

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
 Data File : VU031702.D  
 Acq On : 02 May 2019 11:15  
 Operator : JC/SP  
 Sample : K2603-16DL 10X  
 Misc : 5.10g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 GAJ68DL

## 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<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 1.395       | 88            | 95          | 107          | rBV      | 336936         | 372669        | 0.54%           | 0.102%        |
| 2         | 1.678       | 176           | 183         | 198          | rBV      | 297791         | 388862        | 0.56%           | 0.106%        |
| 3         | 2.264       | 355           | 365         | 379          | rVB      | 636732         | 983213        | 1.43%           | 0.269%        |
| 4         | 2.469       | 421           | 429         | 431          | rBV7     | 3883           | 4404          | 0.01%           | 0.001%        |
| 5         | 2.540       | 448           | 451         | 456          | rVB3     | 1606           | 1160          | 0.00%           | 0.000%        |
| 6         | 2.646       | 481           | 484         | 487          | rBV3     | 1451           | 1122          | 0.00%           | 0.000%        |
| 7         | 2.891       | 556           | 560         | 561          | rBV      | 1585           | 1249          | 0.00%           | 0.000%        |
| 8         | 2.987       | 586           | 590         | 595          | rVB6     | 1774           | 1694          | 0.00%           | 0.000%        |
| 9         | 3.071       | 612           | 616         | 619          | rBV3     | 1387           | 1317          | 0.00%           | 0.000%        |
| 10        | 3.264       | 674           | 676         | 681          | rVV3     | 1222           | 1109          | 0.00%           | 0.000%        |
| 11        | 3.296       | 681           | 686         | 693          | rVB4     | 2087           | 2577          | 0.00%           | 0.001%        |
| 12        | 3.337       | 693           | 699         | 702          | rBV3     | 1838           | 2369          | 0.00%           | 0.001%        |
| 13        | 3.379       | 710           | 712         | 717          | rVB      | 1563           | 1205          | 0.00%           | 0.000%        |
| 14        | 3.495       | 744           | 748         | 750          | rBV4     | 2371           | 1994          | 0.00%           | 0.001%        |
| 15        | 3.543       | 758           | 763         | 764          | rBV3     | 1178           | 933           | 0.00%           | 0.000%        |
| 16        | 3.611       | 782           | 784         | 790          | rVB2     | 1927           | 1131          | 0.00%           | 0.000%        |
| 17        | 3.669       | 799           | 802         | 806          | rVB3     | 1816           | 1486          | 0.00%           | 0.000%        |
| 18        | 3.714       | 810           | 816         | 822          | rBV5     | 1075           | 1305          | 0.00%           | 0.000%        |
| 19        | 3.784       | 833           | 838         | 842          | rBV2     | 1756           | 1952          | 0.00%           | 0.001%        |
| 20        | 3.833       | 849           | 853         | 857          | rBV2     | 1373           | 1281          | 0.00%           | 0.000%        |
| 21        | 4.019       | 908           | 911         | 916          | rVB3     | 1089           | 1040          | 0.00%           | 0.000%        |
| 22        | 4.058       | 916           | 923         | 924          | rBV4     | 1255           | 1414          | 0.00%           | 0.000%        |
| 23        | 4.196       | 952           | 966         | 998          | rBV      | 338741         | 1152650       | 1.67%           | 0.315%        |
| 24        | 4.643       | 1090          | 1105        | 1130         | rBV2     | 559942         | 1324174       | 1.92%           | 0.362%        |
| 25        | 4.845       | 1164          | 1168        | 1172         | rBV2     | 2090           | 1521          | 0.00%           | 0.000%        |
| 26        | 5.148       | 1259          | 1262        | 1266         | rBV3     | 1373           | 1146          | 0.00%           | 0.000%        |
| 27        | 5.235       | 1285          | 1289        | 1293         | rVB3     | 1215           | 1087          | 0.00%           | 0.000%        |
| 28        | 5.334       | 1297          | 1320        | 1356         | rBV2     | 1150866        | 3505957       | 5.09%           | 0.958%        |
| 29        | 5.881       | 1476          | 1490        | 1514         | rBV      | 1233208        | 2498016       | 3.62%           | 0.683%        |
| 30        | 6.180       | 1572          | 1583        | 1596         | rBV9     | 12964          | 30111         | 0.04%           | 0.008%        |
| 31        | 6.328       | 1613          | 1629        | 1655         | rBV      | 793961         | 1660543       | 2.41%           | 0.454%        |
| 32        | 6.604       | 1711          | 1715        | 1718         | rBV4     | 1153           | 999           | 0.00%           | 0.000%        |
| 33        | 6.627       | 1718          | 1722        | 1727         | rVV4     | 1374           | 1449          | 0.00%           | 0.000%        |
| 34        | 6.678       | 1732          | 1738        | 1743         | rBV7     | 1788           | 1843          | 0.00%           | 0.001%        |

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 GAJ68DL

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 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 | 6.762  | 1761 | 1764 | 1768 | rVB2  | 1668    | 1160    | 0.00% | 0.000% |
| 36 | 6.784  | 1768 | 1771 | 1776 | rVB6  | 1060    | 959     | 0.00% | 0.000% |
| 37 | 6.807  | 1776 | 1778 | 1780 | rBV3  | 1724    | 976     | 0.00% | 0.000% |
| 38 | 6.878  | 1797 | 1800 | 1802 | rBV4  | 2024    | 1626    | 0.00% | 0.000% |
| 39 | 7.074  | 1857 | 1861 | 1864 | rVB3  | 1447    | 1160    | 0.00% | 0.000% |
| 40 | 7.106  | 1867 | 1871 | 1875 | rBV5  | 1908    | 1867    | 0.00% | 0.001% |
| 41 | 7.177  | 1888 | 1893 | 1896 | rBV6  | 2572    | 2231    | 0.00% | 0.001% |
| 42 | 7.225  | 1896 | 1908 | 1937 | rBV   | 431776  | 817936  | 1.19% | 0.224% |
| 43 | 7.427  | 1963 | 1971 | 1980 | rBV10 | 3780    | 6890    | 0.01% | 0.002% |
| 44 | 7.562  | 1999 | 2013 | 2026 | rBV   | 1392872 | 2484375 | 3.60% | 0.679% |
| 45 | 7.633  | 2027 | 2035 | 2062 | rVB   | 314340  | 631594  | 0.92% | 0.173% |
| 46 | 7.849  | 2087 | 2102 | 2140 | rVB   | 290346  | 575323  | 0.83% | 0.157% |
| 47 | 8.225  | 2211 | 2219 | 2226 | rBV7  | 4169    | 6001    | 0.01% | 0.002% |
| 48 | 8.315  | 2236 | 2247 | 2284 | rBV   | 1341350 | 2677101 | 3.88% | 0.732% |
| 49 | 8.604  | 2328 | 2337 | 2346 | rBV6  | 20532   | 36563   | 0.05% | 0.010% |
| 50 | 8.752  | 2374 | 2383 | 2391 | rBV9  | 14511   | 26468   | 0.04% | 0.007% |
| 51 | 8.906  | 2424 | 2431 | 2442 | rVB3  | 40543   | 61780   | 0.09% | 0.017% |
| 52 | 9.029  | 2460 | 2469 | 2476 | rBV2  | 50250   | 80625   | 0.12% | 0.022% |
| 53 | 9.086  | 2476 | 2487 | 2524 | rVB   | 1750718 | 3054335 | 4.43% | 0.835% |
| 54 | 9.247  | 2525 | 2537 | 2561 | rBV   | 1648522 | 2778108 | 4.03% | 0.759% |
| 55 | 9.373  | 2566 | 2576 | 2589 | rVV   | 946389  | 1707428 | 2.48% | 0.467% |
| 56 | 9.440  | 2590 | 2597 | 2617 | rVB   | 382807  | 625945  | 0.91% | 0.171% |
| 57 | 9.646  | 2656 | 2661 | 2668 | rBV7  | 10567   | 14258   | 0.02% | 0.004% |
| 58 | 9.781  | 2691 | 2703 | 2737 | rVB3  | 1718186 | 3635869 | 5.27% | 0.994% |
| 59 | 9.948  | 2748 | 2755 | 2766 | rVB2  | 207179  | 312920  | 0.45% | 0.086% |
| 60 | 10.041 | 2776 | 2784 | 2792 | rBV5  | 138865  | 237108  | 0.34% | 0.065% |
| 61 | 10.430 | 2894 | 2905 | 2917 | rBV   | 1173570 | 2153474 | 3.12% | 0.588% |
| 62 | 10.585 | 2946 | 2953 | 2961 | rBV   | 165441  | 239773  | 0.35% | 0.066% |
| 63 | 10.704 | 2976 | 2990 | 2998 | rBV   | 666777  | 1240762 | 1.80% | 0.339% |
| 64 | 10.810 | 3016 | 3023 | 3035 | rVB   | 1639795 | 2404670 | 3.49% | 0.657% |
| 65 | 10.971 | 3064 | 3073 | 3076 | rBV2  | 398744  | 606094  | 0.88% | 0.166% |
| 66 | 11.112 | 3109 | 3117 | 3119 | rBV   | 458207  | 546190  | 0.79% | 0.149% |
| 67 | 11.144 | 3120 | 3127 | 3138 | rVV   | 2899077 | 4450257 | 6.46% | 1.216% |
| 68 | 11.212 | 3139 | 3148 | 3157 | rVV   | 3004125 | 4475597 | 6.49% | 1.223% |
| 69 | 11.263 | 3157 | 3164 | 3175 | rVB   | 994115  | 1479479 | 2.15% | 0.404% |
| 70 | 11.395 | 3197 | 3205 | 3212 | rBV3  | 436858  | 733369  | 1.06% | 0.200% |
| 71 | 11.479 | 3223 | 3231 | 3241 | rBV   | 1888791 | 2850630 | 4.14% | 0.779% |

