%  | 0.092% |
| 71 | 13.080 | 3723 | 3729 | 3732 | rBV  | 236700   | 286265   | 0.15%  | 0.050% |

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
 Data File : VU031663.D  
 Acq On : 01 May 2019 03:21  
 Operator : JC/SP  
 Sample : K2604-06 5X  
 Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 GAJ80

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 0 % 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\SOMULM042719WMA.M  
 Title : VOC Analysis

|     |        |      |      |      |      |           |           |         |         |
|-----|--------|------|------|------|------|-----------|-----------|---------|---------|
| 72  | 13.118 | 3733 | 3741 | 3750 | rVV4 | 1611610   | 2603237   | 1.37%   | 0.453%  |
| 73  | 13.180 | 3750 | 3760 | 3773 | rVB  | 6473855   | 9757625   | 5.15%   | 1.698%  |
| 74  | 13.286 | 3782 | 3793 | 3809 | rVB  | 6929653   | 11333115  | 5.98%   | 1.973%  |
| 75  | 13.382 | 3813 | 3823 | 3840 | rVB3 | 978145    | 1838628   | 0.97%   | 0.320%  |
| 76  | 13.752 | 3921 | 3938 | 3961 | rBV2 | 100770577 | 189451898 | 100.00% | 32.974% |
| 77  | 13.858 | 3963 | 3971 | 3992 | rVB  | 1540926   | 2591740   | 1.37%   | 0.451%  |
| 78  | 14.138 | 4052 | 4058 | 4064 | rBV2 | 330255    | 457383    | 0.24%   | 0.080%  |
| 79  | 14.334 | 4110 | 4119 | 4128 | rBV2 | 892478    | 1572594   | 0.83%   | 0.274%  |
| 80  | 14.459 | 4153 | 4158 | 4175 | rVB2 | 745399    | 1534395   | 0.81%   | 0.267%  |
| 81  | 14.581 | 4190 | 4196 | 4209 | rVB  | 406141    | 615416    | 0.32%   | 0.107%  |
| 82  | 14.726 | 4235 | 4241 | 4258 | rVB  | 779425    | 1318900   | 0.70%   | 0.230%  |
| 83  | 14.816 | 4262 | 4269 | 4275 | rVV  | 388404    | 586604    | 0.31%   | 0.102%  |
| 84  | 14.874 | 4275 | 4287 | 4312 | rVB3 | 46468793  | 72264392  | 38.14%  | 12.577% |
| 85  | 15.057 | 4332 | 4344 | 4363 | rBV  | 28094510  | 43799311  | 23.12%  | 7.623%  |
| 86  | 15.601 | 4503 | 4513 | 4524 | rBV  | 4622113   | 6934366   | 3.66%   | 1.207%  |
| 87  | 15.794 | 4566 | 4573 | 4579 | rBV  | 946804    | 1234477   | 0.65%   | 0.215%  |
| 88  | 15.906 | 4597 | 4608 | 4626 | rBV  | 5549011   | 9809887   | 5.18%   | 1.707%  |
| 89  | 16.054 | 4644 | 4654 | 4661 | rBV  | 7643734   | 12065452  | 6.37%   | 2.100%  |
| 90  | 16.183 | 4684 | 4694 | 4706 | rVV  | 3316963   | 5084729   | 2.68%   | 0.885%  |
| 91  | 16.279 | 4708 | 4724 | 4746 | rVB  | 2153115   | 5534705   | 2.92%   | 0.963%  |
| 92  | 16.427 | 4761 | 4770 | 4786 | rVB  | 1275689   | 2028851   | 1.07%   | 0.353%  |
| 93  | 16.517 | 4787 | 4798 | 4817 | rVV  | 8198284   | 13758085  | 7.26%   | 2.395%  |
| 94  | 16.626 | 4827 | 4832 | 4844 | rVB2 | 550213    | 811159    | 0.43%   | 0.141%  |
| 95  | 16.732 | 4856 | 4865 | 4872 | rBV  | 901822    | 1510330   | 0.80%   | 0.263%  |
| 96  | 16.967 | 4932 | 4938 | 4945 | rVB  | 590758    | 812401    | 0.43%   | 0.141%  |
| 97  | 17.012 | 4946 | 4952 | 4977 | rVB3 | 753736    | 1679708   | 0.89%   | 0.292%  |
| 98  | 17.160 | 4990 | 4998 | 5009 | rVB2 | 874480    | 1380069   | 0.73%   | 0.240%  |
| 99  | 17.221 | 5010 | 5017 | 5028 | rVB2 | 493525    | 769755    | 0.41%   | 0.134%  |
| 100 | 17.527 | 5103 | 5112 | 5125 | rBV2 | 355587    | 669915    | 0.35%   | 0.117%  |

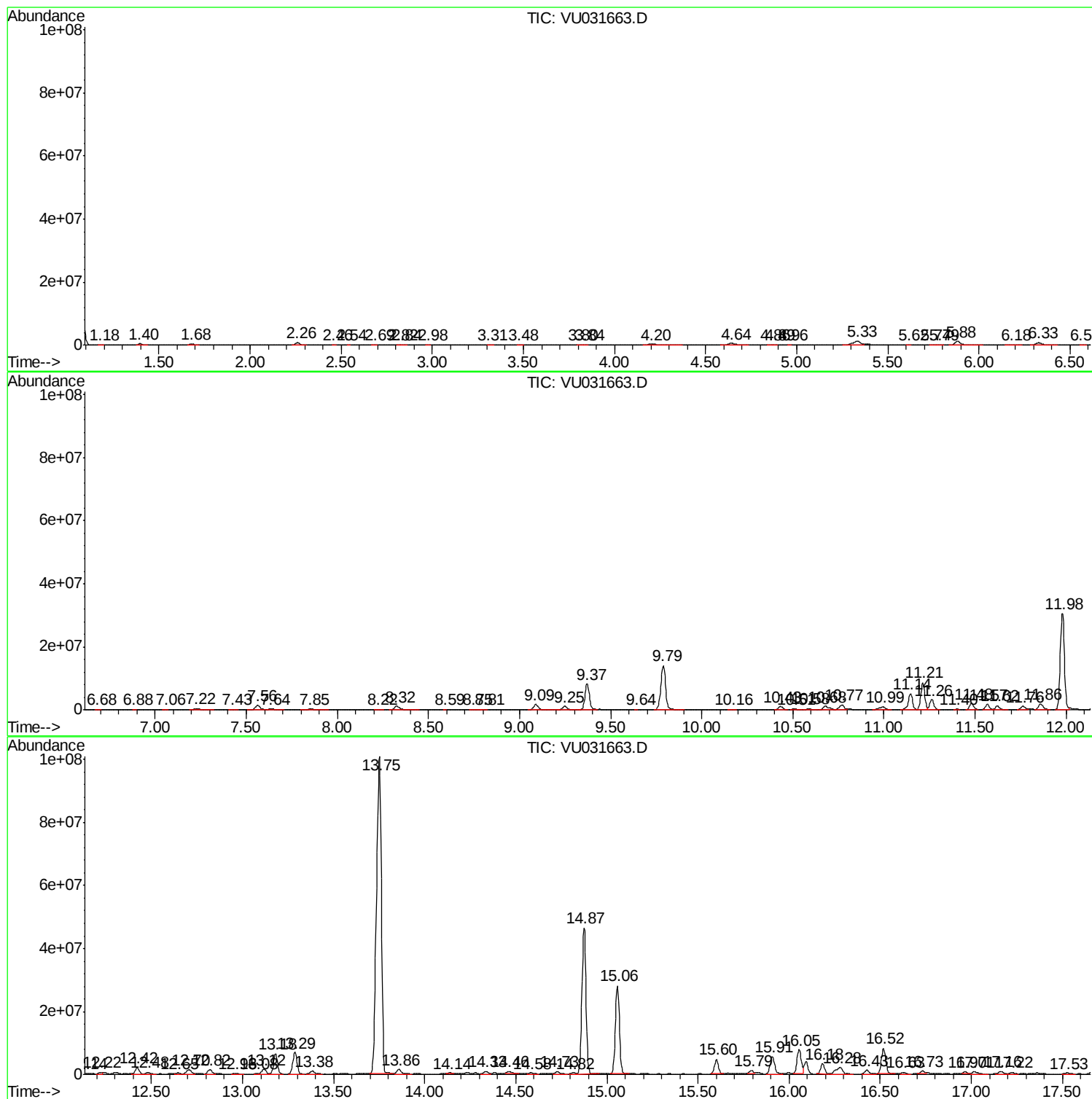
Sum of corrected areas: 574553993

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

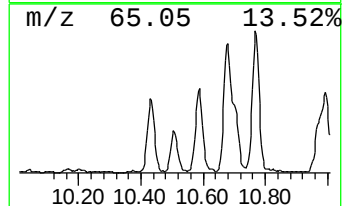
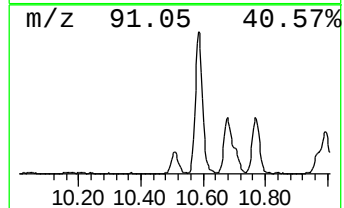
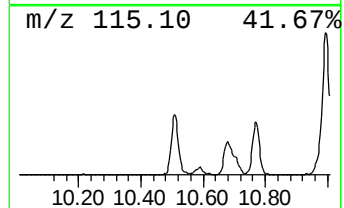
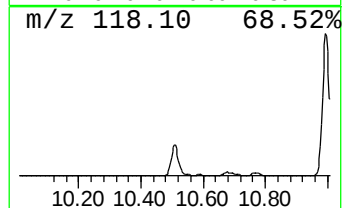
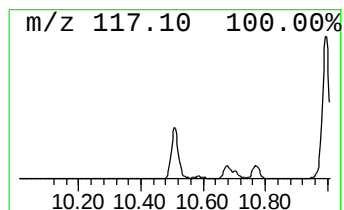
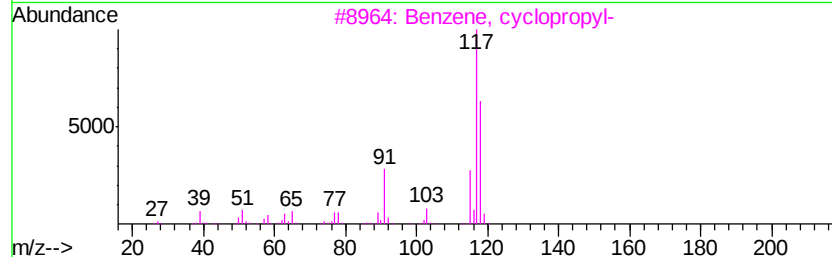
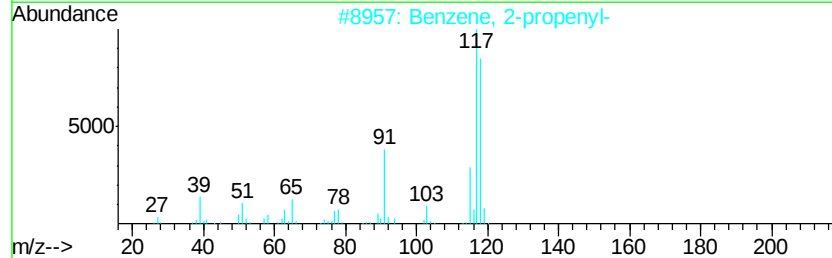
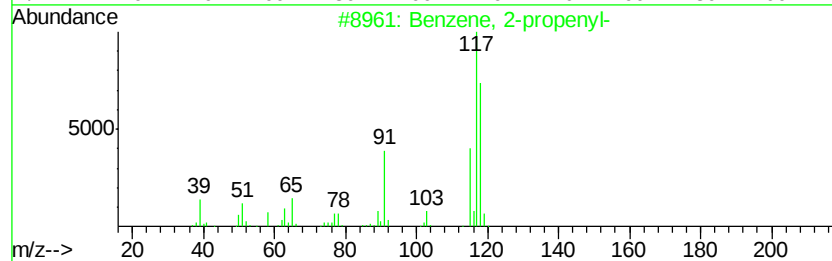
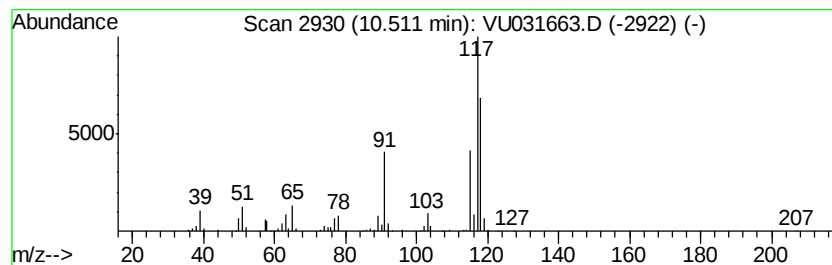
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 2 Benzene, 2-propenyl- Concentration Rank 49

| R.T.      | EstConc                             | Area   | Relative to ISTD       | R.T.         |      |
|-----------|-------------------------------------|--------|------------------------|--------------|------|
| 10.51     | 6.24 ug/L                           | 364763 | 1,4-Dichlorobenzene-d4 | 11.48        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm                | CAS#         | Qual |
| 1         | Benzene, 2-propenyl-                | 118    | C9H10                  | 000300-57-2  | 96   |
| 2         | Benzene, 2-propenyl-                | 118    | C9H10                  | 000300-57-2  | 96   |
| 3         | Benzene, cyclopropyl-               | 118    | C9H10                  | 000873-49-4  | 94   |
| 4         | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118    | C9H10                  | 1000191-13-7 | 94   |
| 5         | Benzene, 2-propenyl-                | 118    | C9H10                  | 000300-57-2  | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

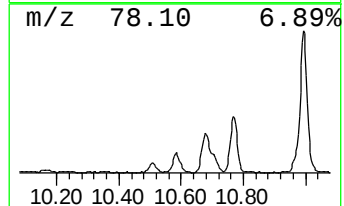
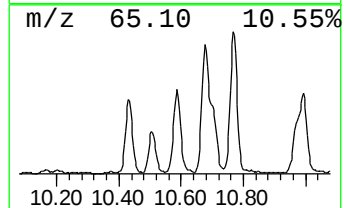
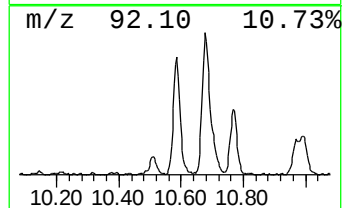
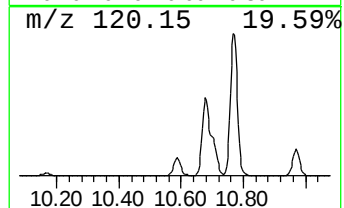
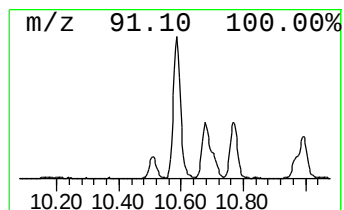
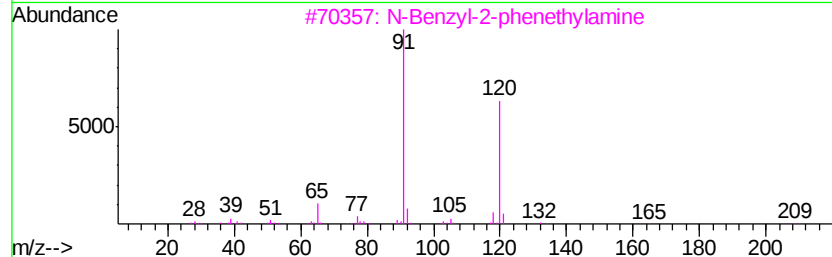
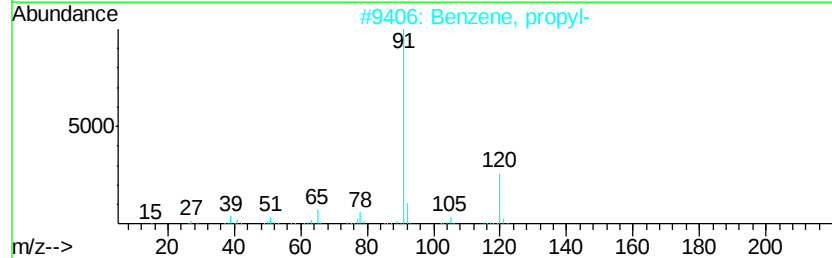
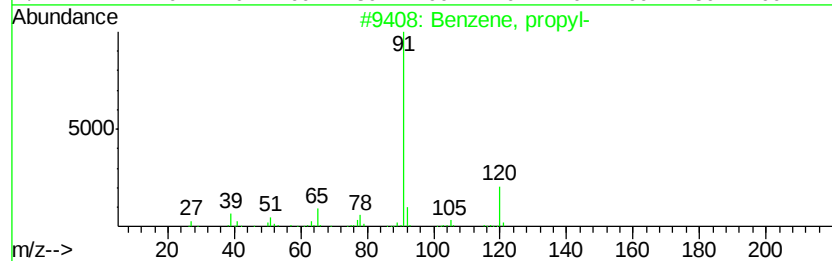
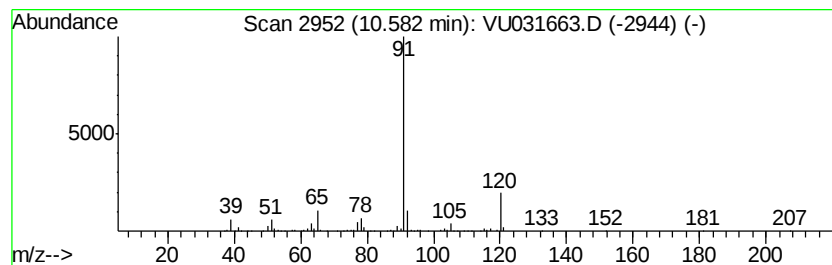
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 3 Benzene, propyl- Concentration Rank 48

| R.T.    | EstConc   | Area                              | Relative to ISTD       | R.T.           |
|---------|-----------|-----------------------------------|------------------------|----------------|
| 10.58   | 7.70 ug/L | 450038                            | 1,4-Dichlorobenzene-d4 | 11.48          |
| Hit# of | 5         | Tentative ID                      | MW MolForm             | CAS# Qual      |
| 1       |           | Benzene, propyl-                  | 120 C9H12              | 000103-65-1 90 |
| 2       |           | Benzene, propyl-                  | 120 C9H12              | 000103-65-1 87 |
| 3       |           | N-Benzyl-2-phenethylamine         | 211 C15H17N            | 003647-71-0 78 |
| 4       |           | Ethanol, 2-[(phenylmethyl)amino]- | 151 C9H13NO            | 000104-63-2 72 |
| 5       |           | Benzene, propyl-                  | 120 C9H12              | 000103-65-1 50 |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