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 GAJ68DL

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 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  | 11.565 | 3251 | 3258 | 3266 | rBV  | 1229500  | 1696817  | 2.46%   | 0.464%  |
| 73  | 11.765 | 3312 | 3320 | 3327 | rBV2 | 542855   | 967421   | 1.40%   | 0.264%  |
| 74  | 11.858 | 3340 | 3349 | 3366 | rVB4 | 2226175  | 4313861  | 6.26%   | 1.179%  |
| 75  | 11.977 | 3375 | 3386 | 3394 | rBV  | 8935162  | 14082719 | 20.43%  | 3.848%  |
| 76  | 12.019 | 3394 | 3399 | 3408 | rVB  | 3463871  | 4615068  | 6.69%   | 1.261%  |
| 77  | 12.353 | 3497 | 3503 | 3510 | rBV3 | 510011   | 797191   | 1.16%   | 0.218%  |
| 78  | 12.485 | 3537 | 3544 | 3547 | rBV5 | 461611   | 660567   | 0.96%   | 0.181%  |
| 79  | 12.607 | 3574 | 3582 | 3587 | rBV  | 991478   | 1505339  | 2.18%   | 0.411%  |
| 80  | 12.700 | 3603 | 3611 | 3630 | rBV5 | 1427876  | 3752488  | 5.44%   | 1.025%  |
| 81  | 12.819 | 3642 | 3648 | 3660 | rVB3 | 957443   | 1421598  | 2.06%   | 0.388%  |
| 82  | 13.118 | 3733 | 3741 | 3751 | rBV2 | 6132896  | 9206371  | 13.35%  | 2.516%  |
| 83  | 13.180 | 3753 | 3760 | 3772 | rVB  | 4229471  | 6456042  | 9.37%   | 1.764%  |
| 84  | 13.286 | 3783 | 3793 | 3812 | rBV2 | 6095669  | 11214746 | 16.27%  | 3.065%  |
| 85  | 13.739 | 3923 | 3934 | 3944 | rVB2 | 38372965 | 61179748 | 88.75%  | 16.719% |
| 86  | 14.138 | 4050 | 4058 | 4065 | rBV2 | 3214086  | 4614796  | 6.69%   | 1.261%  |
| 87  | 14.334 | 4111 | 4119 | 4129 | rBV6 | 2052518  | 3677101  | 5.33%   | 1.005%  |
| 88  | 14.874 | 4276 | 4287 | 4300 | rBV2 | 44416279 | 68935787 | 100.00% | 18.839% |
| 89  | 15.057 | 4334 | 4344 | 4362 | rBV  | 21937125 | 35473190 | 51.46%  | 9.694%  |
| 90  | 15.794 | 4566 | 4573 | 4581 | rBV  | 3173430  | 4675622  | 6.78%   | 1.278%  |
| 91  | 15.909 | 4598 | 4609 | 4622 | rBV  | 12909658 | 21814166 | 31.64%  | 5.961%  |
| 92  | 16.054 | 4644 | 4654 | 4661 | rBV  | 11464386 | 18045728 | 26.18%  | 4.931%  |
| 93  | 16.093 | 4662 | 4666 | 4683 | rVB  | 6116747  | 8346161  | 12.11%  | 2.281%  |
| 94  | 16.183 | 4686 | 4694 | 4706 | rVB  | 2662901  | 4202348  | 6.10%   | 1.148%  |
| 95  | 16.279 | 4709 | 4724 | 4734 | rBV  | 3298738  | 7378616  | 10.70%  | 2.016%  |
| 96  | 16.427 | 4763 | 4770 | 4781 | rVB  | 1471337  | 2200806  | 3.19%   | 0.601%  |
| 97  | 16.520 | 4790 | 4799 | 4815 | rVB2 | 3253388  | 6007658  | 8.71%   | 1.642%  |
| 98  | 17.012 | 4946 | 4952 | 4976 | rVB2 | 1422240  | 2682519  | 3.89%   | 0.733%  |
| 99  | 17.160 | 4991 | 4998 | 5008 | rVB  | 1162522  | 1792078  | 2.60%   | 0.490%  |
| 100 | 17.224 | 5010 | 5018 | 5028 | rBV2 | 823664   | 1295820  | 1.88%   | 0.354%  |