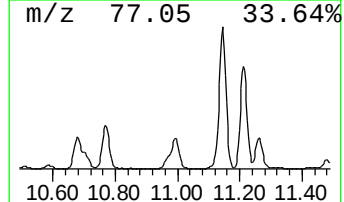
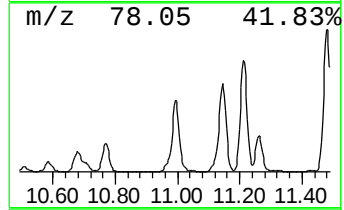
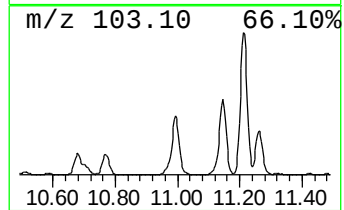
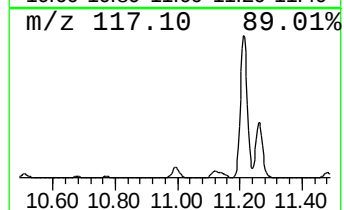
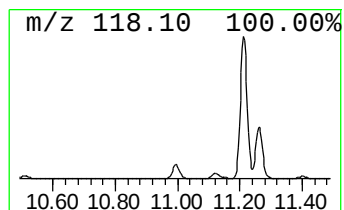
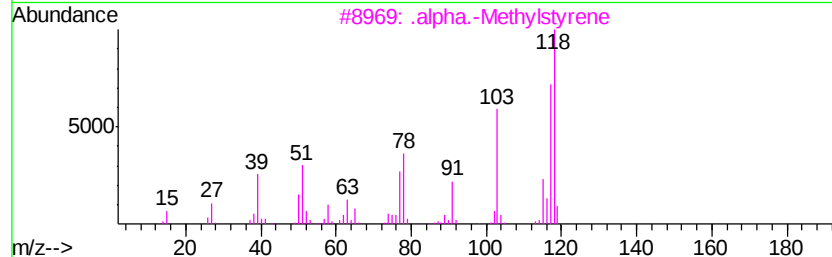
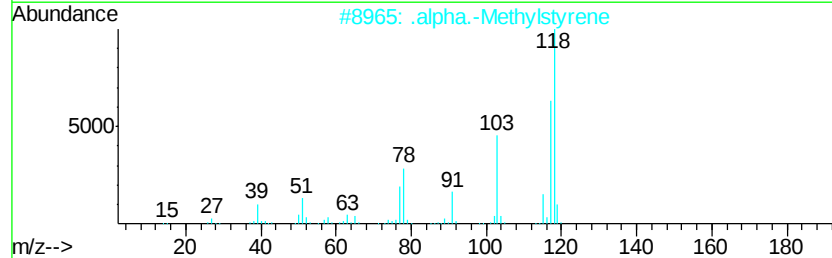
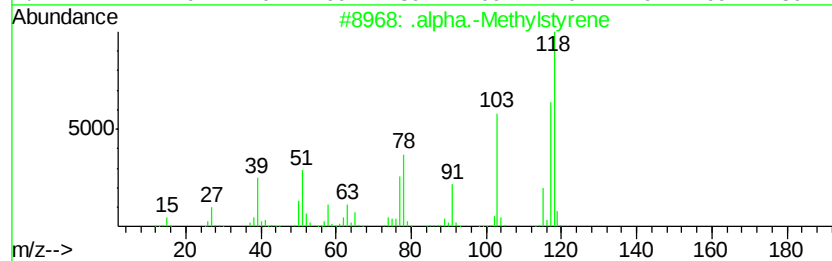
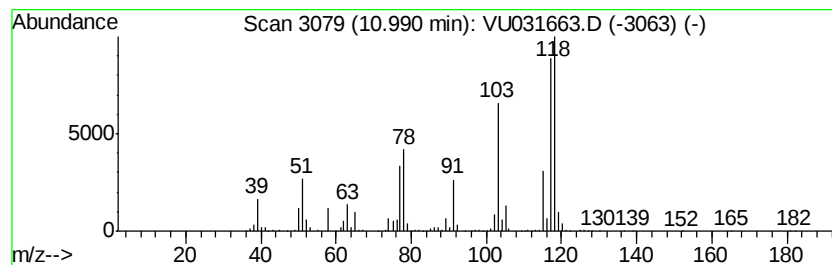
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 6 .alpha.-Methylstyrene Concentration Rank 24

| R.T.      | EstConc                      | Area    | Relative to ISTD       | R.T.        |      |
|-----------|------------------------------|---------|------------------------|-------------|------|
| 10.99     | 35.76 ug/L                   | 2089510 | 1,4-Dichlorobenzene-d4 | 11.48       |      |
| Hit# of 5 | Tentative ID                 | MW      | MolForm                | CAS#        | Qual |
| 1         | .alpha.-Methylstyrene        | 118     | C9H10                  | 000098-83-9 | 96   |
| 2         | .alpha.-Methylstyrene        | 118     | C9H10                  | 000098-83-9 | 93   |
| 3         | .alpha.-Methylstyrene        | 118     | C9H10                  | 000098-83-9 | 93   |
| 4         | Benzene, 1-ethenyl-4-methyl- | 118     | C9H10                  | 000622-97-9 | 49   |
| 5         | Benzene, 1-ethenyl-3-methyl- | 118     | C9H10                  | 000100-80-1 | 49   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

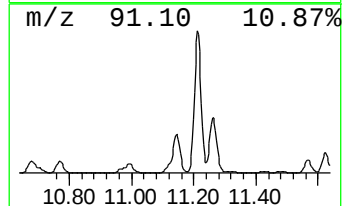
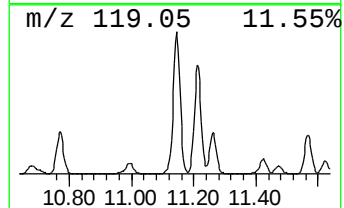
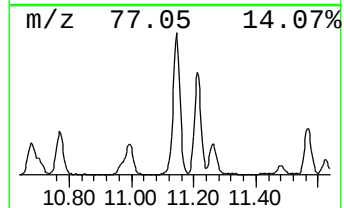
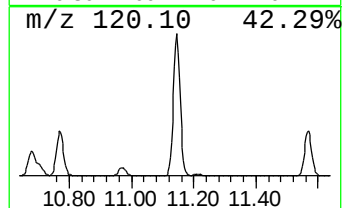
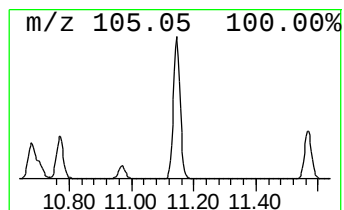
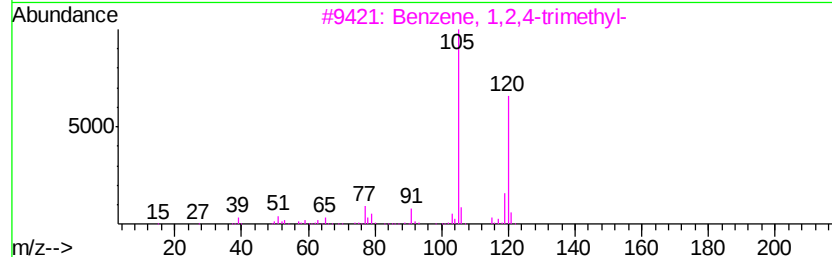
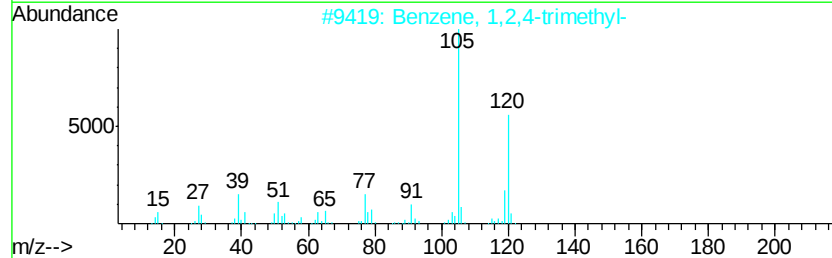
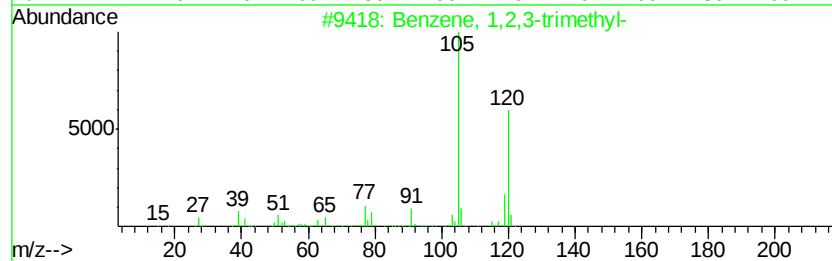
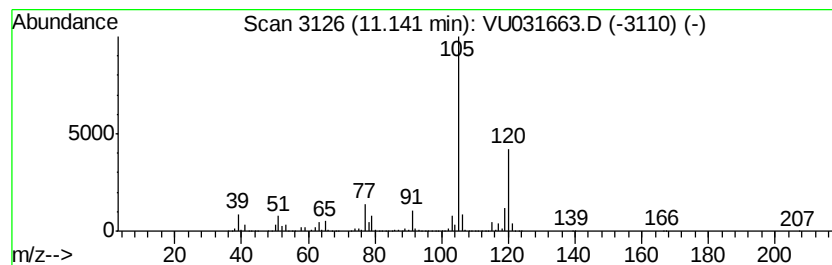
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 11

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 11.14 | 141.25 ug/L | 8252230 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 95   |
| 3    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 4    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 5    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

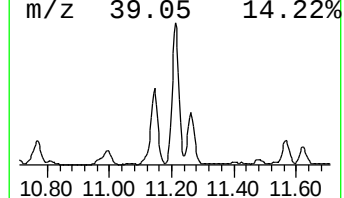
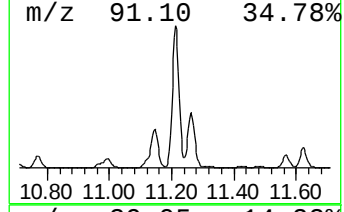
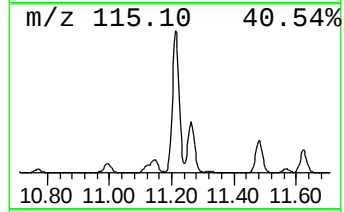
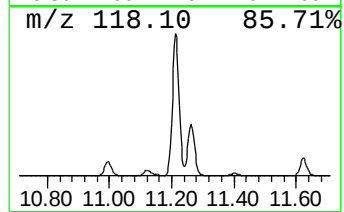
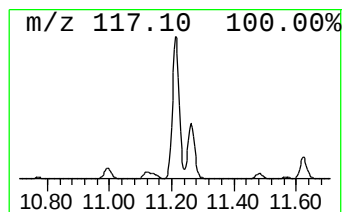
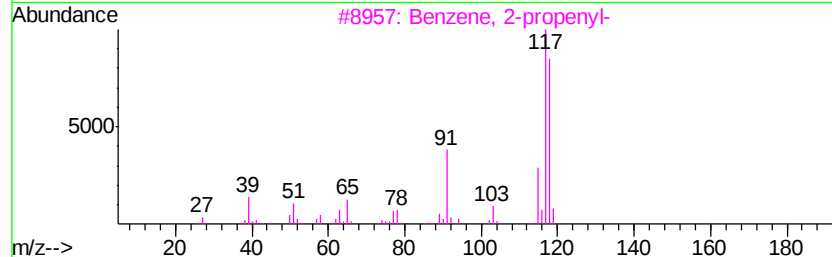
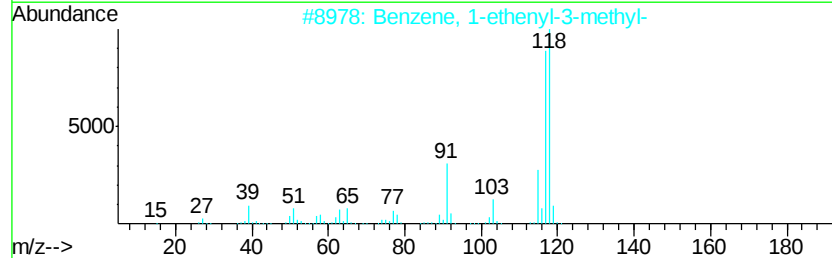
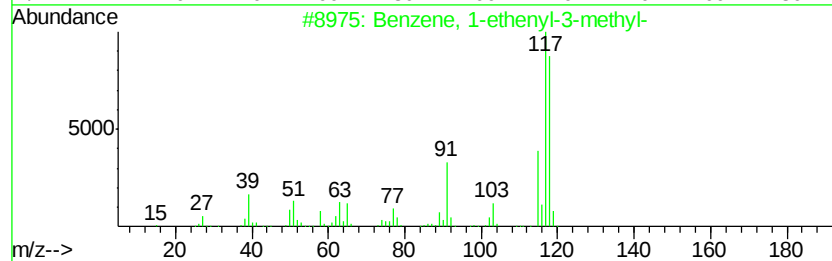
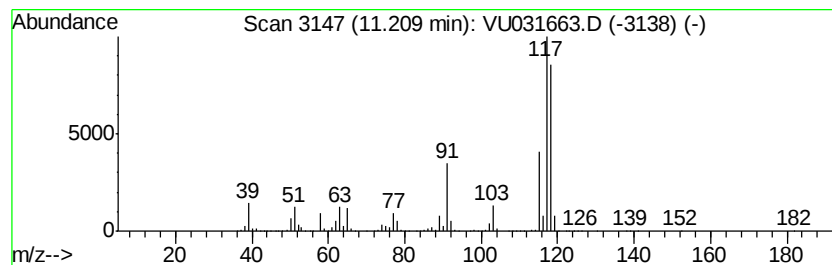
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Benzene, 1-ethenyl-3-methyl- Concentration Rank 6

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 11.21 | 220.87 ug/L | 12903900 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 96   |
| 2    |    | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 96   |
| 3    |    | Benzene, 2-propenyl-         | 118 | C9H10   | 000300-57-2 | 95   |
| 4    |    | Benzene, 1-propenyl-         | 118 | C9H10   | 000637-50-3 | 94   |
| 5    |    | Benzene, 1-ethenyl-2-methyl- | 118 | C9H10   | 000611-15-4 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

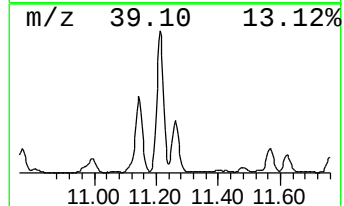
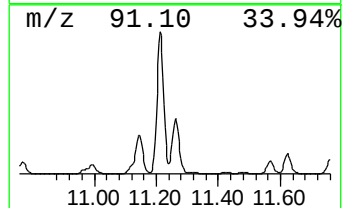
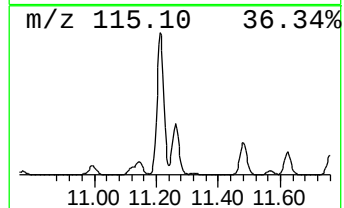
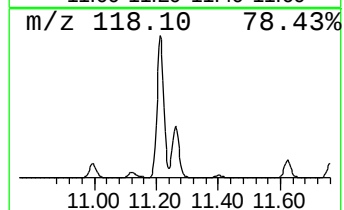
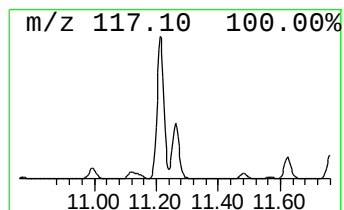
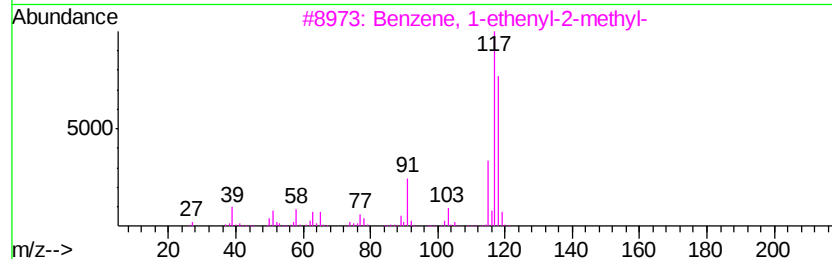
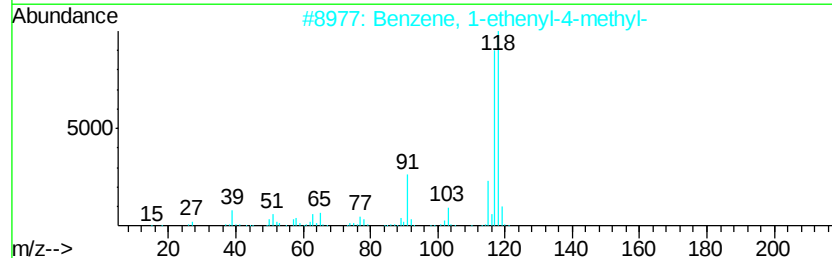
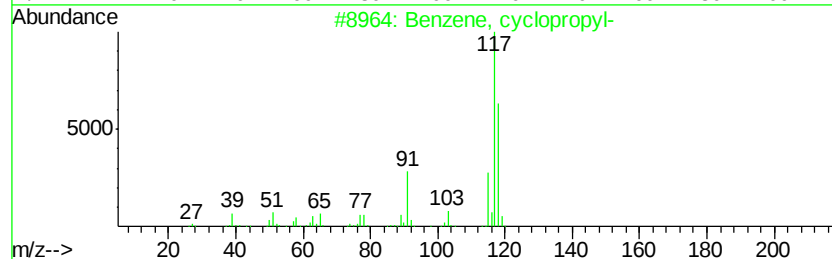
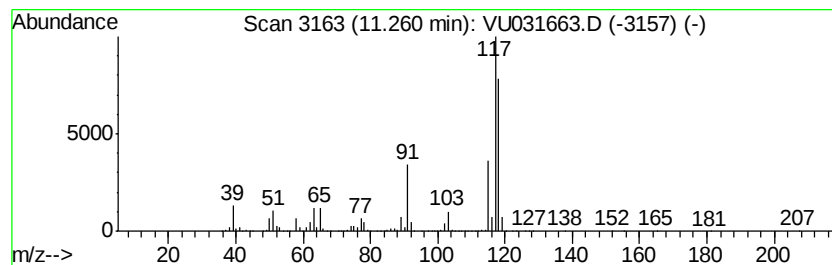
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Benzene, cyclopropyl- Concentration Rank 15

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 11.26 | 80.19 ug/L | 4684750 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, cvclopropyl-        | 118 | C9H10   | 000873-49-4 | 94   |
| 2    |    | Benzene, 1-ethenyl-4-methyl- | 118 | C9H10   | 000622-97-9 | 93   |
| 3    |    | Benzene, 1-ethenyl-2-methyl- | 118 | C9H10   | 000611-15-4 | 93   |
| 4    |    | Benzene, 2-propenyl-         | 118 | C9H10   | 000300-57-2 | 93   |
| 5    |    | Benzene, 1-propenyl-         | 118 | C9H10   | 000637-50-3 | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