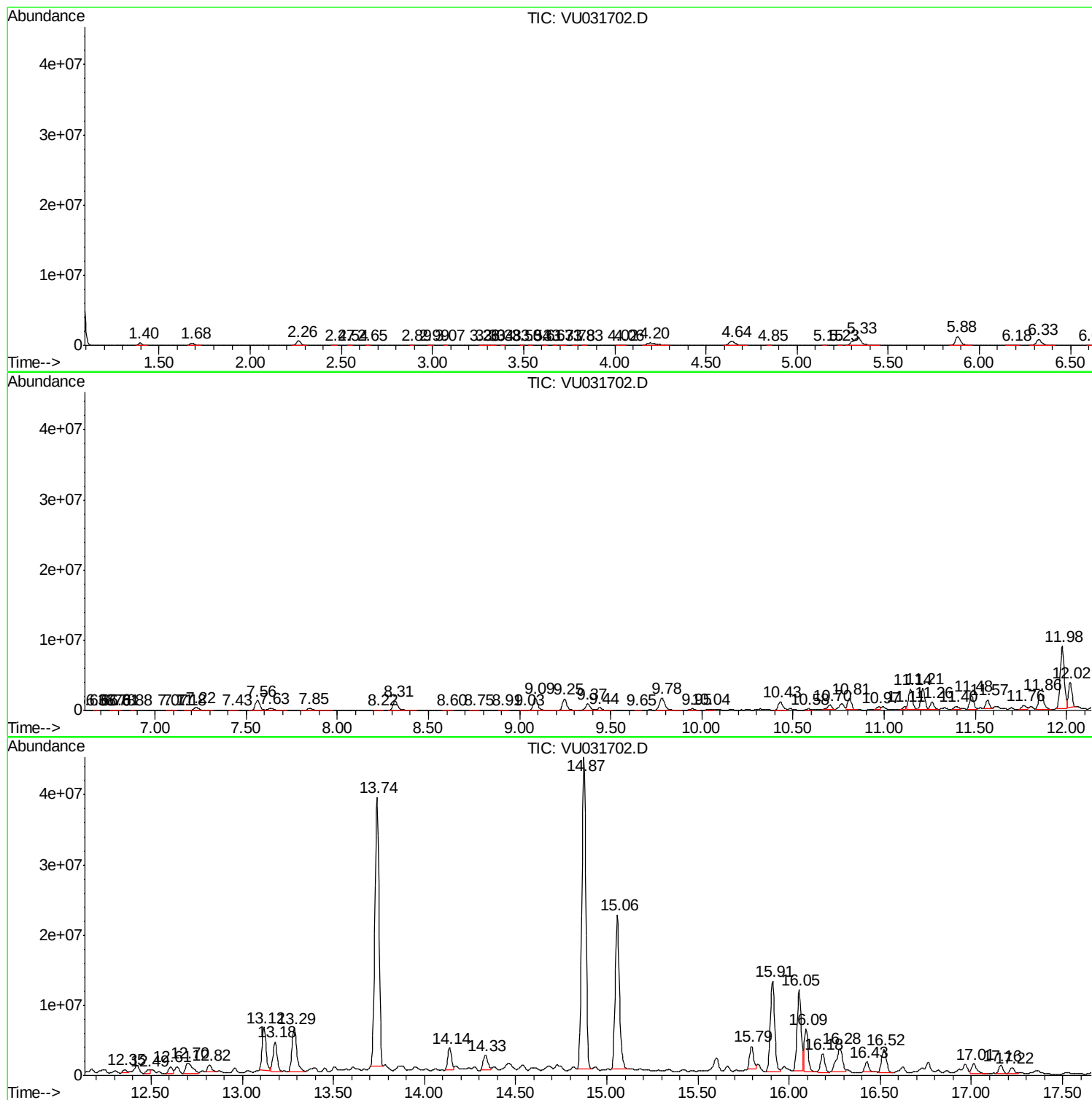
Sum of corrected areas: 365928255

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ClientSampleId :  
GAJ68DL

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM042719WMA.M  
Quant Title : VOC Analysis

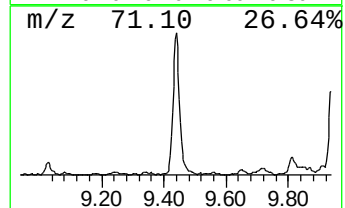
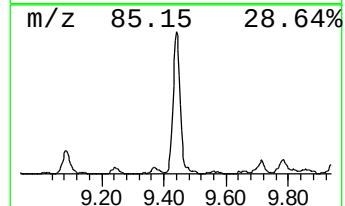
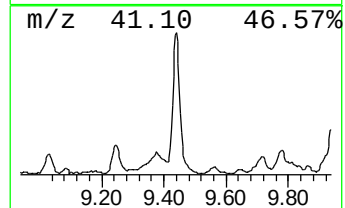
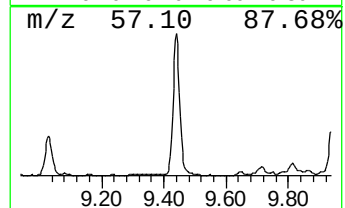
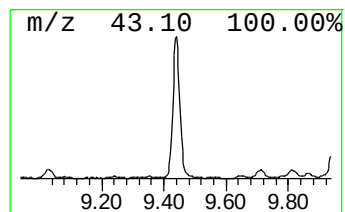
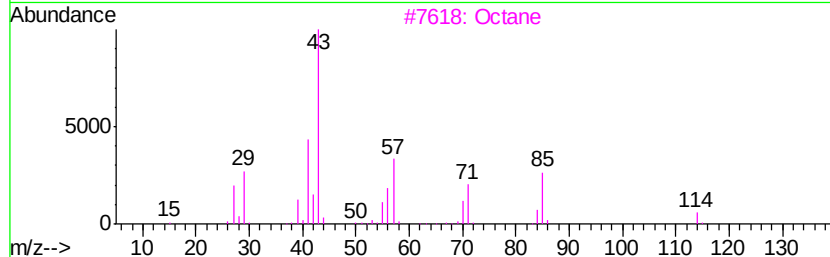
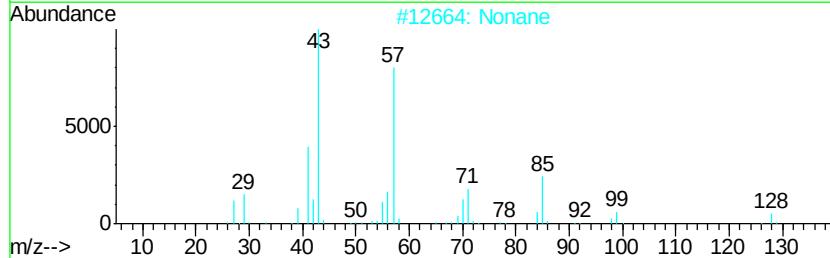
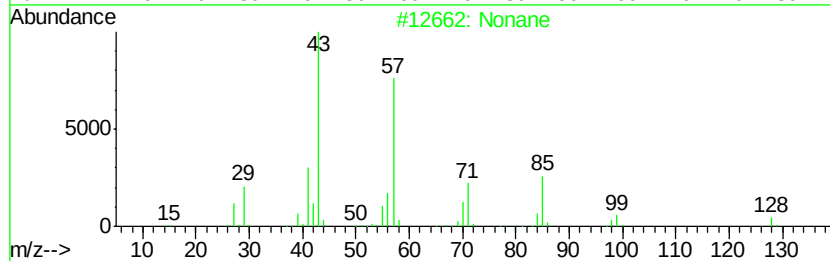
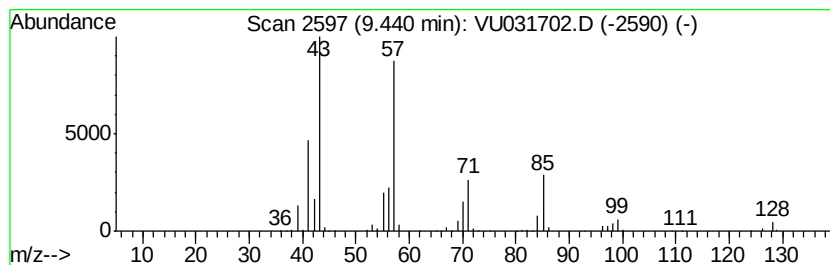
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 37

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 9.44 | 10.25 ug/L | 625945 | Chlorobenzene-d5 | 9.09 |

| Hit# | of | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|----|------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Nonane                             | 128 | C9H20    | 000111-84-2  | 97   |
| 2    |    | Nonane                             | 128 | C9H20    | 000111-84-2  | 90   |
| 3    |    | Octane                             | 114 | C8H18    | 000111-65-9  | 72   |
| 4    |    | Oxalic acid, isohexyl pentyl ester | 244 | C13H24O4 | 1000309-32-8 | 64   |
| 5    |    | Pentadecane                        | 212 | C15H32   | 000629-62-9  | 59   |



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

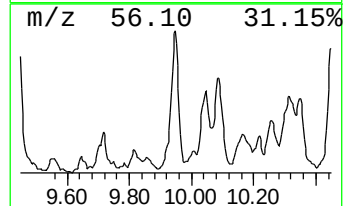
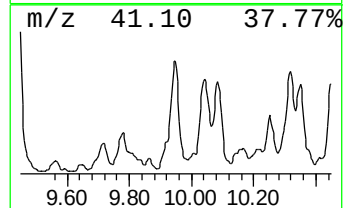
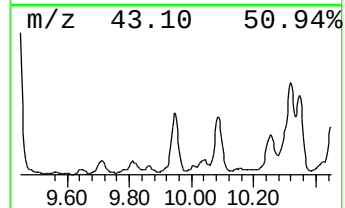
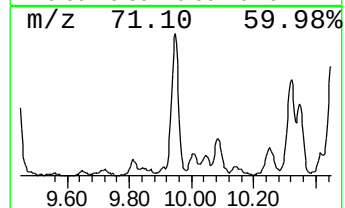
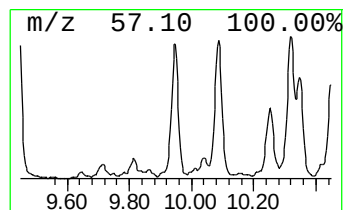
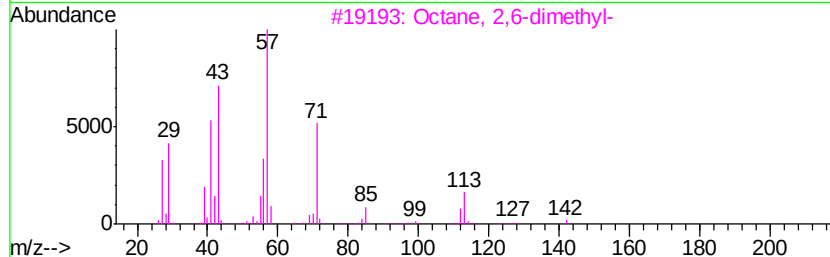
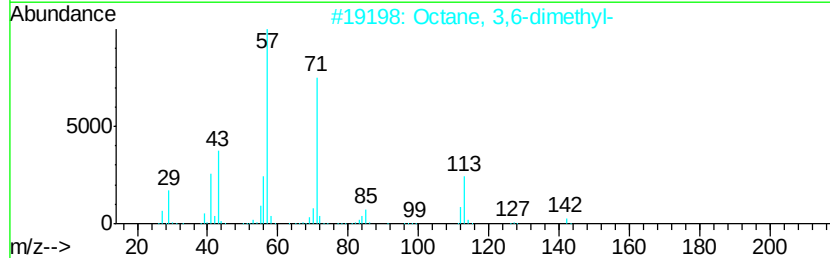
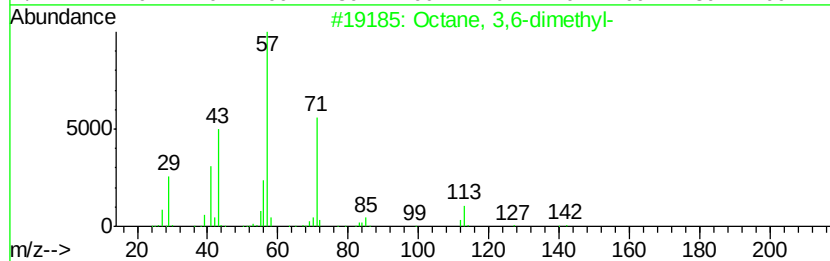
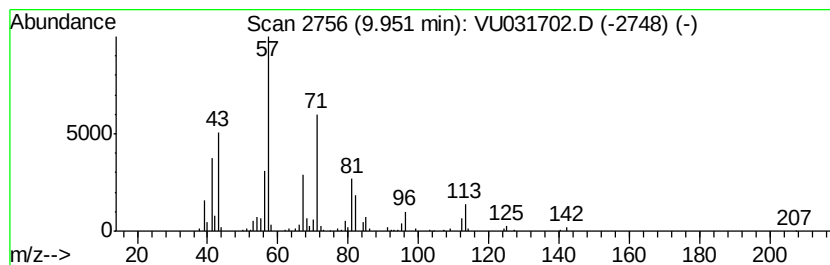
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 39

| R.T. | EstConc   | Area   | Relative to ISTD | R.T. |
|------|-----------|--------|------------------|------|
| 9.95 | 5.12 ug/L | 312920 | Chlorobenzene-d5 | 9.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Octane, 3,6-dimethyl-               | 142 | C10H22    | 015869-94-0  | 62   |
| 2    |    | Octane, 3,6-dimethyl-               | 142 | C10H22    | 015869-94-0  | 60   |
| 3    |    | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 59   |
| 4    |    | Tridecane, 7-methyl-                | 198 | C14H30    | 026730-14-3  | 47   |
| 5    |    | Sulfurous acid, 2-ethylhexyl hex... | 278 | C14H30O3S | 1000309-20-2 | 47   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

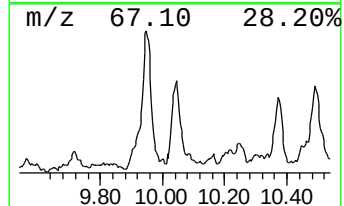
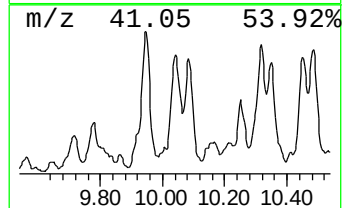
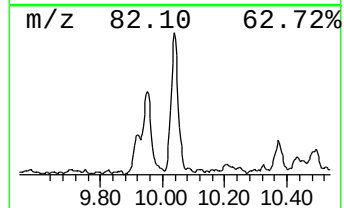
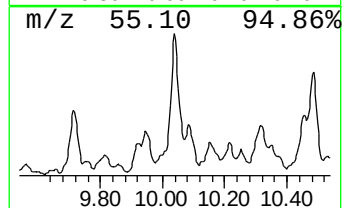
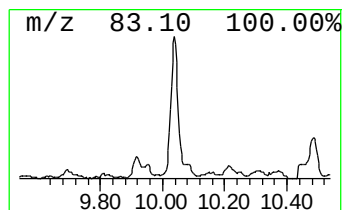
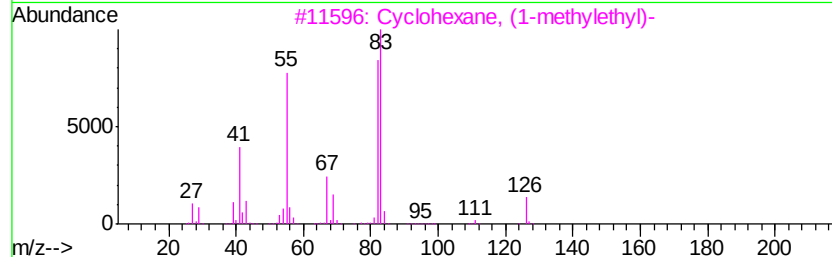
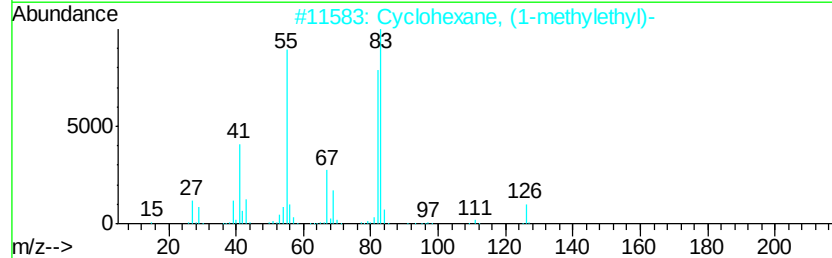
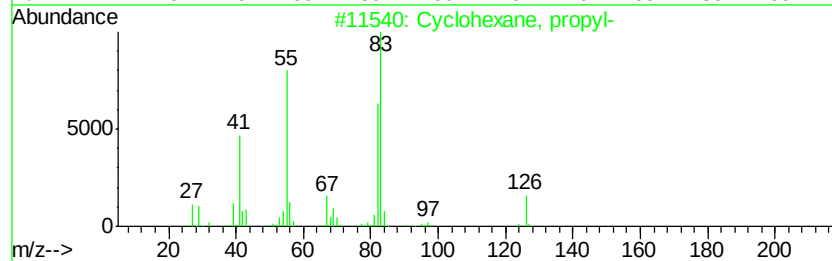
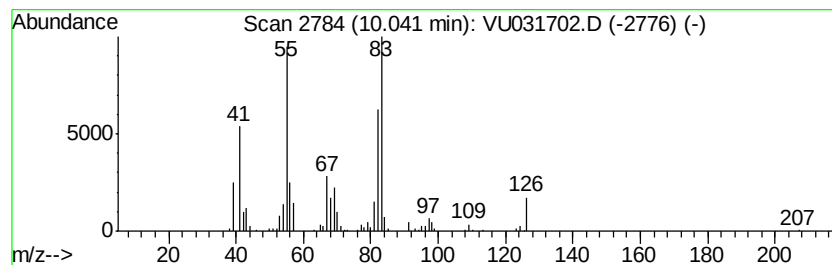
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ClientSampleId :  
GAJ68DL