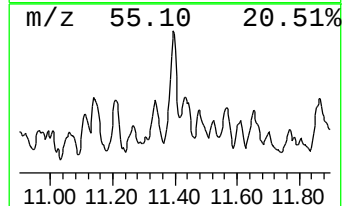
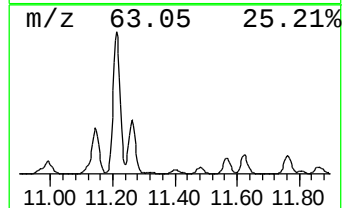
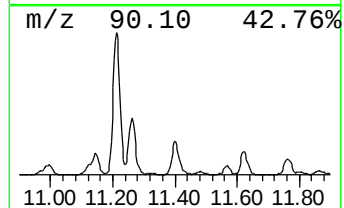
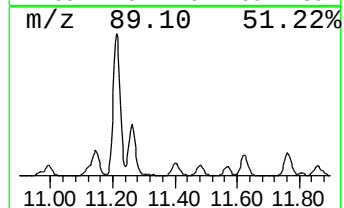
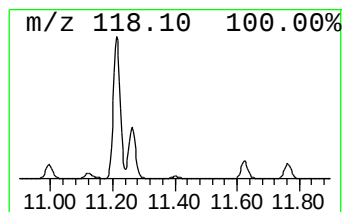
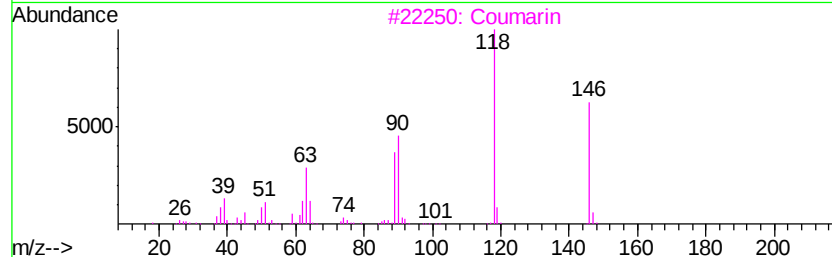
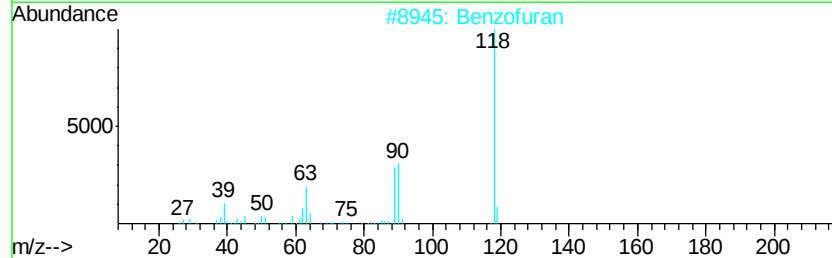
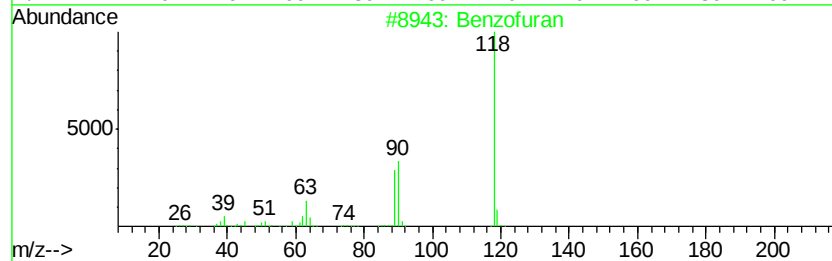
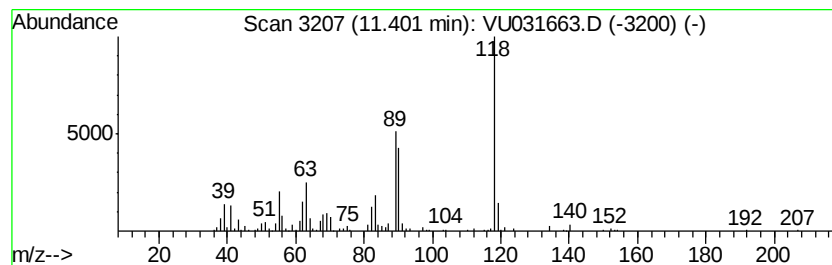
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 10 Benzofuran Concentration Rank 52

| R.T.    | EstConc   | Area                     | Relative to ISTD       | R.T.           |
|---------|-----------|--------------------------|------------------------|----------------|
| 11.40   | 2.78 ug/L | 162423                   | 1,4-Dichlorobenzene-d4 | 11.48          |
| Hit# of | 5         | Tentative ID             | MW MolForm             | CAS# Qual      |
| 1       |           | Benzofuran               | 118 C8H6O              | 000271-89-6 93 |
| 2       |           | Benzofuran               | 118 C8H6O              | 000271-89-6 87 |
| 3       |           | Coumarin                 | 146 C9H6O2             | 000091-64-5 72 |
| 4       |           | Coumarin                 | 146 C9H6O2             | 000091-64-5 72 |
| 5       |           | 3-Hydroxyphenylacetylene | 118 C8H6O              | 010401-11-3 68 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

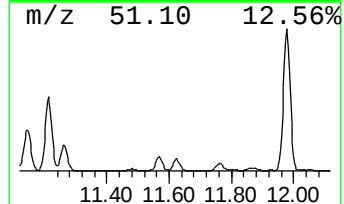
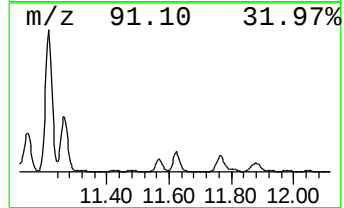
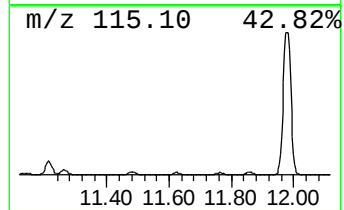
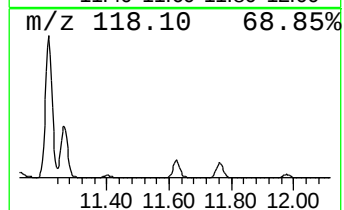
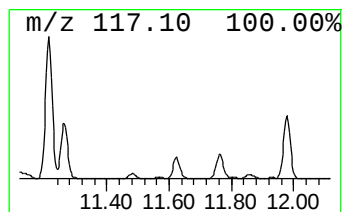
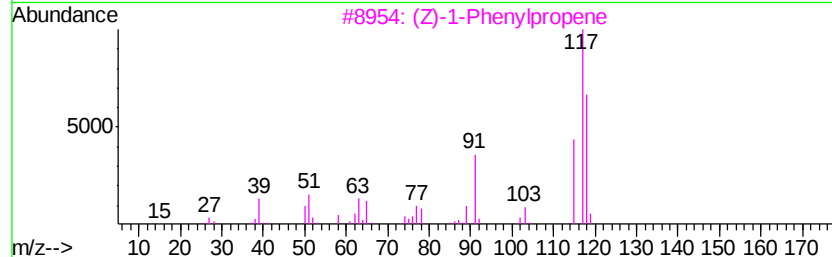
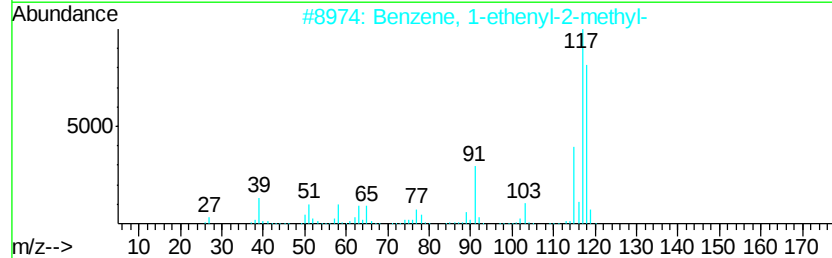
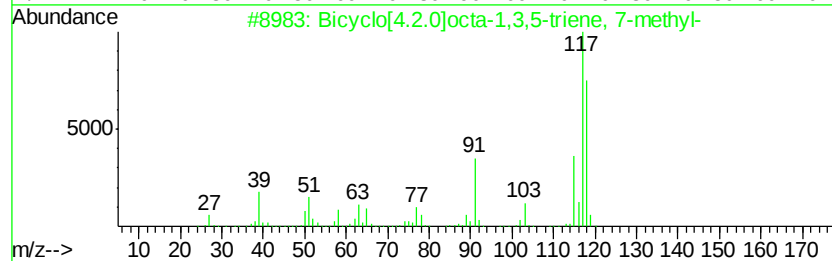
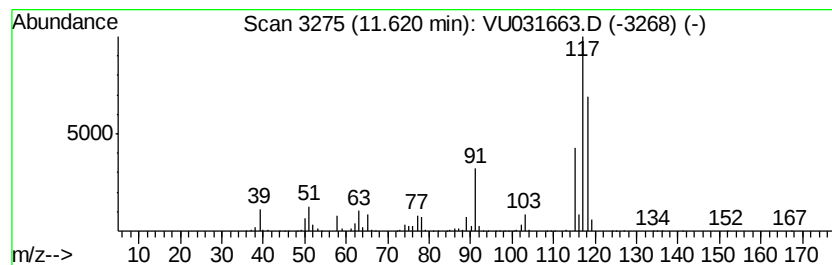
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 28

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 11.62 | 31.08 ug/L | 1816020 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Bicyclo[4.2.0]octa-1,3,5-triene,... | 118 | C9H10   | 055337-80-9 | 96   |
| 2    |    | Benzene, 1-ethenyl-2-methyl-        | 118 | C9H10   | 000611-15-4 | 95   |
| 3    |    | (Z)-1-Phenylpropene                 | 118 | C9H10   | 000766-90-5 | 95   |
| 4    |    | Benzene, cyclopropyl-               | 118 | C9H10   | 000873-49-4 | 95   |
| 5    |    | Benzene, 1-propenyl-                | 118 | C9H10   | 000637-50-3 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

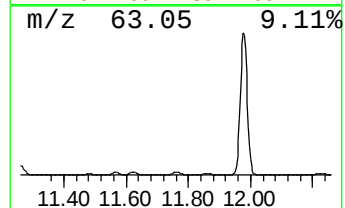
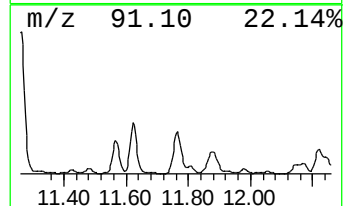
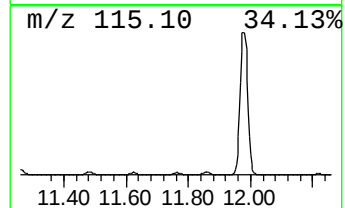
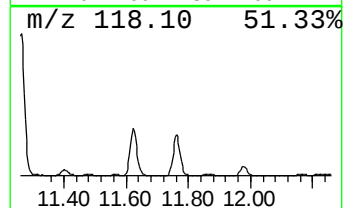
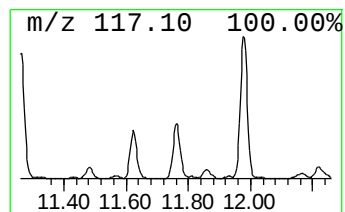
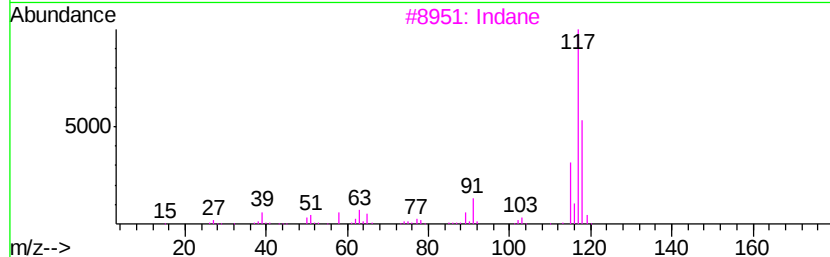
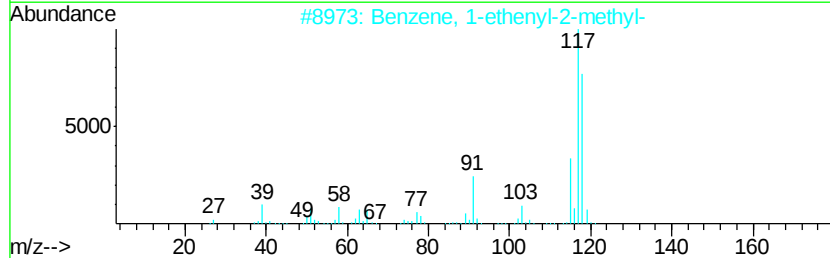
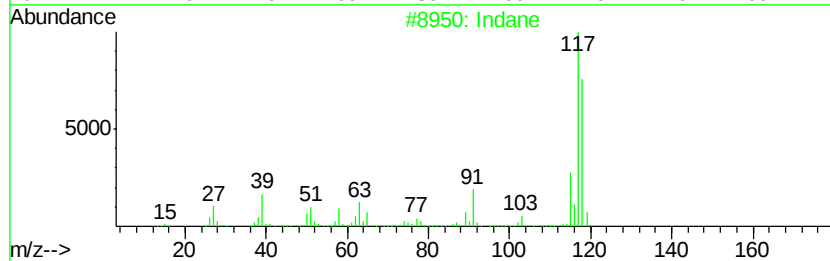
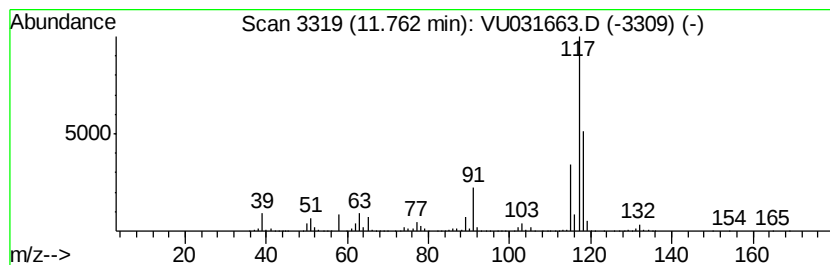
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 13 Indane Concentration Rank 26

| R.T.    | EstConc                             | Area         | Relative to ISTD       | R.T.            |
|---------|-------------------------------------|--------------|------------------------|-----------------|
| 11.76   | 32.59 ug/L                          | 1904000      | 1,4-Dichlorobenzene-d4 | 11.48           |
| Hit# of | 5                                   | Tentative ID | MW MolForm             | CAS# Qual       |
| 1       | Indane                              |              | 118 C9H10              | 000496-11-7 91  |
| 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 76  |
| 5       | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... |              | 118 C9H10              | 1000191-13-7 74 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

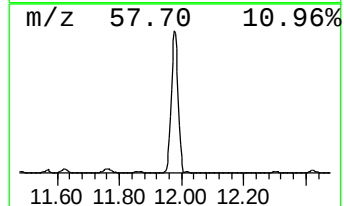
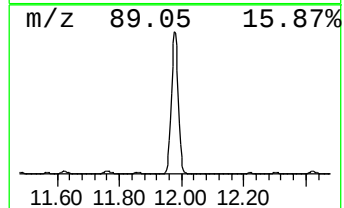
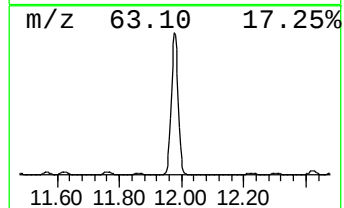
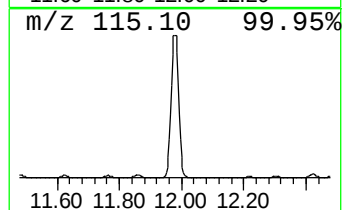
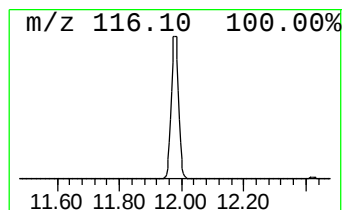
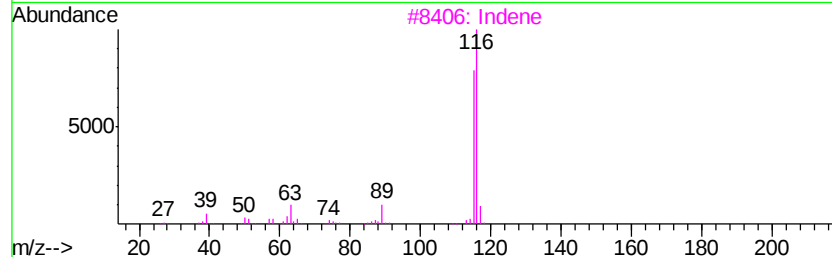
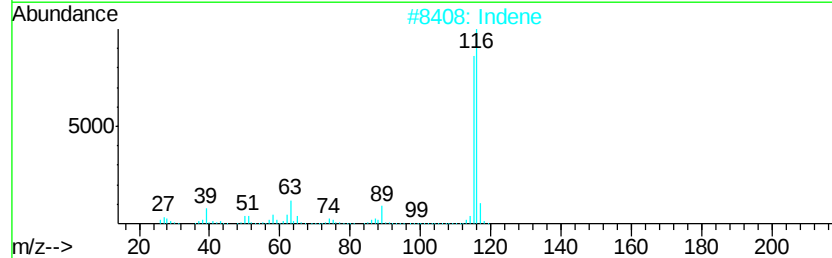
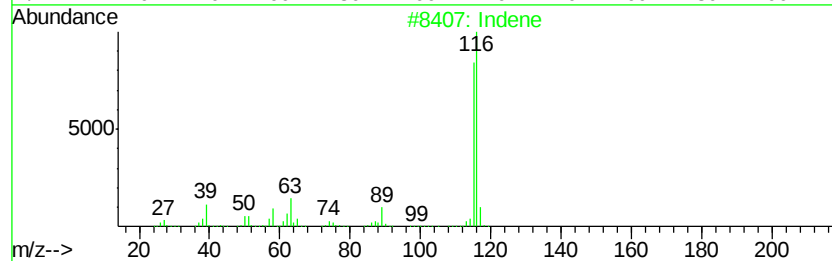
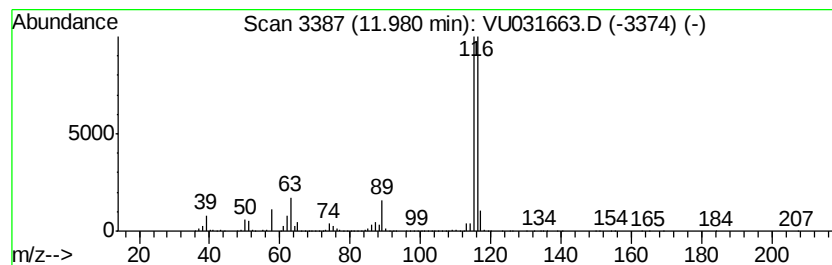
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Indene Concentration Rank 3

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 11.98 | 820.27 ug/L | 47923500 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Indene                  | 116 | C9H8    | 000095-13-6 | 97   |
| 2    |    | Indene                  | 116 | C9H8    | 000095-13-6 | 96   |
| 3    |    | Indene                  | 116 | C9H8    | 000095-13-6 | 93   |
| 4    |    | Benzene, 1-propynyl-    | 116 | C9H8    | 000673-32-5 | 91   |
| 5    |    | 3-Methylphenylacetylene | 116 | C9H8    | 000766-82-5 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

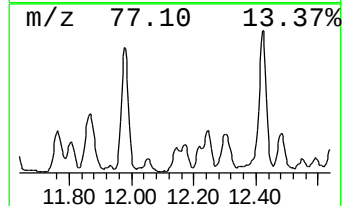
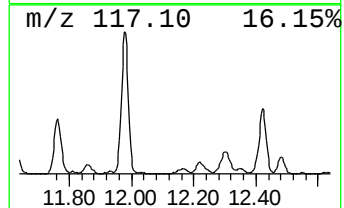
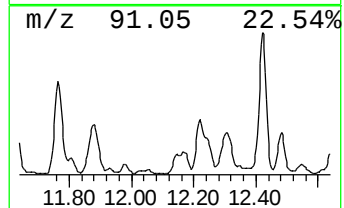
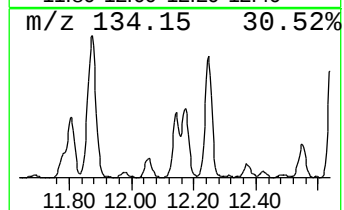
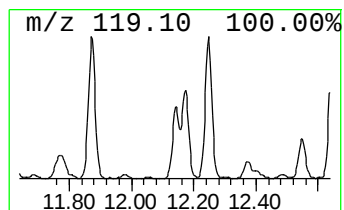
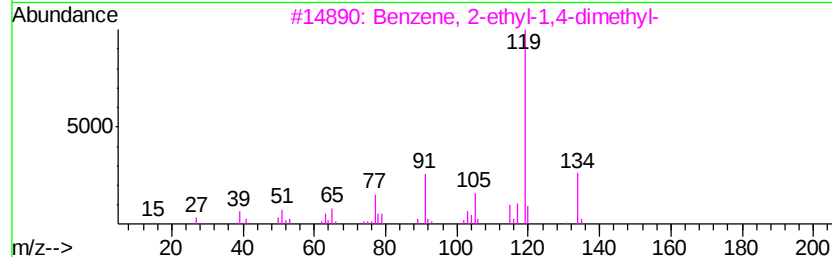
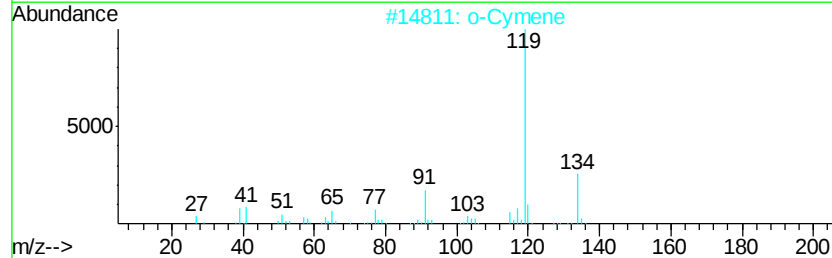
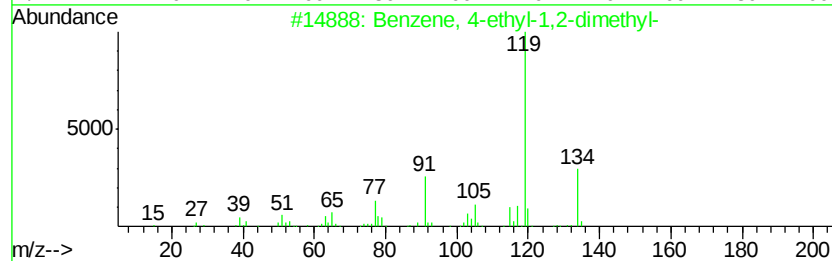
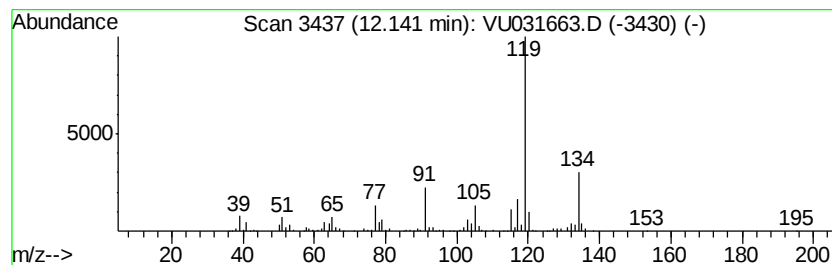
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 50

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.14 | 5.60 ug/L | 327161 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 93   |
| 2    |    | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 93   |
| 3    |    | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 91   |
| 4    |    | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 90   |
| 5    |    | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 87   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

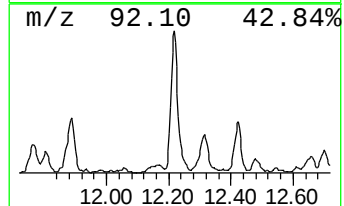
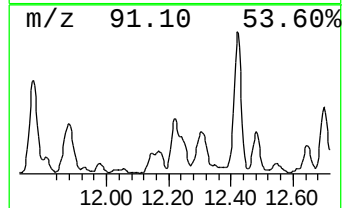
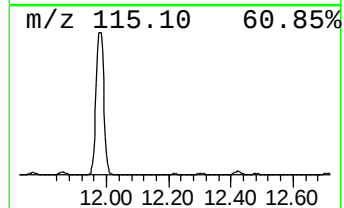
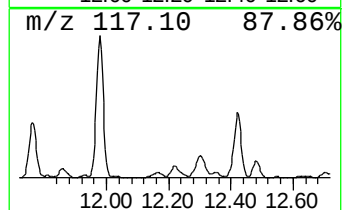
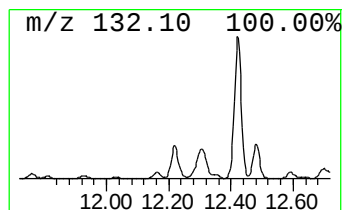
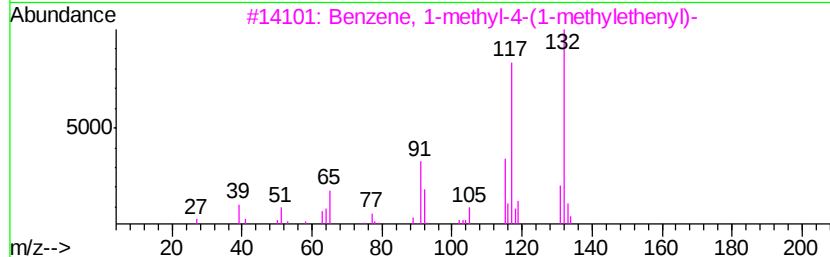
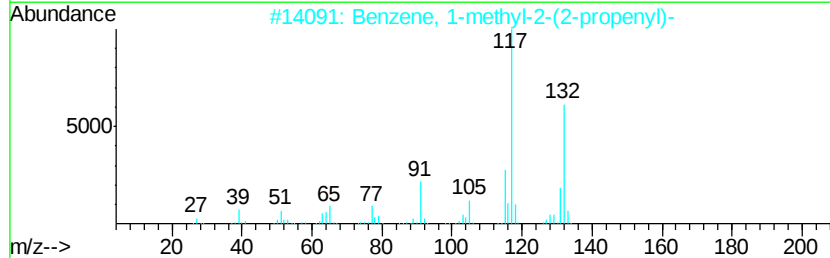
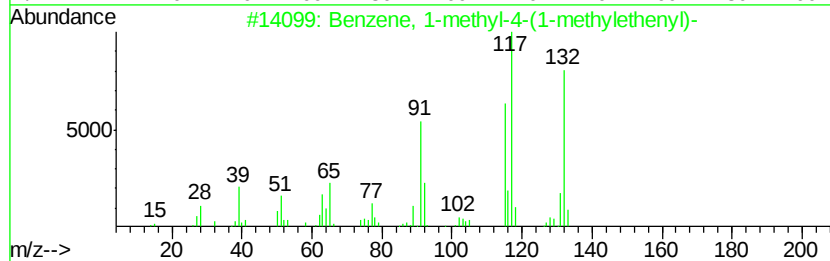
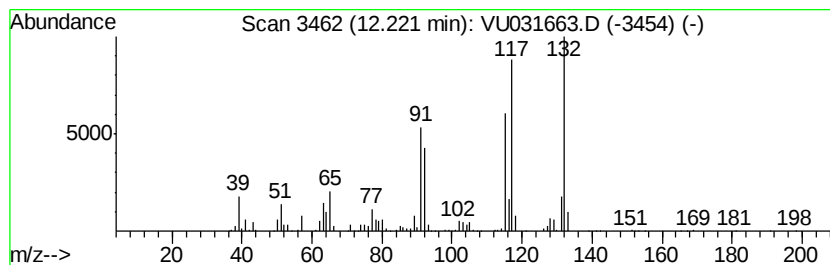
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Benzene, 1-methyl-4-(1-meth... Concentration Rank 41

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 12.22 | 12.48 ug/L | 728979 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Benzene, 1-methyl-4-(1-methyleth... | 132 | C10H12  | 001195-32-0  | 95   |
| 2    |    |   | Benzene, 1-methyl-2-(2-propenyl)-   | 132 | C10H12  | 001587-04-8  | 89   |
| 3    |    |   | Benzene, 1-methyl-4-(1-methyleth... | 132 | C10H12  | 001195-32-0  | 81   |
| 4    |    |   | Benzene, 1-methyl-4-(1-methyleth... | 132 | C10H12  | 001195-32-0  | 81   |
| 5    |    |   | 5,8-Dimethylenebicyclo[2.2.2]oct... | 132 | C10H12  | 1000210-64-8 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

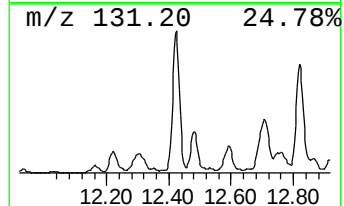
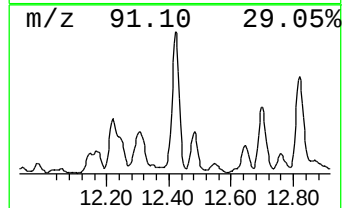
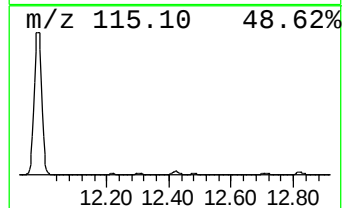
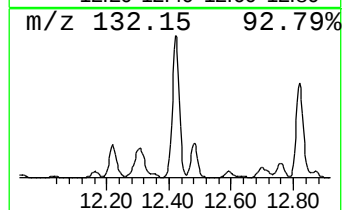
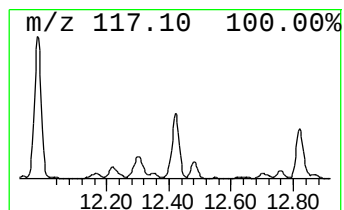
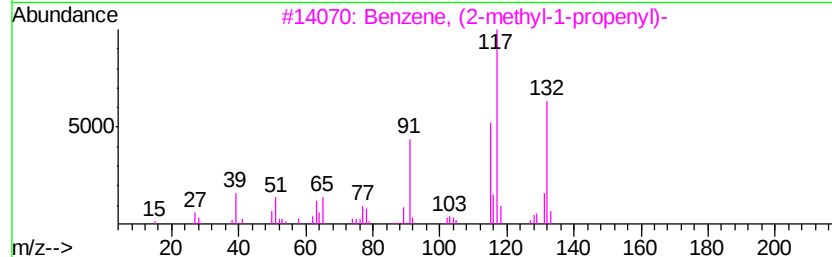
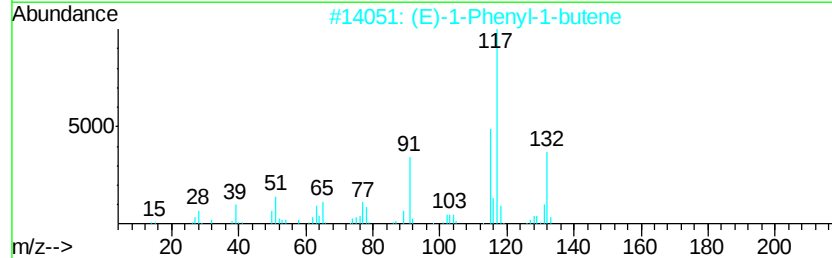
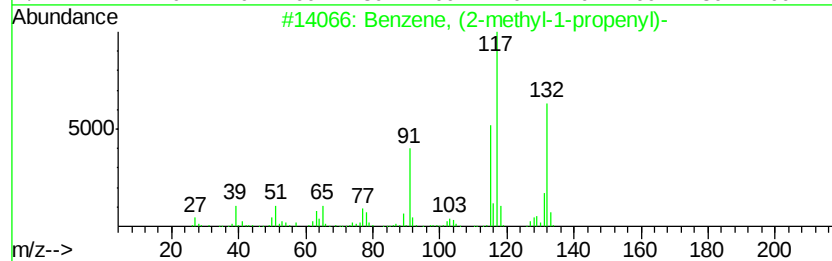
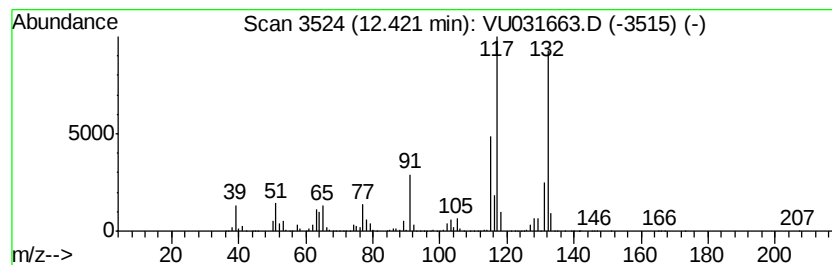
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Benzene, (2-methyl-1-propen... Concentration Rank 16

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.42 | 53.65 ug/L | 3134360 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (2-methyl-1-propenyl)-  | 132 | C10H12  | 000768-49-0 | 95   |
| 2    |    | (E)-1-Phenyl-1-butene            | 132 | C10H12  | 001005-64-7 | 94   |
| 3    |    | Benzene, (2-methyl-1-propenyl)-  | 132 | C10H12  | 000768-49-0 | 94   |
| 4    |    | o-Isopropenyltoluene             | 132 | C10H12  | 007399-49-7 | 94   |
| 5    |    | Benzene, 2-ethenyl-1,3-dimethyl- | 132 | C10H12  | 002039-90-9 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

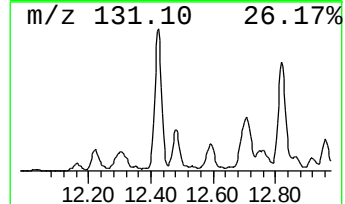
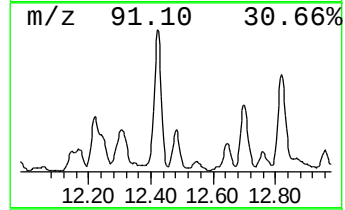
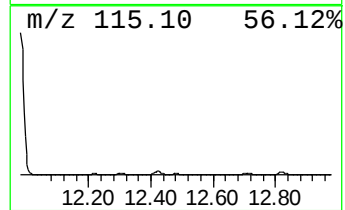
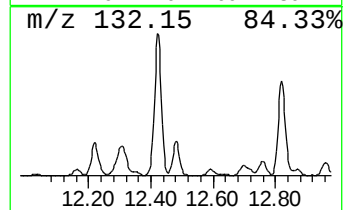
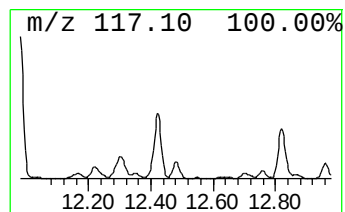
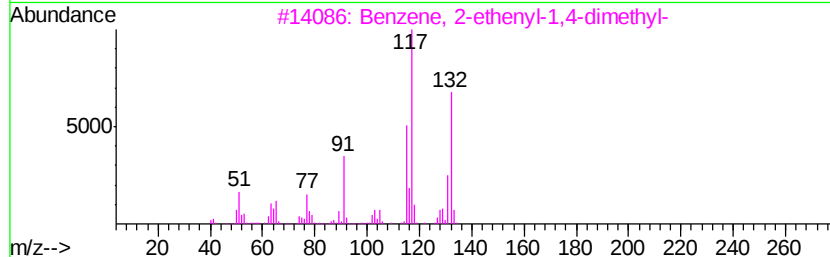
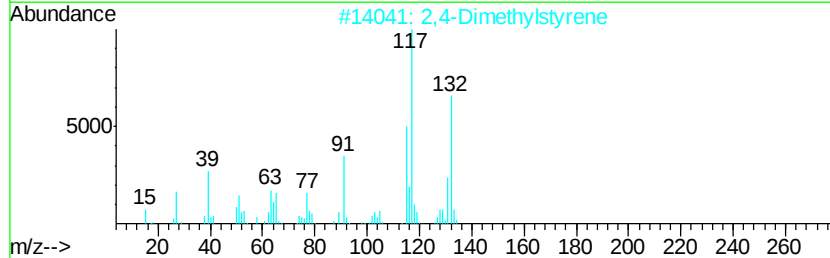
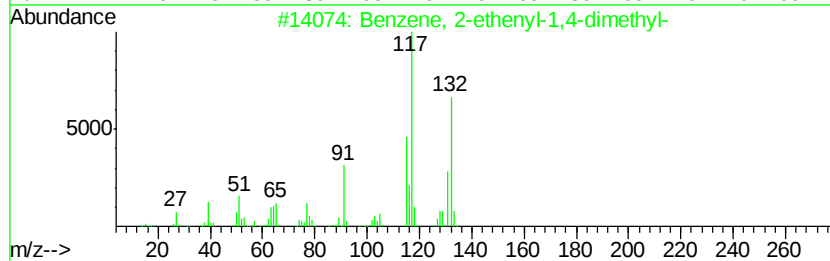
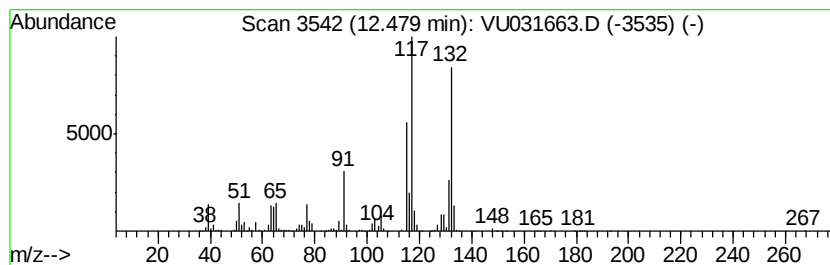
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 37

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 12.48 | 14.20 ug/L | 829554 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethenyl-1,4-dimethyl-  | 132 | C10H12  | 002039-89-6 | 98   |
| 2    |    | 2,4-Dimethylstyrene               | 132 | C10H12  | 002234-20-0 | 97   |
| 3    |    | Benzene, 2-ethenyl-1,4-dimethyl-  | 132 | C10H12  | 002039-89-6 | 96   |
| 4    |    | Benzene, 2-ethenyl-1,3-dimethyl-  | 132 | C10H12  | 002039-90-9 | 95   |
| 5    |    | Benzene, 1-methyl-2-(2-propenyl)- | 132 | C10H12  | 001587-04-8 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