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 4 (DEL) Alkane: Cyclic10.04 Concentration Rank 41

| R.T.      | EstConc                       | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------|--------|------------------|-------------|------|
| 10.04     | 3.88 ug/L                     | 237108 | Chlorobenzene-d5 | 9.09        |      |
| Hit# of 5 | Tentative ID                  | MW     | MolForm          | CAS#        | Qual |
| 1         | Cyclohexane, propyl-          | 126    | C9H18            | 001678-92-8 | 80   |
| 2         | Cyclohexane, (1-methylethyl)- | 126    | C9H18            | 000696-29-7 | 80   |
| 3         | Cyclohexane, (1-methylethyl)- | 126    | C9H18            | 000696-29-7 | 74   |
| 4         | Cyclohexane, propyl-          | 126    | C9H18            | 001678-92-8 | 72   |
| 5         | Cyclohexane, propyl-          | 126    | C9H18            | 001678-92-8 | 72   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

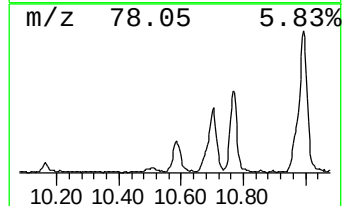
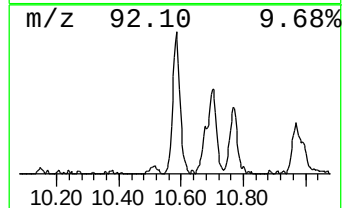
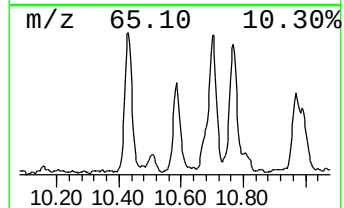
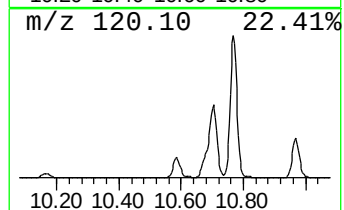
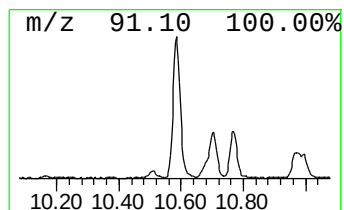
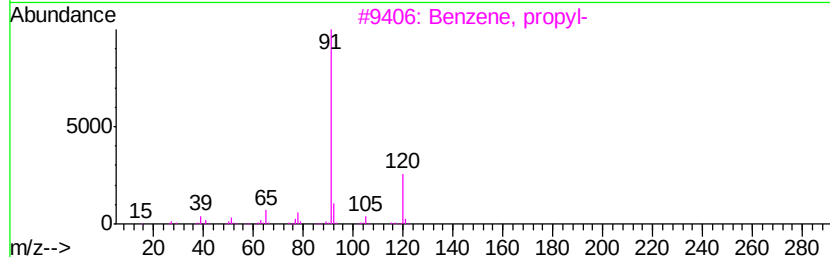
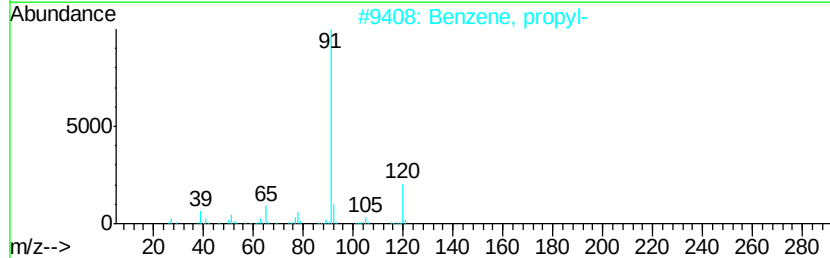
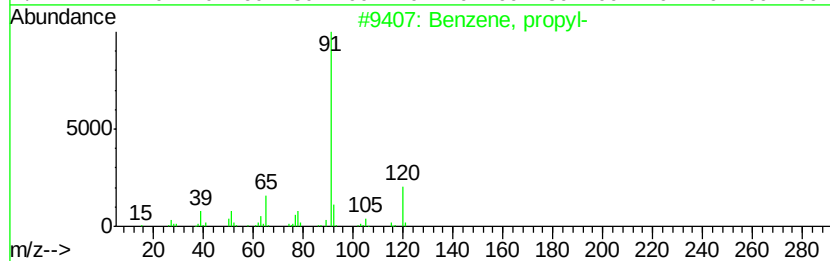
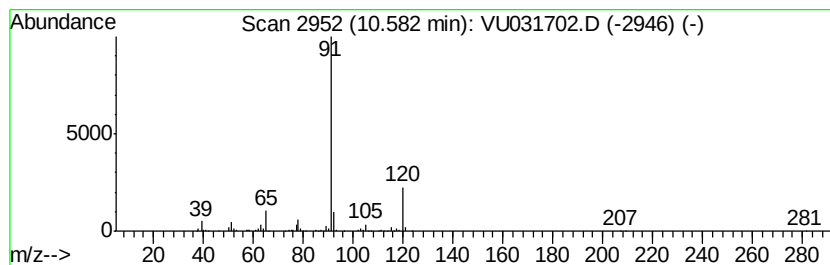
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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 5 Benzene, propyl- Concentration Rank 40

| R.T.    | EstConc   | Area                              | Relative to ISTD       | R.T.           |
|---------|-----------|-----------------------------------|------------------------|----------------|
| 10.58   | 4.21 ug/L | 239773                            | 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       |           | Benzene, propyl-                  | 120 C9H12              | 000103-65-1 83 |
| 4       |           | N-Benzyl-2-phenethylamine         | 211 C15H17N            | 003647-71-0 78 |
| 5       |           | Ethanol, 2-[(phenylmethyl)amino]- | 151 C9H13NO            | 000104-63-2 64 |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

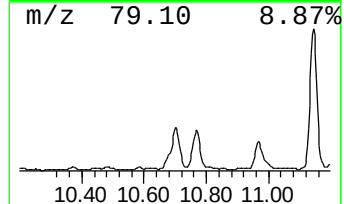
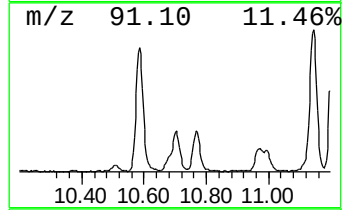
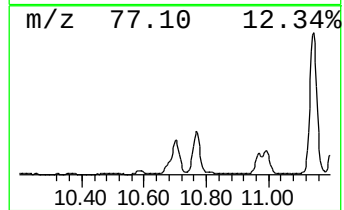
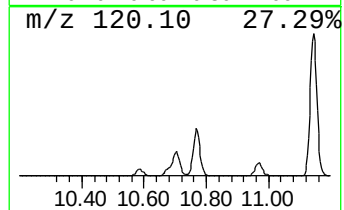
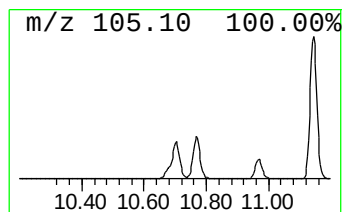
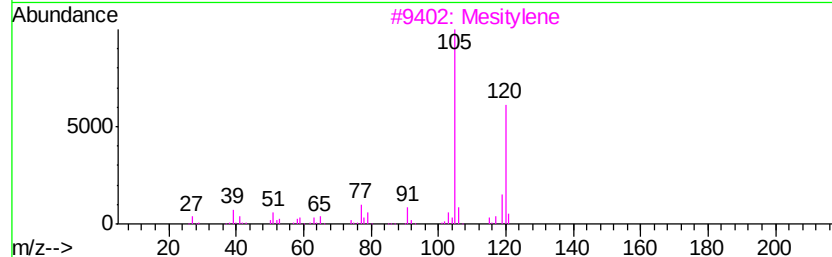
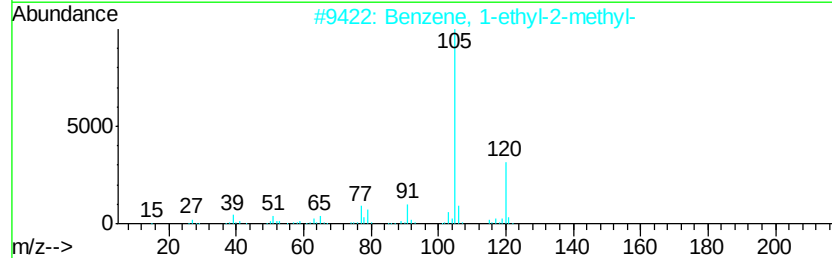
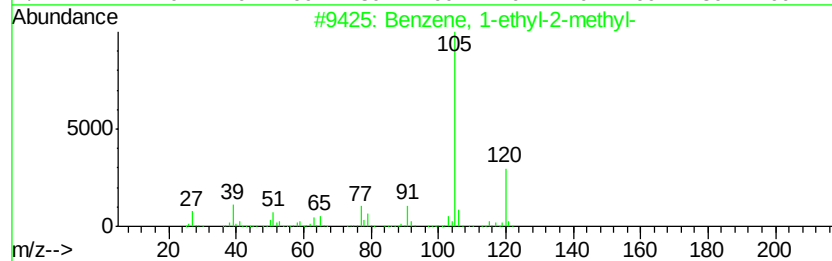
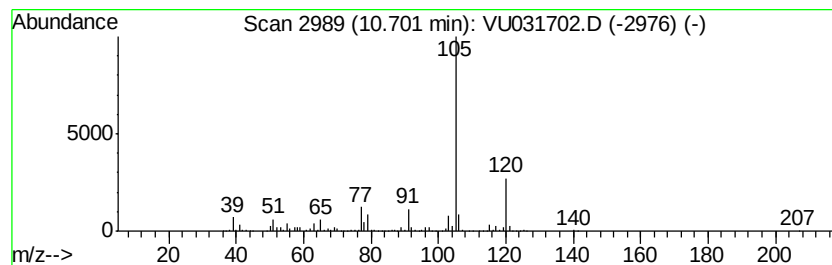
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MSVOA\_U  
ClientSampleId :  
GAJ68DL

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 Benzene, 1-ethyl-2-methyl- Concentration Rank 30

| R.T.    | EstConc    | Area                       | Relative to ISTD       | R.T.           |
|---------|------------|----------------------------|------------------------|----------------|
| 10.70   | 21.76 ug/L | 1240760                    | 1,4-Dichlorobenzene-d4 | 11.48          |
| Hit# of | 5          | Tentative ID               | MW MolForm             | CAS# Qual      |
| 1       |            | Benzene, 1-ethyl-2-methyl- | 120 C9H12              | 000611-14-3 95 |
| 2       |            | Benzene, 1-ethyl-2-methyl- | 120 C9H12              | 000611-14-3 94 |
| 3       |            | Mesitylene                 | 120 C9H12              | 000108-67-8 91 |
| 4       |            | Benzene, 1-ethyl-4-methyl- | 120 C9H12              | 000622-96-8 91 |
| 5       |            | Benzene, 1,2,4-trimethyl-  | 120 C9H12              | 000095-63-6 91 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

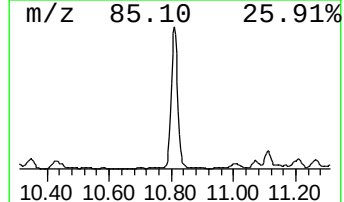
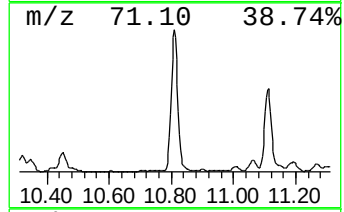
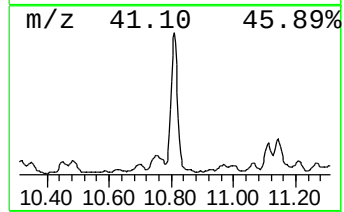
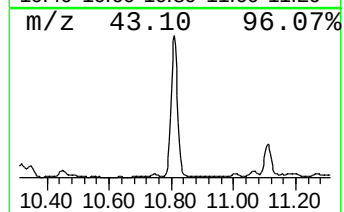
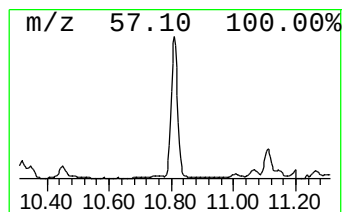
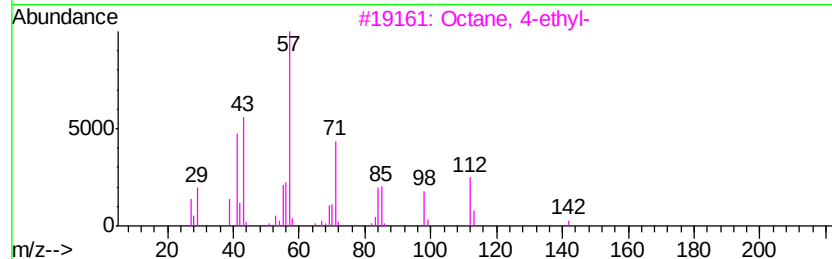
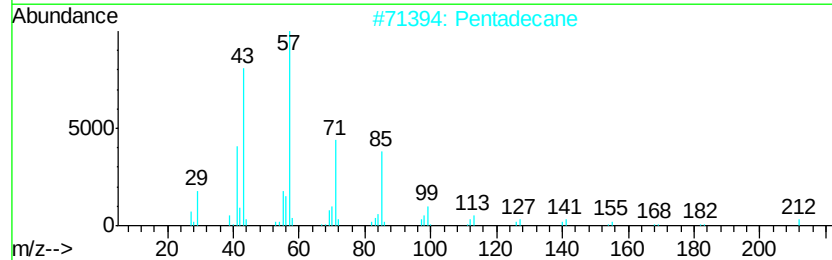
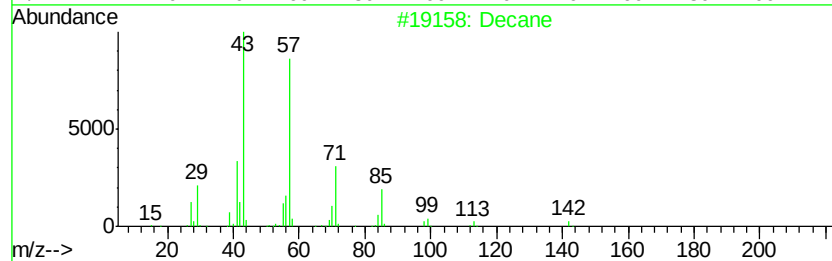
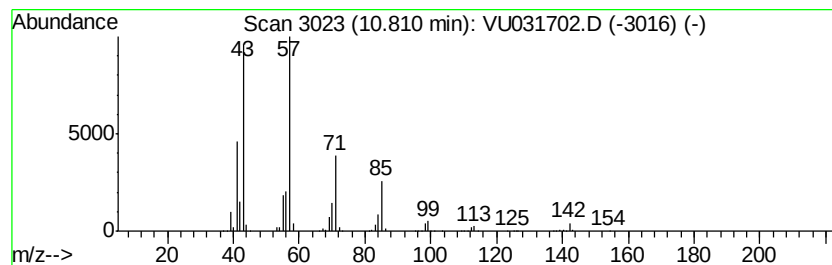
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T.    | EstConc               | Area         | Relative to ISTD       | R.T.           |
|---------|-----------------------|--------------|------------------------|----------------|
| 10.81   | 42.18 ug/L            | 2404670      | 1,4-Dichlorobenzene-d4 | 11.48          |
| Hit# of | 5                     | Tentative ID | MW MolForm             | CAS# Qual      |
| 1       | Decane                |              | 142 C10H22             | 000124-18-5 87 |
| 2       | Pentadecane           |              | 212 C15H32             | 000629-62-9 78 |
| 3       | Octane, 4-ethyl-      |              | 142 C10H22             | 015869-86-0 74 |
| 4       | Hexadecane            |              | 226 C16H34             | 000544-76-3 72 |
| 5       | Decane, 2,9-dimethyl- |              | 170 C12H26             | 001002-17-1 72 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