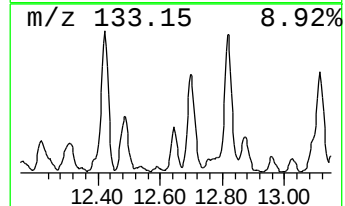
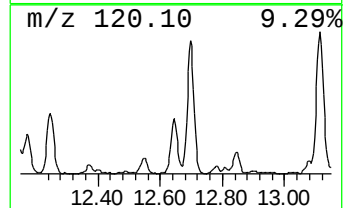
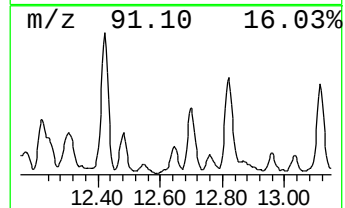
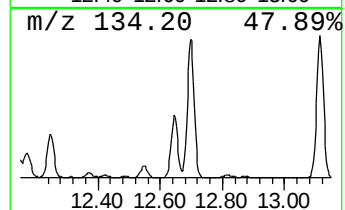
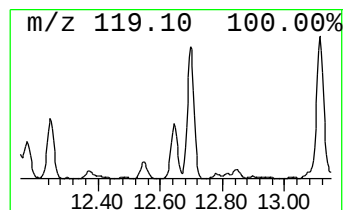
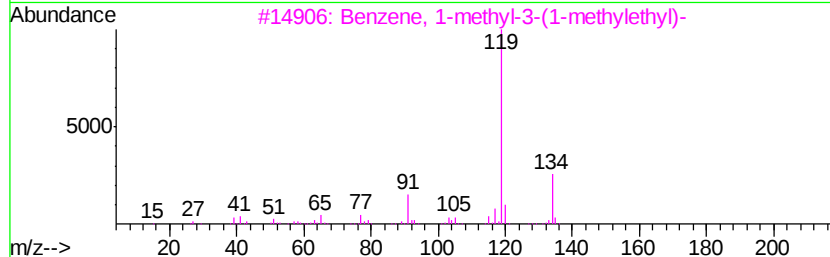
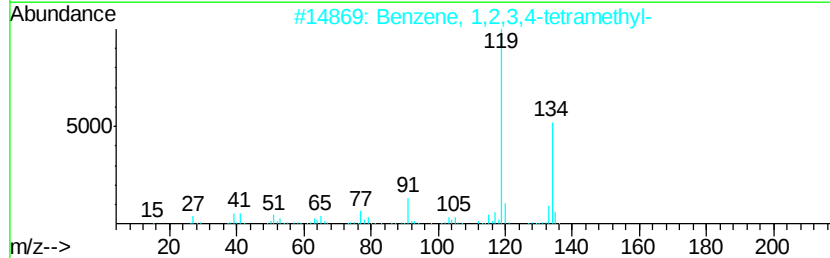
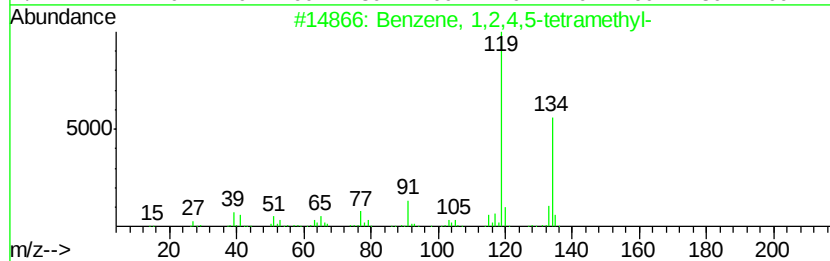
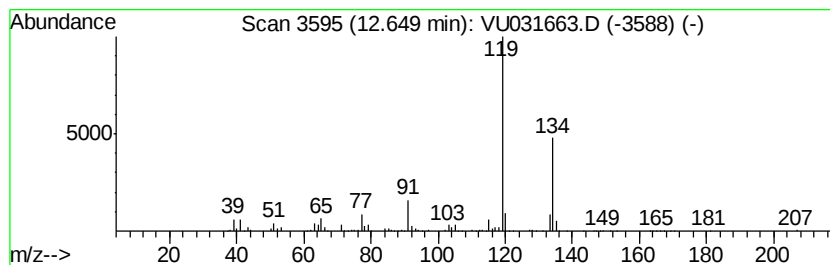
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 46

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.65 | 8.94 ug/L | 522070 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl-        | 134 | C10H14  | 000095-93-2 | 94   |
| 2    |    |   | Benzene, 1,2,3,4-tetramethyl-        | 134 | C10H14  | 000488-23-3 | 94   |
| 3    |    |   | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 94   |
| 4    |    |   | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 93   |
| 5    |    |   | Benzene, 1-ethyl-3,5-dimethyl-       | 134 | C10H14  | 000934-74-7 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

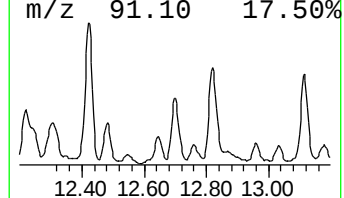
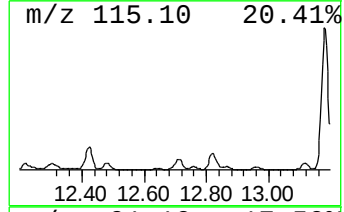
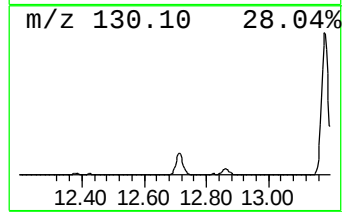
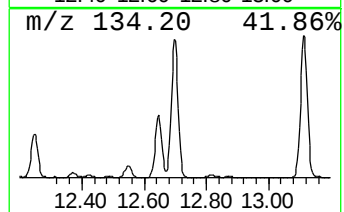
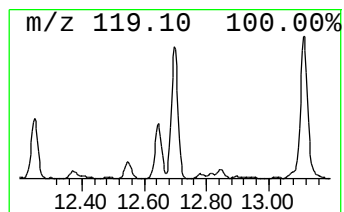
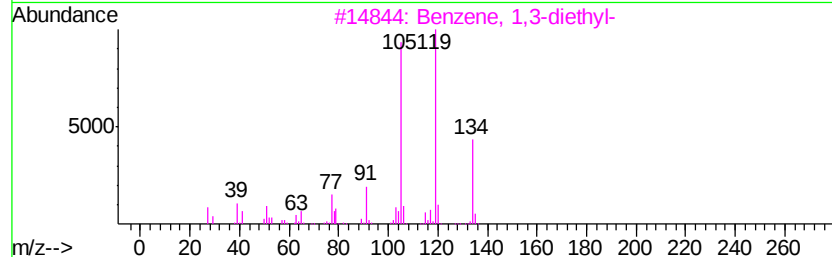
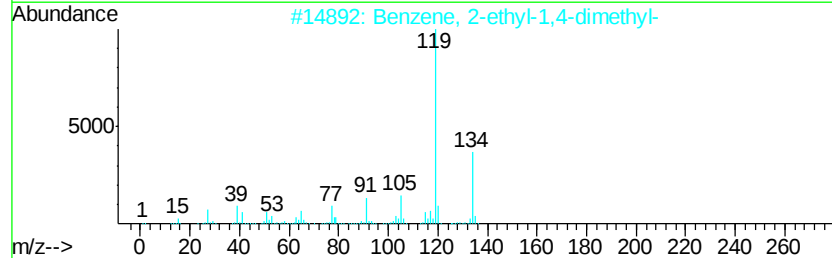
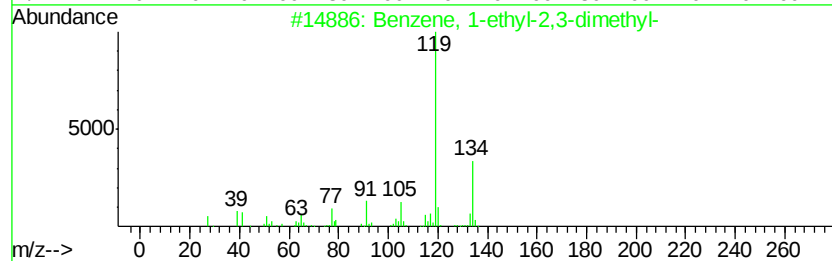
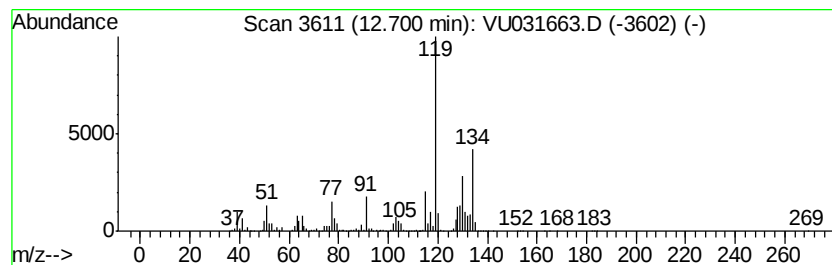
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 17

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.70 | 48.21 ug/L | 2816390 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 55   |
| 2    |    | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 55   |
| 3    |    | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 55   |
| 4    |    | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 55   |
| 5    |    | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 55   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

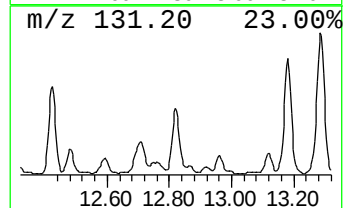
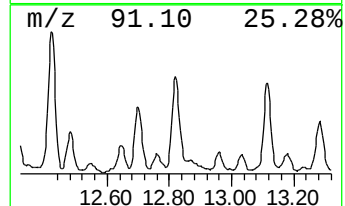
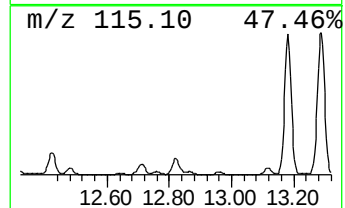
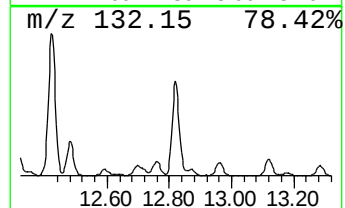
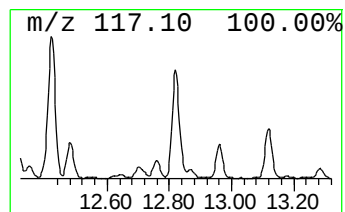
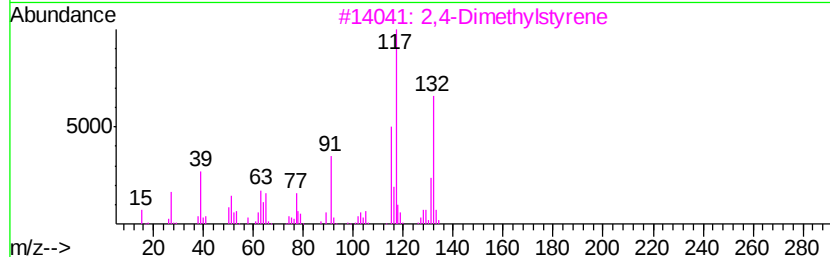
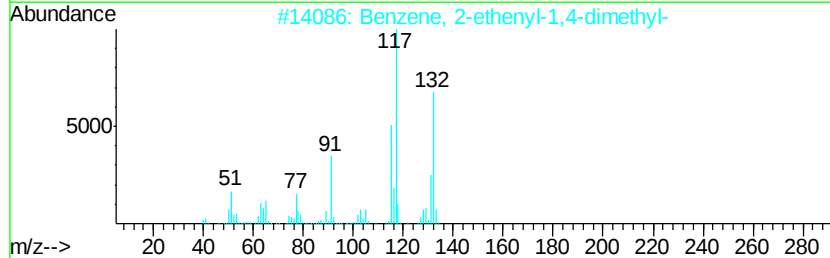
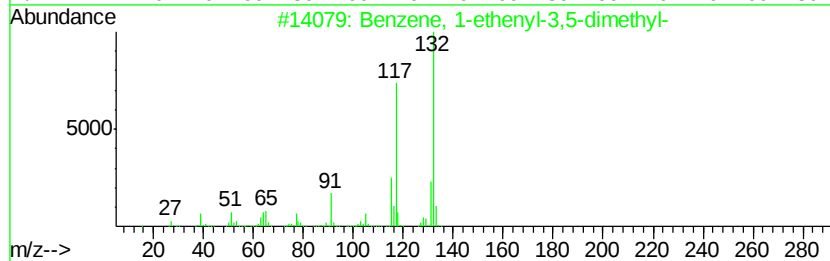
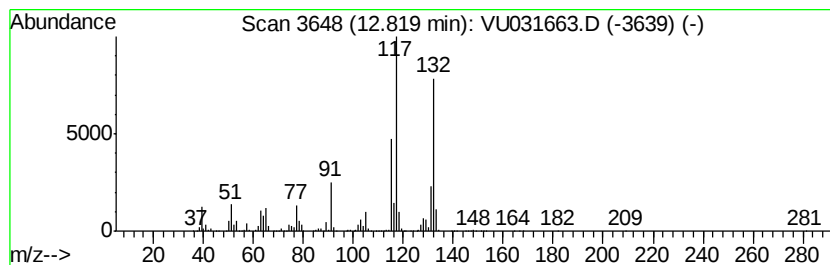
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 Benzene, 1-ethenyl-3,5-dime... Concentration Rank 22

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.82 | 41.66 ug/L | 2433960 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethenyl-3,5-dimethyl-  | 132 | C10H12  | 005379-20-4 | 97   |
| 2    |    | Benzene, 2-ethenyl-1,4-dimethyl-  | 132 | C10H12  | 002039-89-6 | 96   |
| 3    |    | 2,4-Dimethylstyrene               | 132 | C10H12  | 002234-20-0 | 96   |
| 4    |    | Benzene, 4-ethenyl-1,2-dimethyl-  | 132 | C10H12  | 027831-13-6 | 95   |
| 5    |    | Benzene, 1-methyl-4-(2-propenyl)- | 132 | C10H12  | 003333-13-9 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

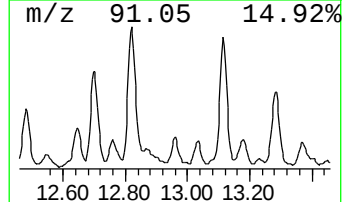
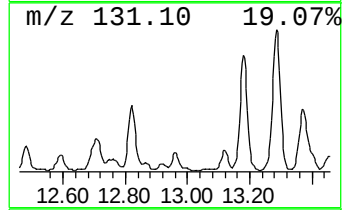
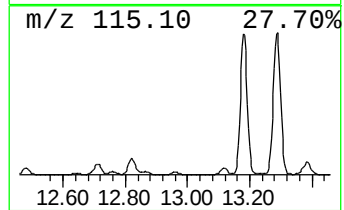
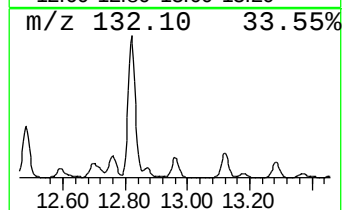
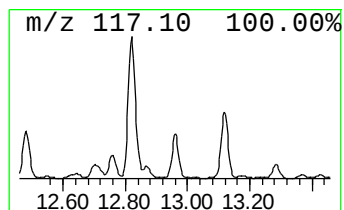
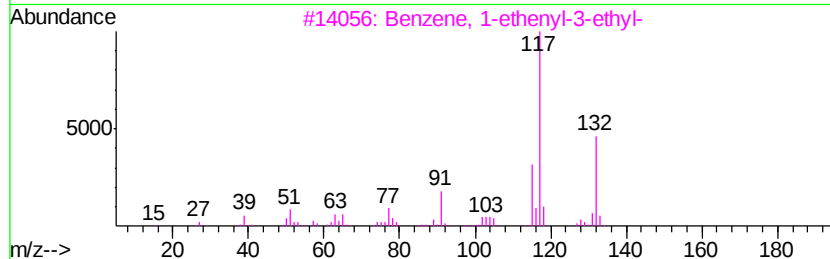
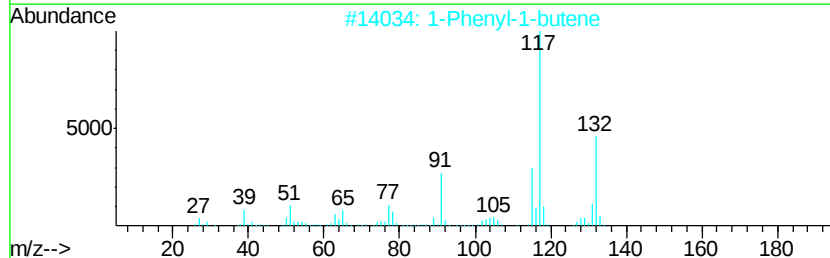
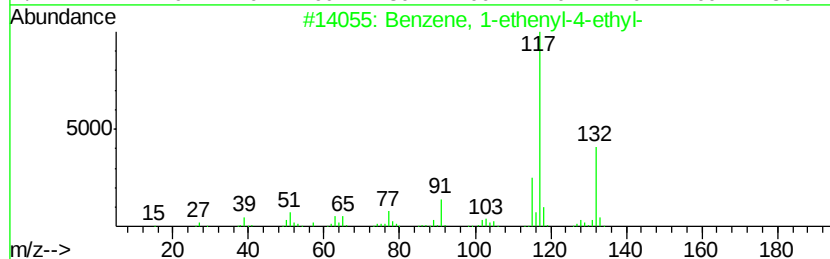
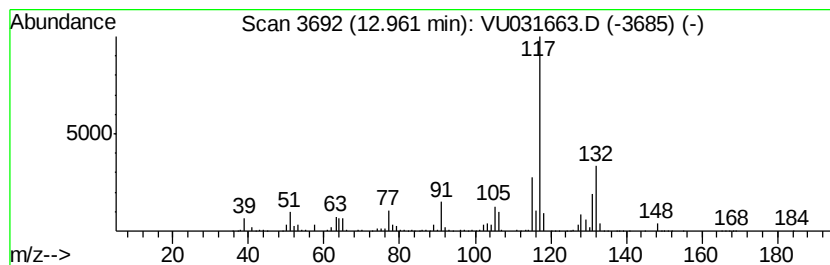
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 45

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.96 | 9.04 ug/L | 528213 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethenyl-4-ethyl-      | 132 | C10H12  | 003454-07-7 | 94   |
| 2    |    |   | 1-Phenyl-1-butene                | 132 | C10H12  | 000824-90-8 | 94   |
| 3    |    |   | Benzene, 1-ethenyl-3-ethyl-      | 132 | C10H12  | 007525-62-4 | 93   |
| 4    |    |   | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12  | 002039-89-6 | 92   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-5-methyl- | 132 | C10H12  | 000874-35-1 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

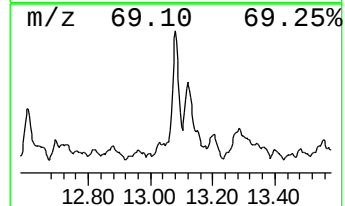
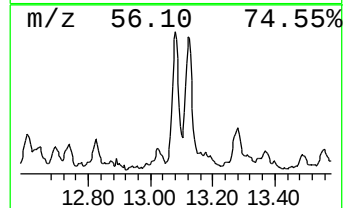
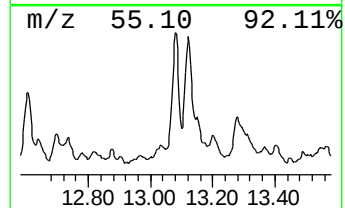
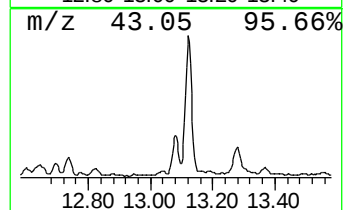
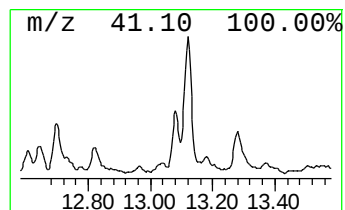
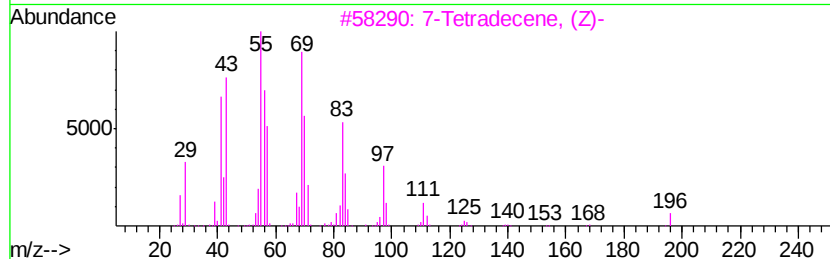
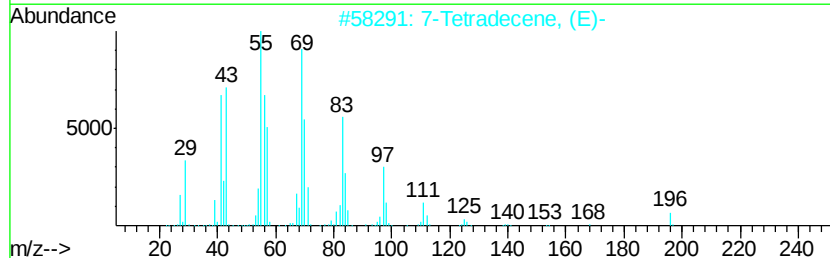
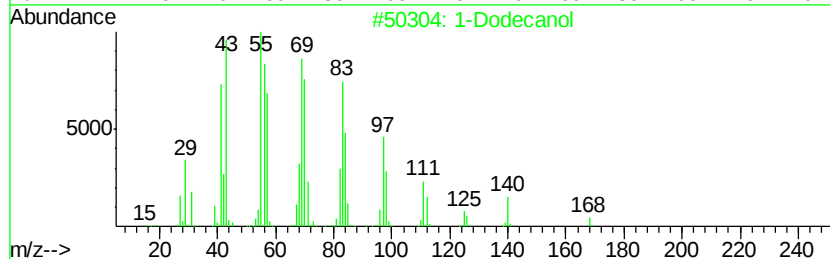
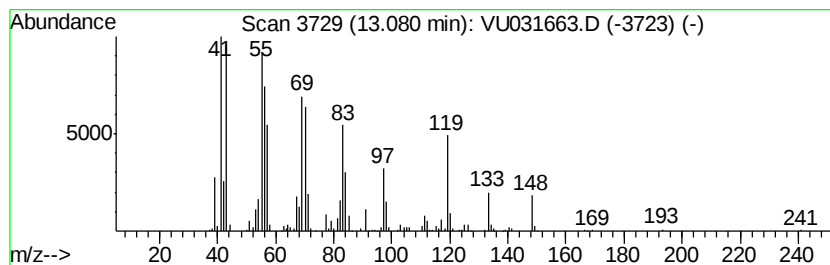
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 23 1-Dodecanol Concentration Rank 51

| R.T.      | EstConc             | Area   | Relative to ISTD       | R.T.        |      |
|-----------|---------------------|--------|------------------------|-------------|------|
| 13.08     | 4.90 ug/L           | 286265 | 1,4-Dichlorobenzene-d4 | 11.48       |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm                | CAS#        | Qual |
| 1         | 1-Dodecanol         | 186    | C12H26O                | 000112-53-8 | 68   |
| 2         | 7-Tetradecene, (E)- | 196    | C14H28                 | 041446-63-3 | 68   |
| 3         | 7-Tetradecene, (Z)- | 196    | C14H28                 | 041446-60-0 | 68   |
| 4         | 2-Dodecene, (Z)-    | 168    | C12H24                 | 007206-26-0 | 68   |
| 5         | 5-Tetradecene, (E)- | 196    | C14H28                 | 041446-66-6 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

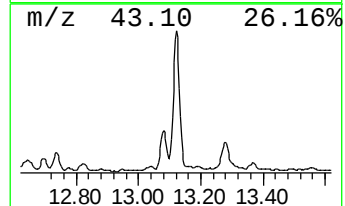
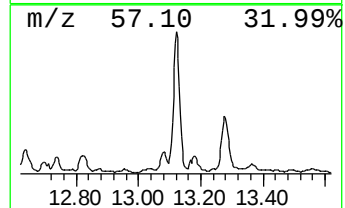
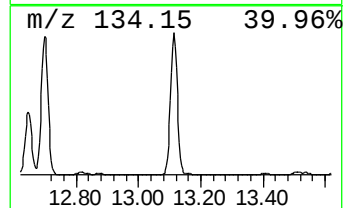
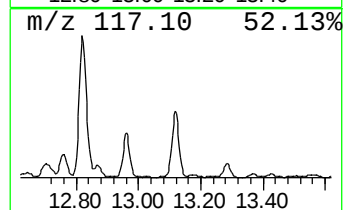
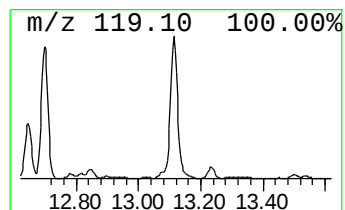
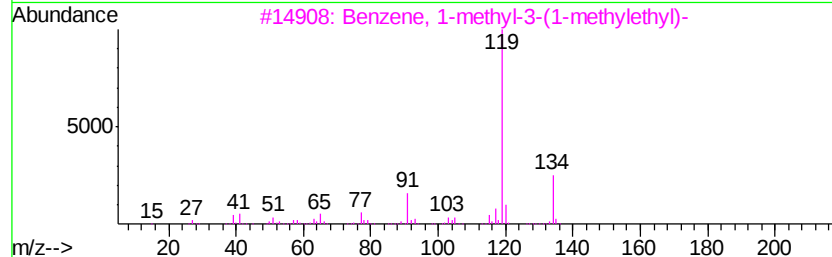
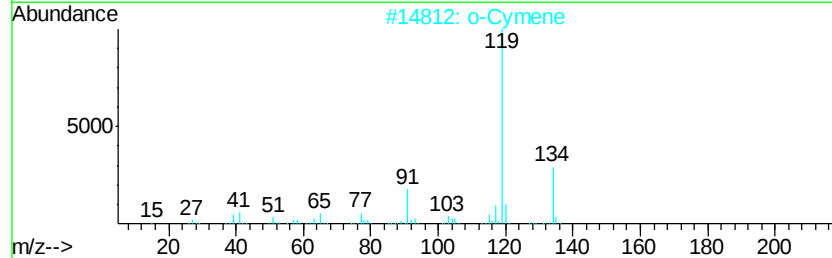
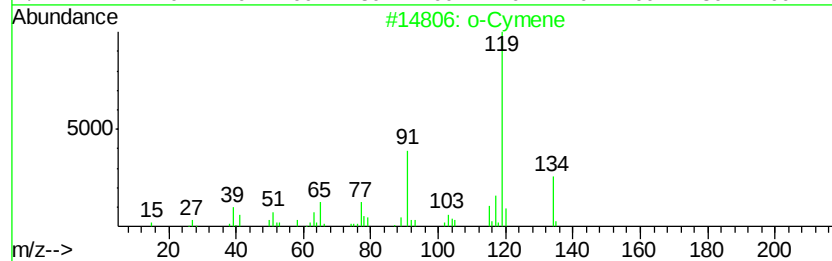
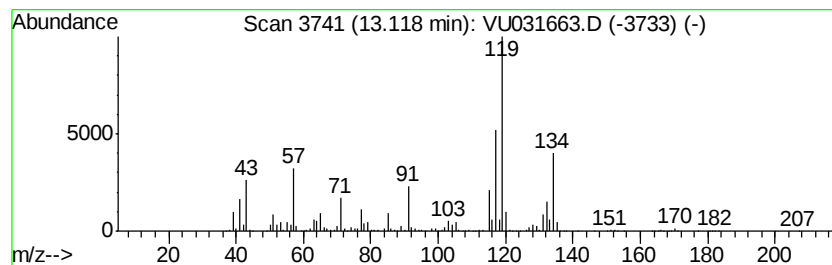
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 o-Cymene Concentration Rank 18

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 13.12 | 44.56 ug/L | 2603240 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 70   |
| 2    |    | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 64   |
| 3    |    | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 64   |
| 4    |    | p-Cymene                             | 134 | C10H14  | 000099-87-6 | 64   |
| 5    |    | Benzene, 1,2,4,5-tetramethyl-        | 134 | C10H14  | 000095-93-2 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

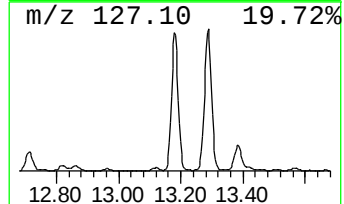
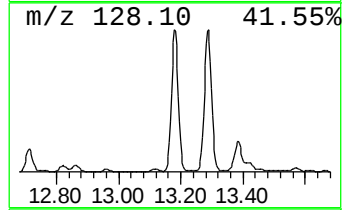
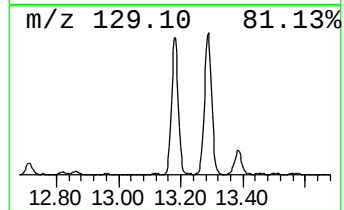
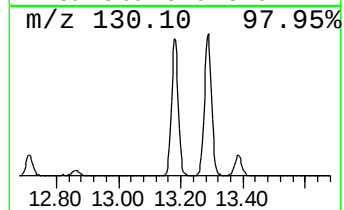
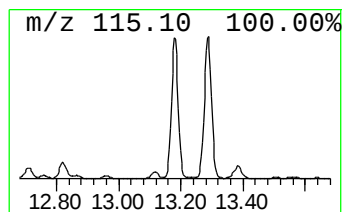
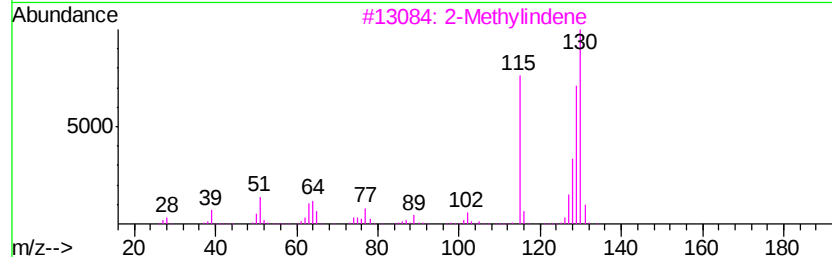
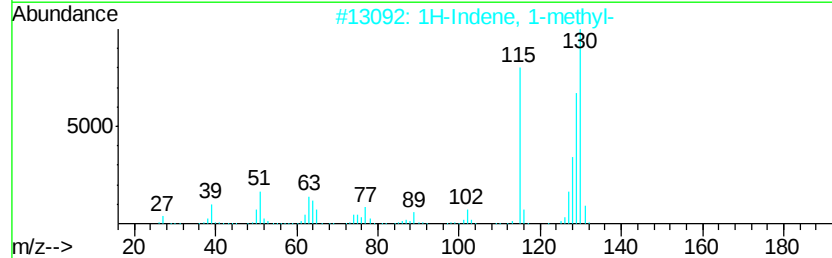
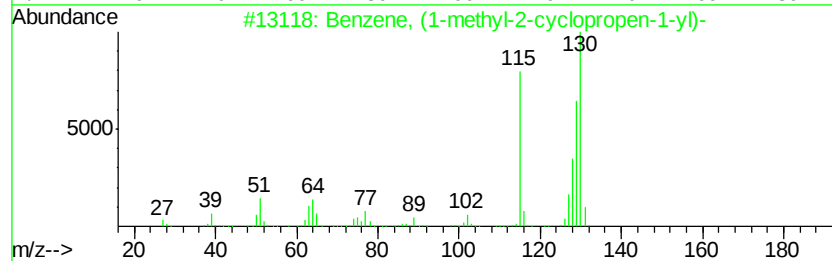
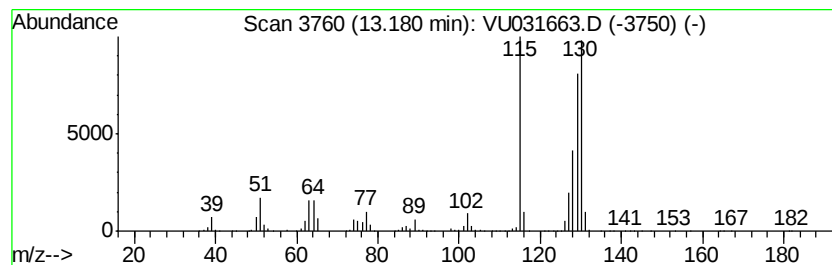
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 Benzene, (1-methyl-2-cyclop... Concentration Rank 10

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 13.18 | 167.02 ug/L | 9757630 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (1-methyl-2-cyclopropen... | 130 | C10H10  | 065051-83-4 | 97   |
| 2    |    | 1H-Indene, 1-methyl-                | 130 | C10H10  | 000767-59-9 | 96   |
| 3    |    | 2-Methylindene                      | 130 | C10H10  | 002177-47-1 | 96   |
| 4    |    | Cycloprop[alindene, 1,1a,6,6a-te... | 130 | C10H10  | 015677-15-3 | 95   |
| 5    |    | 1H-Indene, 3-methyl-                | 130 | C10H10  | 000767-60-2 | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

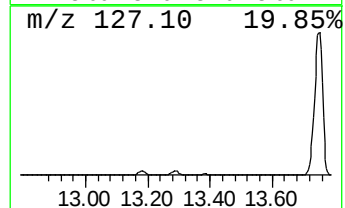
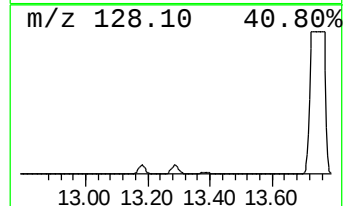
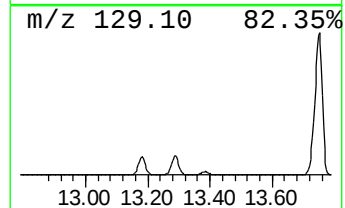
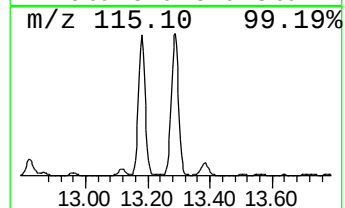
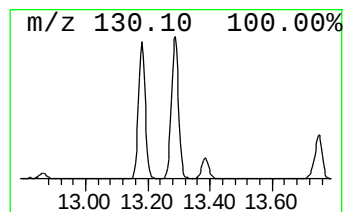
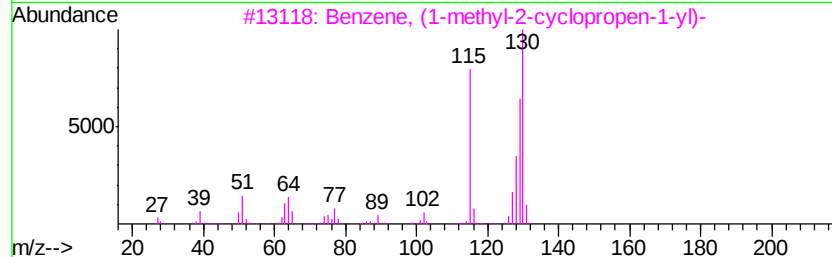
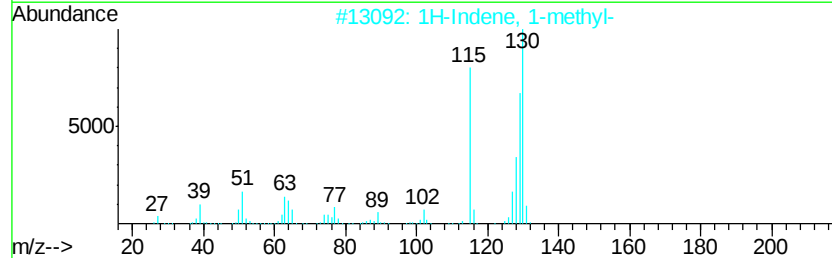
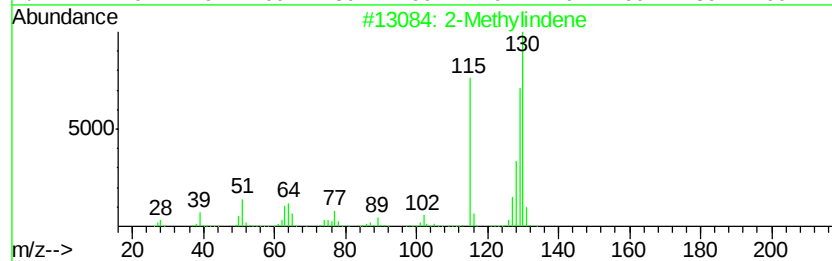
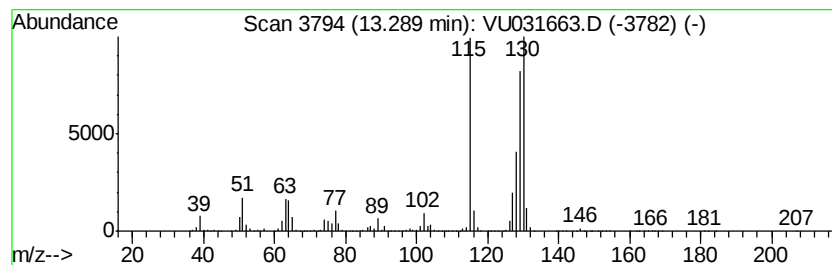
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 2-Methylindene Concentration Rank 8

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 13.29 | 193.98 ug/L | 11333100 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                              | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Methylindene                            | 130 | C10H10  | 002177-47-1 | 96   |
| 2    |    | 1H-Indene, 1-methyl-                      | 130 | C10H10  | 000767-59-9 | 96   |
| 3    |    | Benzene, (1-methyl-2-cyclopropen-1-yl)-   | 130 | C10H10  | 065051-83-4 | 96   |
| 4    |    | 1H-Indene, 1-methyl-                      | 130 | C10H10  | 000767-59-9 | 95   |
| 5    |    | Cycloprop[a]indene, 1,1a,6,6a-tetrahydro- | 130 | C10H10  | 015677-15-3 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