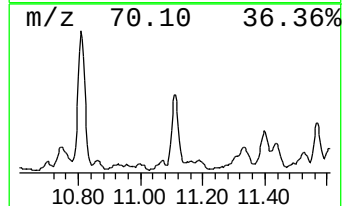
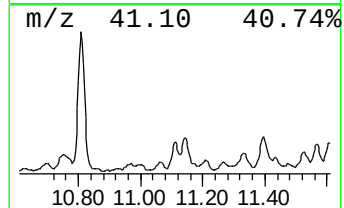
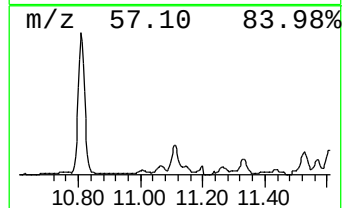
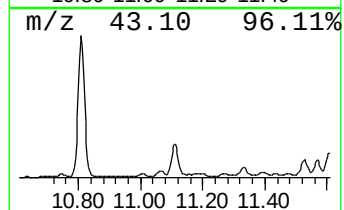
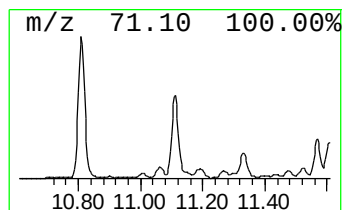
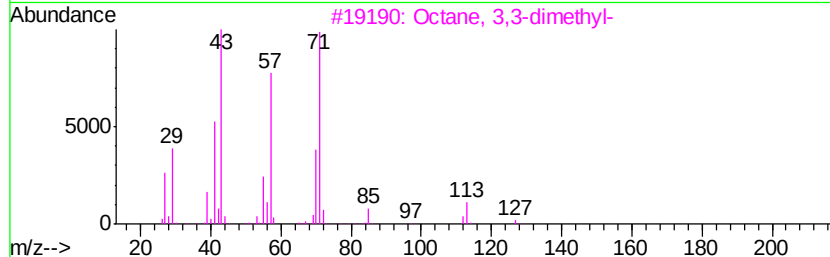
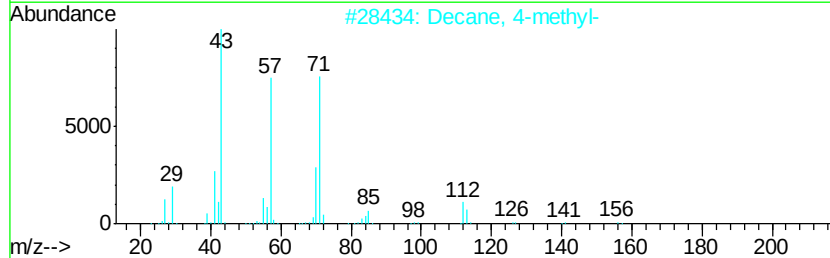
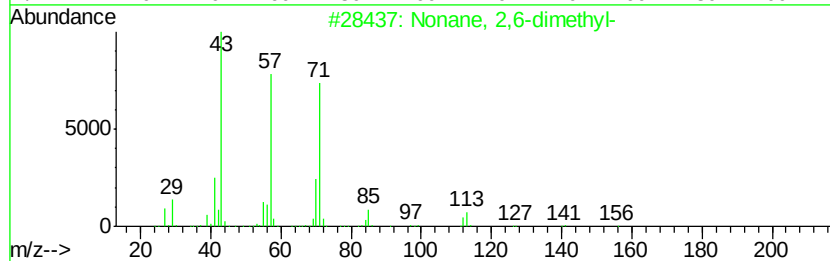
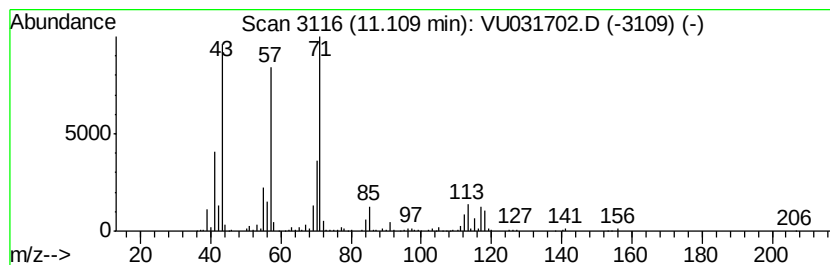
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 38

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 11.11 | 9.58 ug/L | 546190 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 2,6-dimethyl-     | 156 | C11H24  | 017302-28-2 | 87   |
| 2    |    | Decane, 4-methyl-         | 156 | C11H24  | 002847-72-5 | 81   |
| 3    |    | Octane, 3,3-dimethyl-     | 142 | C10H22  | 004110-44-5 | 72   |
| 4    |    | Heptane, 3,3,5-trimethyl- | 142 | C10H22  | 007154-80-5 | 64   |
| 5    |    | Decane, 4-methyl-         | 156 | C11H24  | 002847-72-5 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