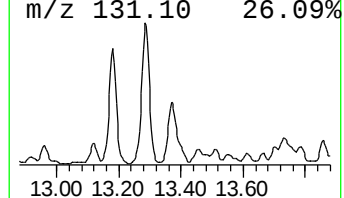
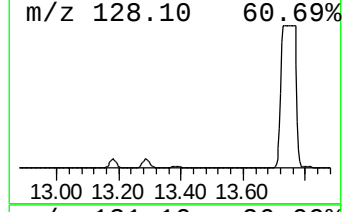
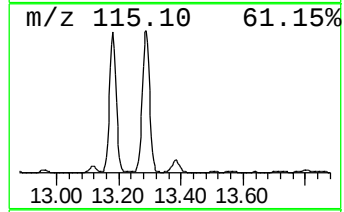
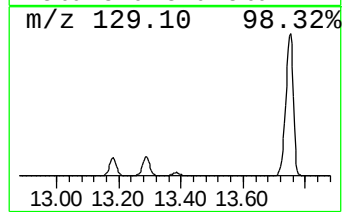
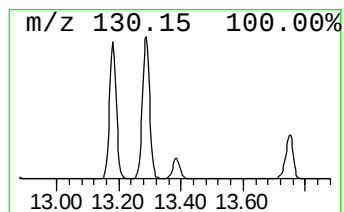
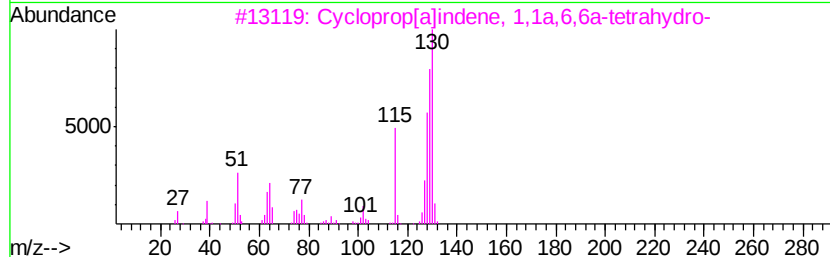
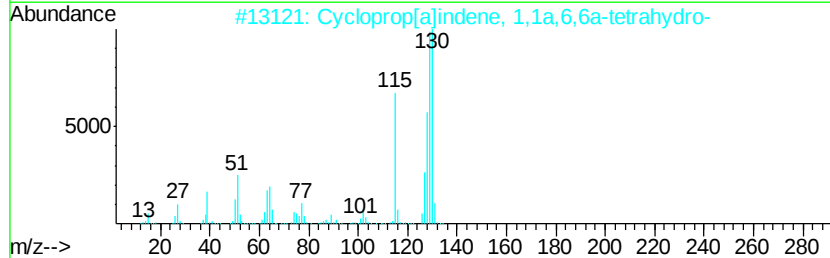
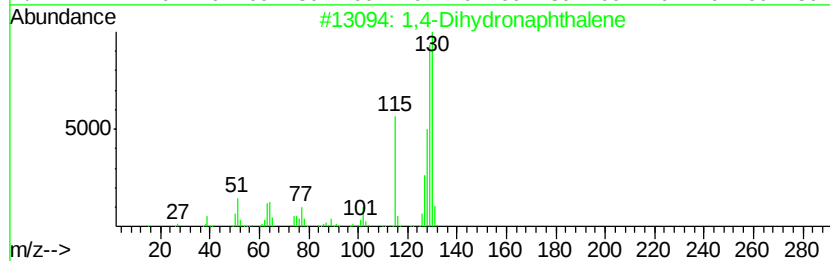
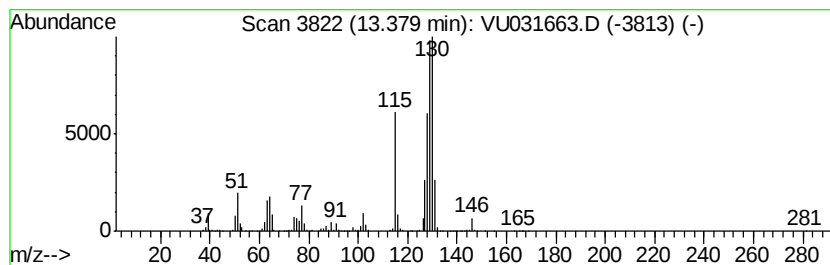
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 27 1,4-Dihydronaphthalene Concentration Rank 27

| R.T.    | EstConc                             | Area         | Relative to ISTD       | R.T.      |
|---------|-------------------------------------|--------------|------------------------|-----------|
| 13.38   | 31.47 ug/L                          | 1838630      | 1,4-Dichlorobenzene-d4 | 11.48     |
| Hit# of | 5                                   | Tentative ID | MW MolForm             | CAS# Qual |
| 1       | 1,4-Dihydronaphthalene              | 130 C10H10   | 000612-17-9            | 96        |
| 2       | Cycloprop[alindene, 1,1a,6,6a-te... | 130 C10H10   | 015677-15-3            | 96        |
| 3       | Cycloprop[alindene, 1,1a,6,6a-te... | 130 C10H10   | 015677-15-3            | 94        |
| 4       | Cycloprop[alindene, 1,1a,6,6a-te... | 130 C10H10   | 015677-15-3            | 93        |
| 5       | Naphthalene, 1,2-dihydro-           | 130 C10H10   | 000447-53-0            | 93        |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

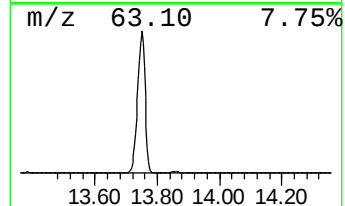
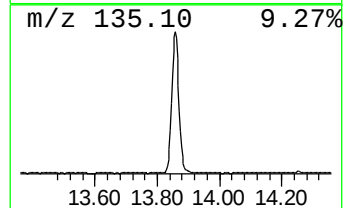
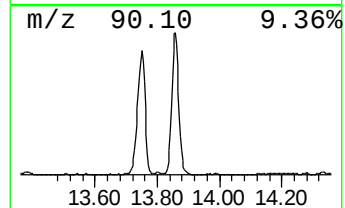
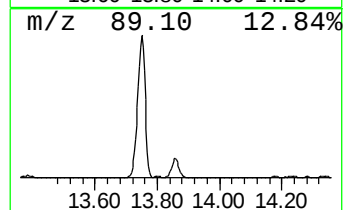
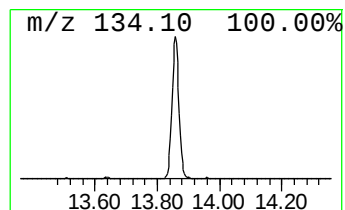
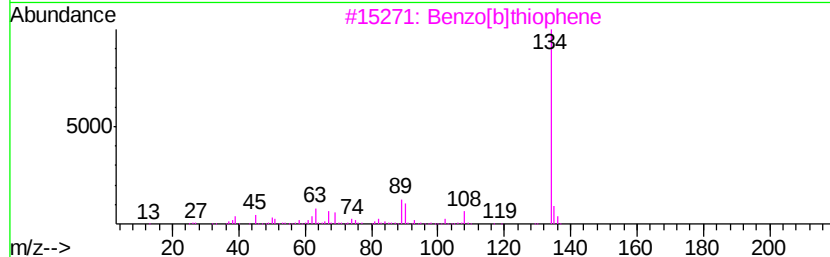
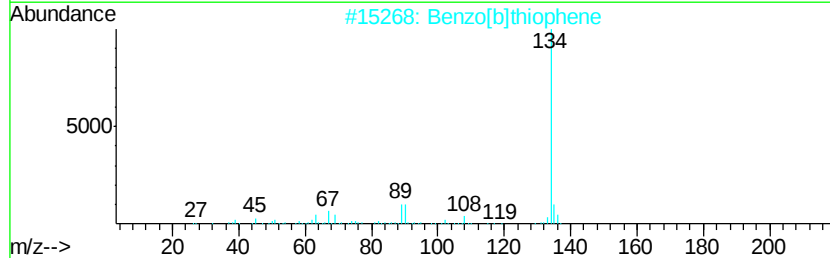
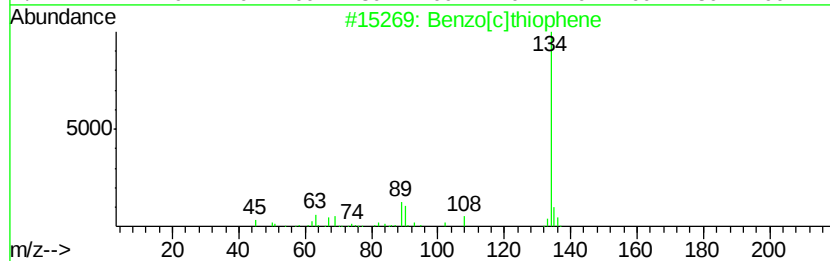
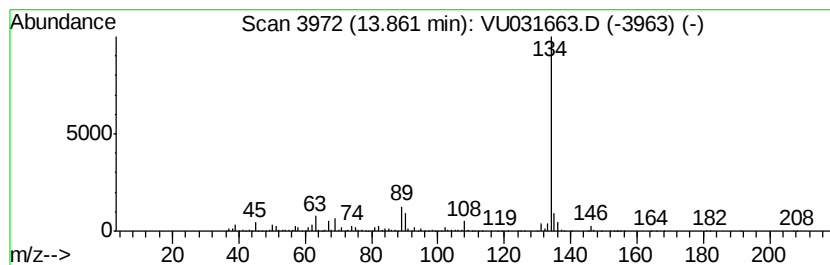
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 29 Benzo[c]thiophene Concentration Rank 19

| R.T.    | EstConc    | Area                   | Relative to ISTD       | R.T.           |
|---------|------------|------------------------|------------------------|----------------|
| 13.86   | 44.36 ug/L | 2591740                | 1,4-Dichlorobenzene-d4 | 11.48          |
| Hit# of | 5          | Tentative ID           | MW MolForm             | CAS# Qual      |
| 1       |            | Benzo[c]thiophene      | 134 C8H6S              | 000270-82-6 97 |
| 2       |            | Benzo[b]thiophene      | 134 C8H6S              | 000095-15-8 95 |
| 3       |            | Benzo[b]thiophene      | 134 C8H6S              | 000095-15-8 94 |
| 4       |            | Cyclopenta[c]thiapyran | 134 C8H6S              | 000270-63-3 94 |
| 5       |            | Benzo[b]thiophene      | 134 C8H6S              | 000095-15-8 91 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

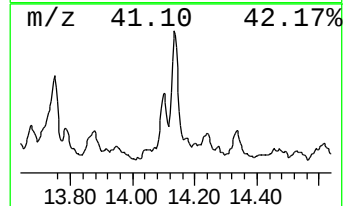
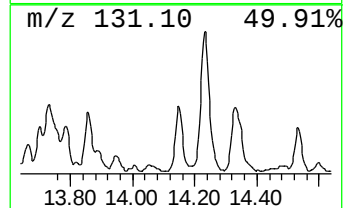
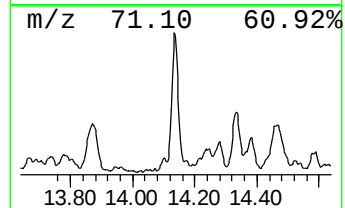
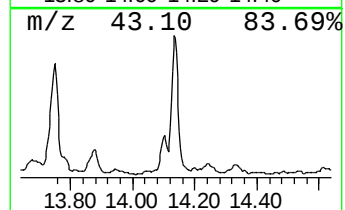
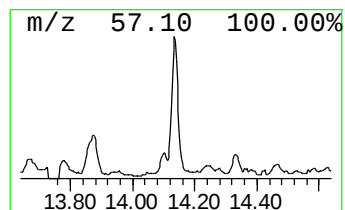
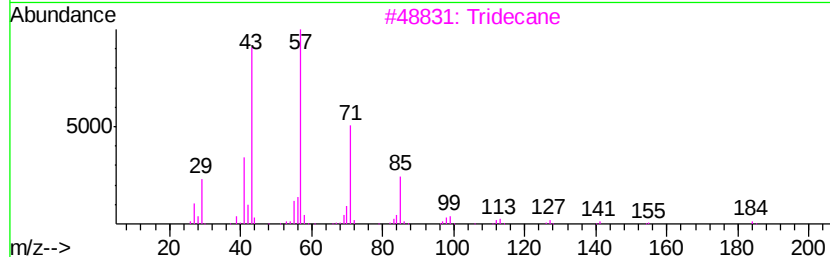
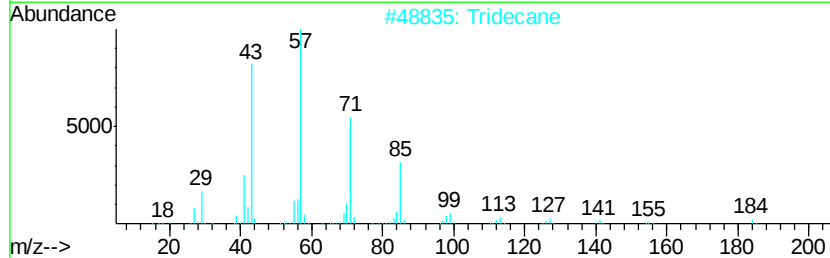
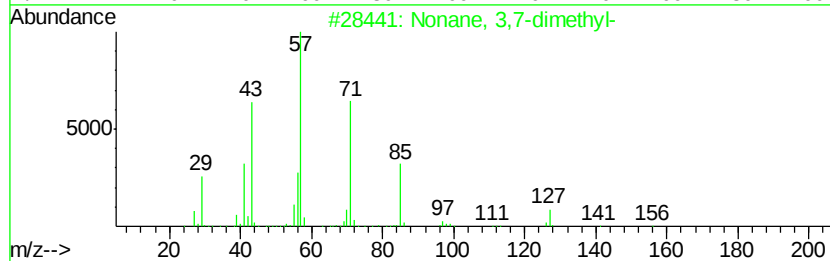
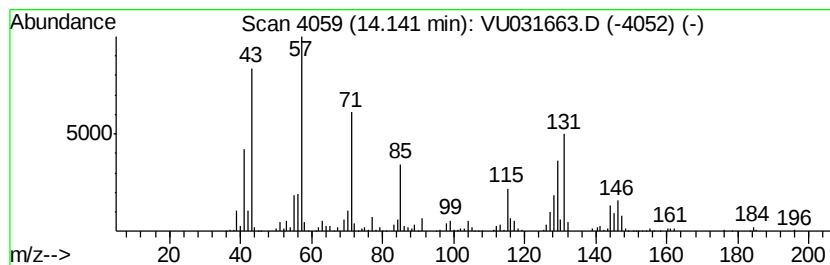
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 47

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 14.14 | 7.83 ug/L | 457383 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 3,7-dimethyl- | 156 | C11H24  | 017302-32-8 | 55   |
| 2    |    |   | Tridecane             | 184 | C13H28  | 000629-50-5 | 50   |
| 3    |    |   | Tridecane             | 184 | C13H28  | 000629-50-5 | 50   |
| 4    |    |   | Decane, 3-methyl-     | 156 | C11H24  | 013151-34-3 | 46   |
| 5    |    |   | 4-Octanone            | 128 | C8H16O  | 000589-63-9 | 43   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

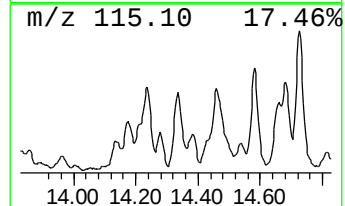
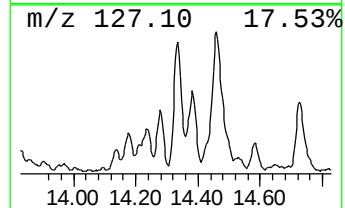
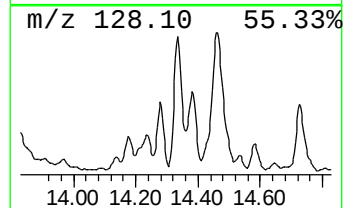
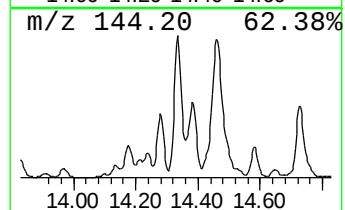
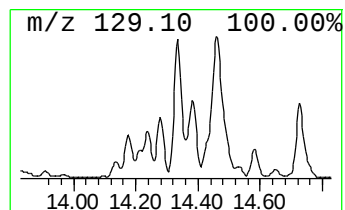
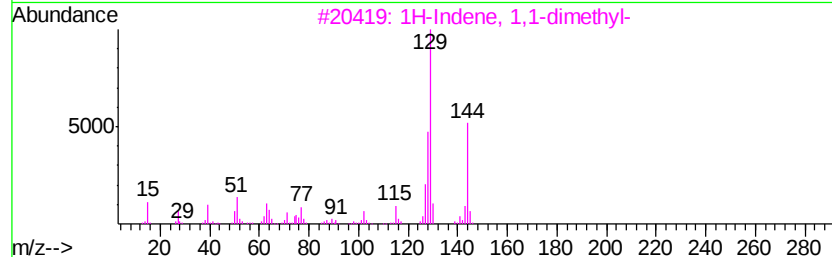
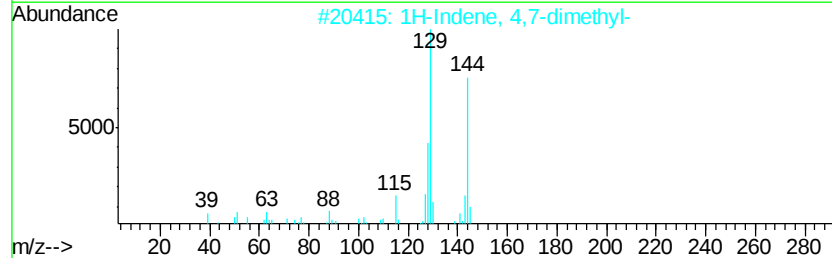
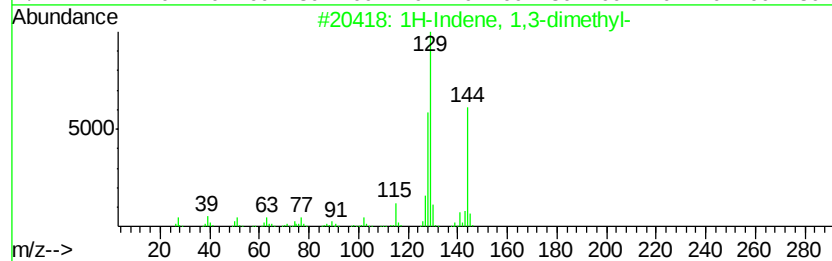
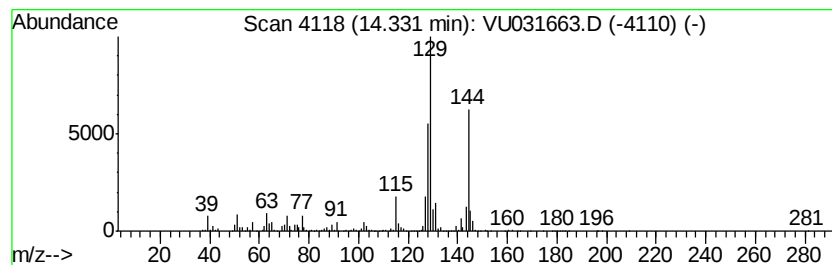
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 31 1H-Indene, 1,3-dimethyl- Concentration Rank 30

| R.T.    | EstConc                  | Area         | Relative to ISTD       | R.T.      |
|---------|--------------------------|--------------|------------------------|-----------|
| 14.33   | 26.92 ug/L               | 1572590      | 1,4-Dichlorobenzene-d4 | 11.48     |
| Hit# of | 5                        | Tentative ID | MW MolForm             | CAS# Qual |
| 1       | 1H-Indene, 1,3-dimethyl- | 144 C11H12   | 002177-48-2            | 94        |
| 2       | 1H-Indene, 4,7-dimethyl- | 144 C11H12   | 006974-97-6            | 93        |
| 3       | 1H-Indene, 1,1-dimethyl- | 144 C11H12   | 018636-55-0            | 93        |
| 4       | 1H-Indene, 2,3-dimethyl- | 144 C11H12   | 004773-82-4            | 91        |
| 5       | 2-Ethyl-1-H-indene       | 144 C11H12   | 017059-50-6            | 87        |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

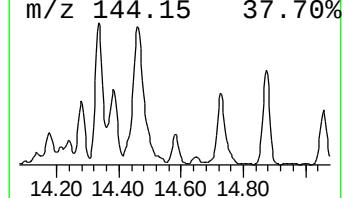
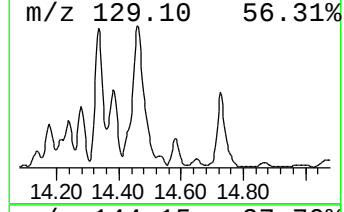
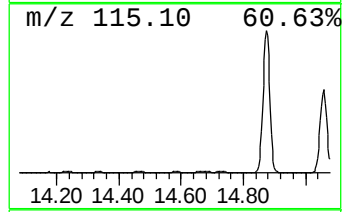
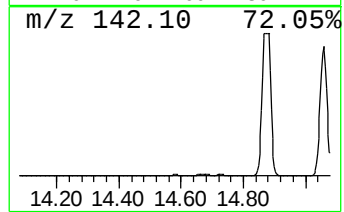
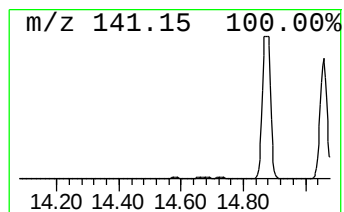
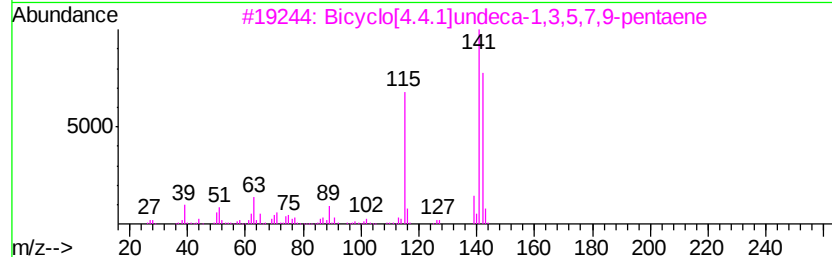
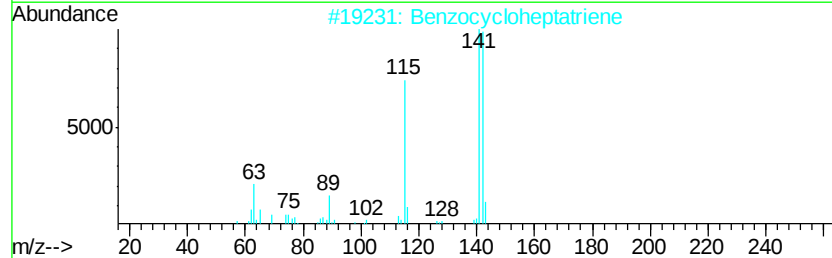
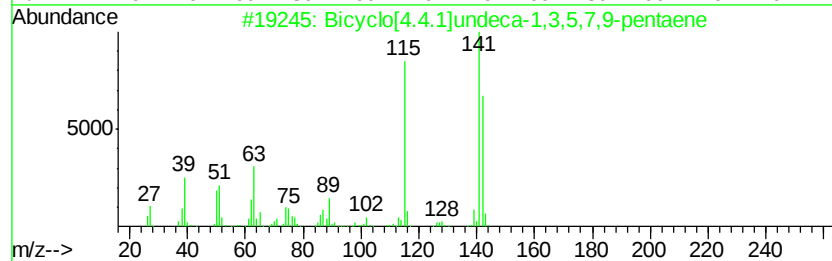
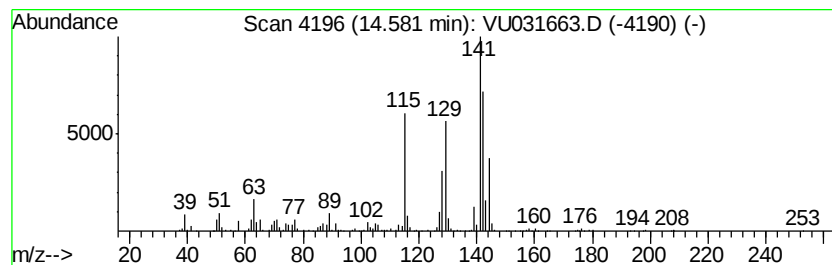
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 33 Bicyclo[4.4.1]undeca-1,3,5,... Concentration Rank 43

| R.T.    | EstConc    | Area                                | Relative to ISTD       | R.T.           |
|---------|------------|-------------------------------------|------------------------|----------------|
| 14.58   | 10.53 ug/L | 615416                              | 1,4-Dichlorobenzene-d4 | 11.48          |
| Hit# of | 5          | Tentative ID                        | MW MolForm             | CAS# Qual      |
| 1       |            | Bicyclo[4.4.1]undeca-1,3,5,7,9-p... | 142 C11H10             | 002443-46-1 91 |
| 2       |            | Benzocycloheptatriene               | 142 C11H10             | 000264-09-5 90 |
| 3       |            | Bicyclo[4.4.1]undeca-1,3,5,7,9-p... | 142 C11H10             | 002443-46-1 83 |
| 4       |            | 1H-Indene, 1-ethylidene-            | 142 C11H10             | 002471-83-2 64 |
| 5       |            | 1,4-Methanonaphthalene, 1,4-dihy... | 142 C11H10             | 004453-90-1 64 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

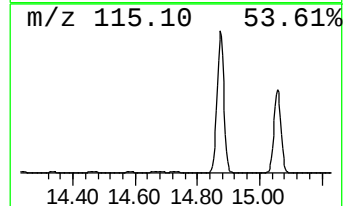
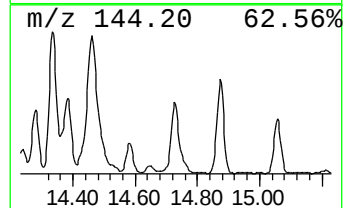
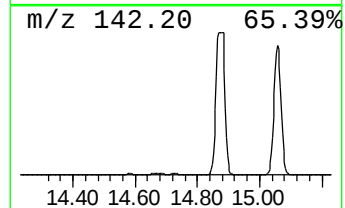
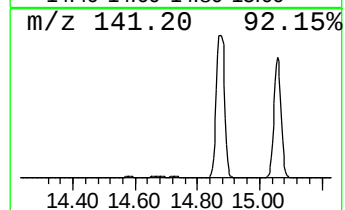
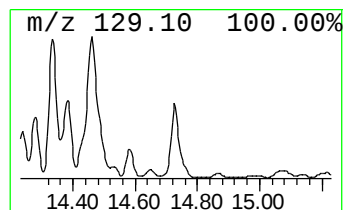
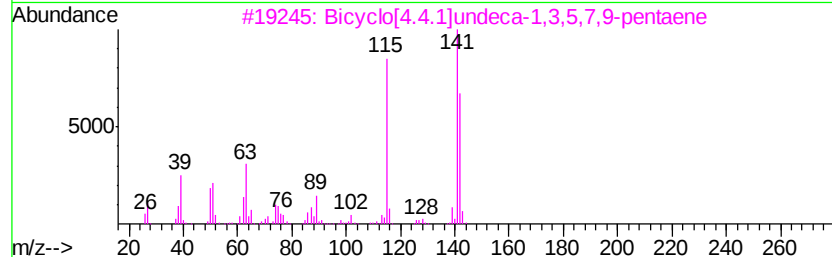
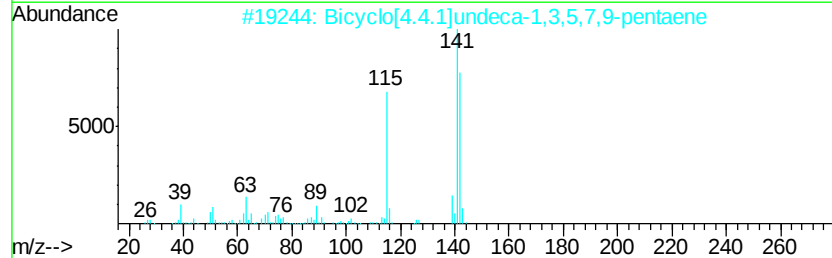
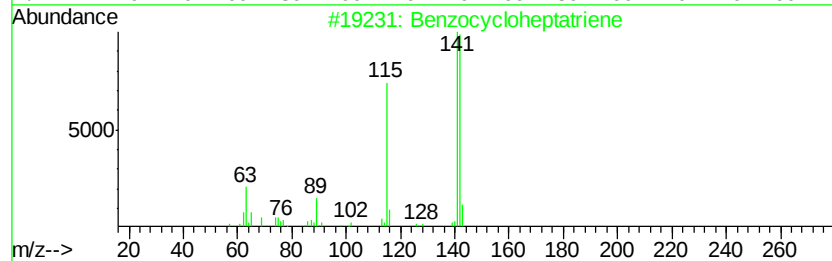
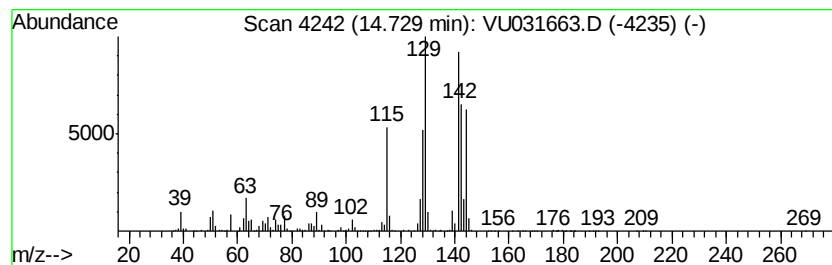
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 Benzocycloheptatriene Concentration Rank 34

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 14.73 | 22.57 ug/L | 1318900 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzocycloheptatriene               | 142 | C11H10  | 000264-09-5 | 94   |
| 2    |    | Bicyclo[4.4.1]undeca-1,3,5,7,9-p... | 142 | C11H10  | 002443-46-1 | 89   |
| 3    |    | Bicyclo[4.4.1]undeca-1,3,5,7,9-p... | 142 | C11H10  | 002443-46-1 | 87   |
| 4    |    | 1H-Indene, 1,1-dimethyl-            | 144 | C11H12  | 018636-55-0 | 70   |
| 5    |    | 1H-Indene, 1-ethenyl-2,3-dihydro-   | 144 | C11H12  | 051783-46-1 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU043019\  
Data File : VU031663.D  
Acq On : 01 May 2019 03:21  
Operator : JC/SP  
Sample : K2604-06 5X  
Misc : 5.01µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ80

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

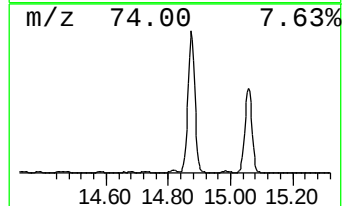
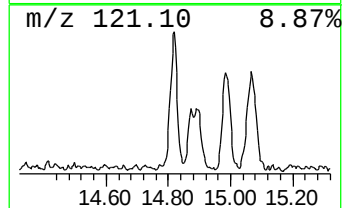
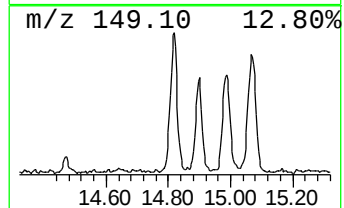
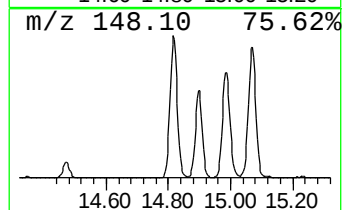
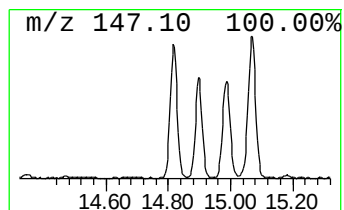
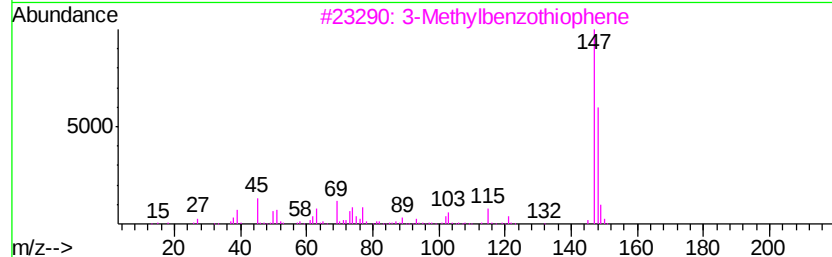
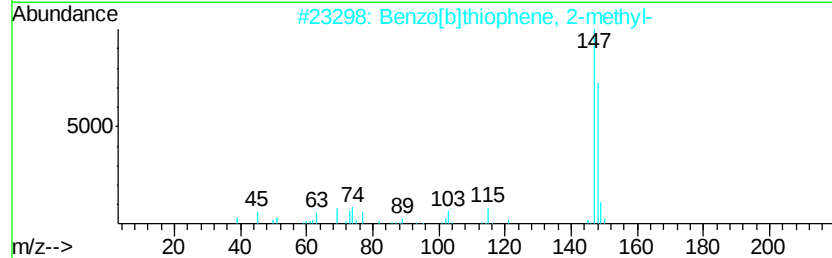
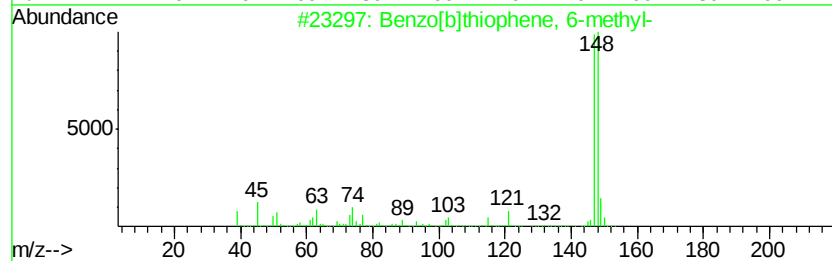
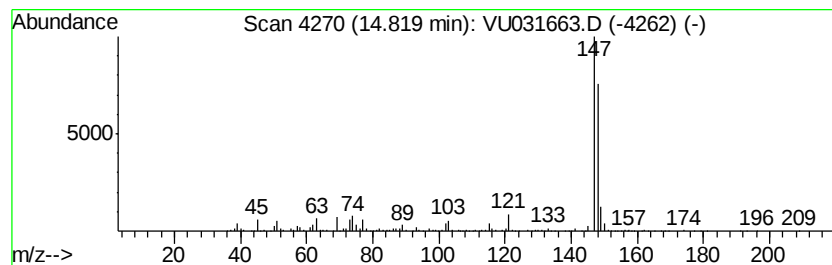
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 Benzo[b]thiophene, 6-methyl- Concentration Rank 44

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 14.82 | 10.04 ug/L | 586604 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[b]thiophene, 6-methyl- | 148 | C9H8S   | 016587-47-6 | 94   |
| 2    |    | Benzo[b]thiophene, 2-methyl- | 148 | C9H8S   | 001195-14-8 | 91   |
| 3    |    | 3-Methylbenzothiophene       | 148 | C9H8S   | 001455-18-1 | 91   |
| 4    |    | Benzo[b]thiophene, 4-methyl- | 148 | C9H8S   | 014315-11-8 | 91   |
| 5    |    | Benzo[b]thiophene, 5-methyl- | 148 | C9H8S   | 014315-14-1 | 91   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU043019\  
 Data File : VU031663.D  
 Acq On : 01 May 2019 03:21  
 Operator : JC/SP  
 Sample : K2604-06 5X  
 Misc : 5.01g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 GAJ80

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
 Quant Title : VOC Analysis

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Benzene, 2-propenyl- | 10.51 | 6.2     | ug/L  | 364763   | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, propyl-     | 10.58 | 7.7     | ug/L  | 450038   | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, 1-ethyl-... | 10.68 | 42.4    | ug/L  | 2479620  | 3                     | 11.48 | 2921190 | 50.0 |
| .alpha.-Methylsty... | 10.99 | 35.8    | ug/L  | 2089510  | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, 1,2,3-tr... | 11.14 | 141.3   | ug/L  | 8252230  | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, 1-etheny... | 11.21 | 220.9   | ug/L  | 12903900 | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, cyclopro... | 11.26 | 80.2    | ug/L  | 4684750  | 3                     | 11.48 | 2921190 | 50.0 |
| Benzofuran           | 11.40 | 2.8     | ug/L  | 162423   | 3                     | 11.48 | 2921190 | 50.0 |
| Bicyclo[4.2.0]oct... | 11.62 | 31.1    | ug/L  | 1816020  | 3                     | 11.48 | 2921190 | 50.0 |
| Indane               | 11.76 | 32.6    | ug/L  | 1904000  | 3                     | 11.48 | 2921190 | 50.0 |
| Indene               | 11.98 | 820.3   | ug/L  | 47923500 | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, 4-ethyl-... | 12.14 | 5.6     | ug/L  | 327161   | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, 1-methyl... | 12.22 | 12.5    | ug/L  | 728979   | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, (2-methy... | 12.42 | 53.6    | ug/L  | 3134360  | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, 2-etheny... | 12.48 | 14.2    | ug/L  | 829554   | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, 1,2,4,5-... | 12.65 | 8.9     | ug/L  | 522070   | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, 1-ethyl-... | 12.70 | 48.2    | ug/L  | 2816390  | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, 1-etheny... | 12.82 | 41.7    | ug/L  | 2433960  | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, 1-etheny... | 12.96 | 9.0     | ug/L  | 528213   | 3                     | 11.48 | 2921190 | 50.0 |
| 1-Dodecanol          | 13.08 | 4.9     | ug/L  | 286265   | 3                     | 11.48 | 2921190 | 50.0 |
| o-Cymene             | 13.12 | 44.6    | ug/L  | 2603240  | 3                     | 11.48 | 2921190 | 50.0 |
| Benzene, (1-methy... | 13.18 | 167.0   | ug/L  | 9757630  | 3                     | 11.48 | 2921190 | 50.0 |
| 2-Methylindene       | 13.29 | 194.0   | ug/L  | 11333100 | 3                     | 11.48 | 2921190 | 50.0 |
| 1,4-Dihydronaphth... | 13.38 | 31.5    | ug/L  | 1838630  | 3                     | 11.48 | 2921190 | 50.0 |
| Benzo[c]thiophene    | 13.86 | 44.4    | ug/L  | 2591740  | 3                     | 11.48 | 2921190 | 50.0 |
| (DEL) Alkane: Str... | 14.14 | 7.8     | ug/L  | 457383   | 3                     | 11.48 | 2921190 | 50.0 |
| 1H-Indene, 1,3-di... | 14.33 | 26.9    | ug/L  | 1572590  | 3                     | 11.48 | 2921190 | 50.0 |
| Bicyclo[4.4.1]und... | 14.58 | 10.5    | ug/L  | 615416   | 3                     | 11.48 | 2921190 | 50.0 |
| Benzocycloheptatr... | 14.73 | 22.6    | ug/L  | 1318900  | 3                     | 11.48 | 2921190 | 50.0 |
| Benzo[b]thiophene... | 14.82 | 10.0    | ug/L  | 586604   | 3                     | 11.48 | 2921190 | 50.0 |