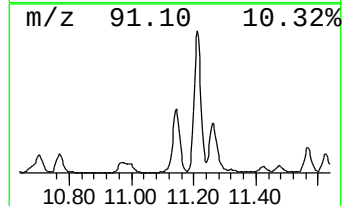
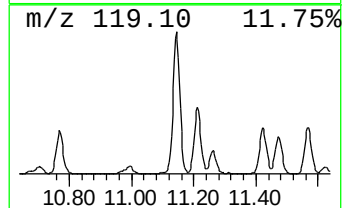
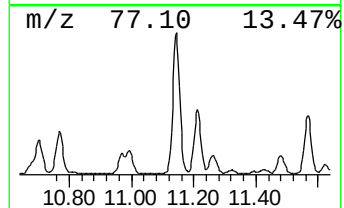
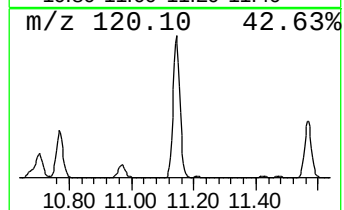
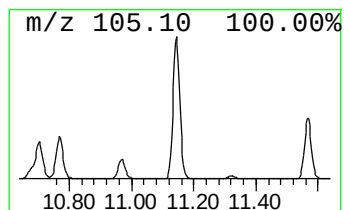
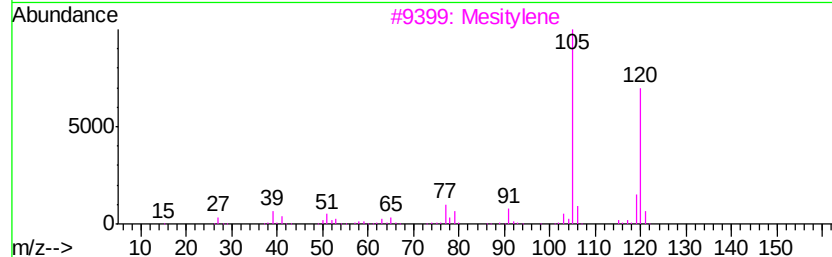
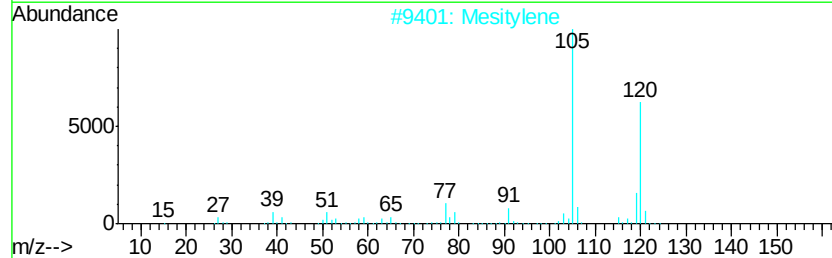
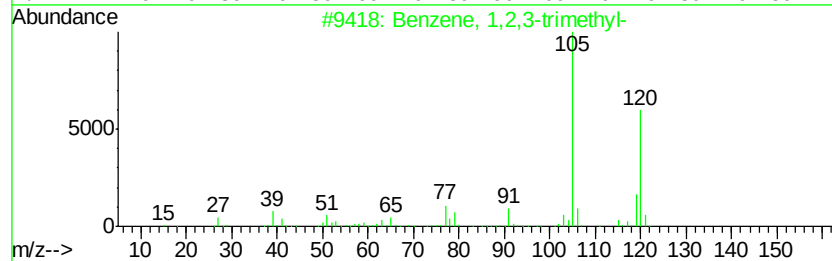
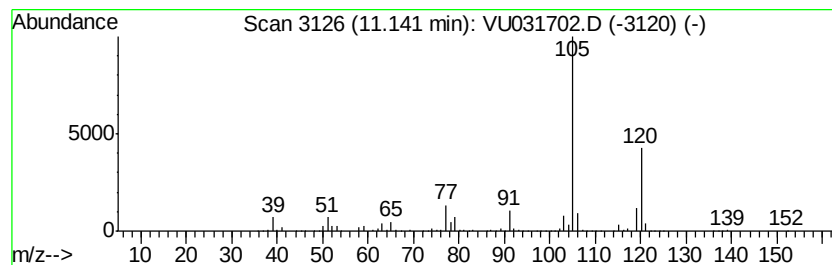
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ClientSampleId :  
GAJ68DL

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 Benzene, 1,2,3-trimethyl- Concentration Rank 17

| R.T.    | EstConc    | Area                      | Relative to ISTD       | R.T.           |
|---------|------------|---------------------------|------------------------|----------------|
| 11.14   | 78.06 ug/L | 4450260                   | 1,4-Dichlorobenzene-d4 | 11.48          |
| Hit# of | 5          | Tentative ID              | MW MolForm             | CAS# Qual      |
| 1       |            | Benzene, 1,2,3-trimethyl- | 120 C9H12              | 000526-73-8 95 |
| 2       |            | Mesitylene                | 120 C9H12              | 000108-67-8 94 |
| 3       |            | Mesitylene                | 120 C9H12              | 000108-67-8 94 |
| 4       |            | Mesitylene                | 120 C9H12              | 000108-67-8 91 |
| 5       |            | Benzene, 1,2,4-trimethyl- | 120 C9H12              | 000095-63-6 91 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

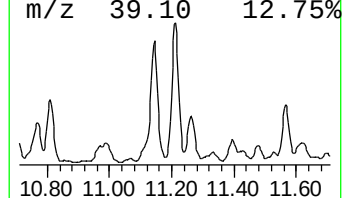
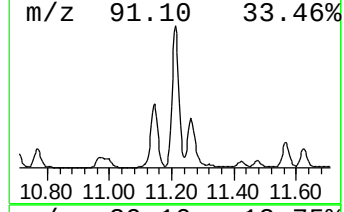
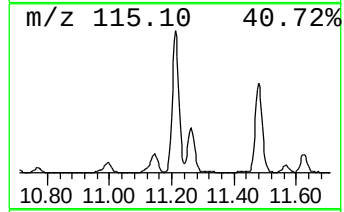
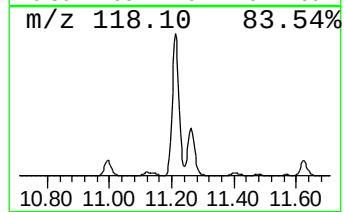
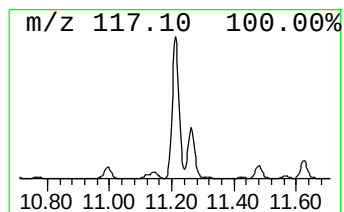
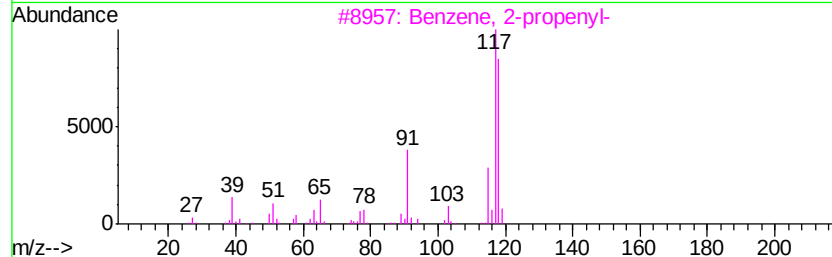
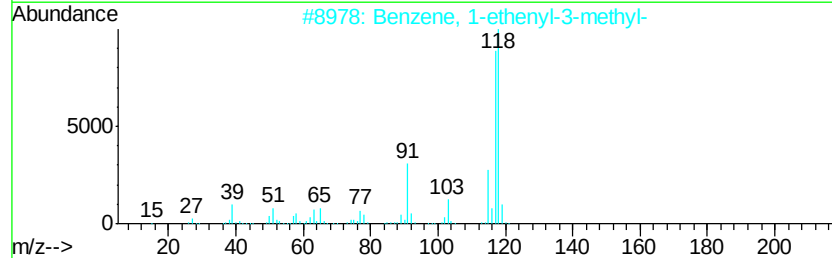
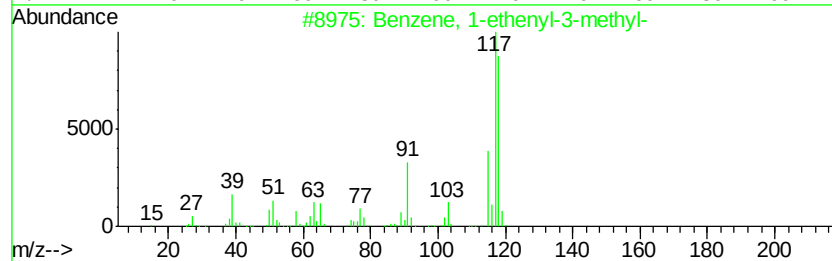
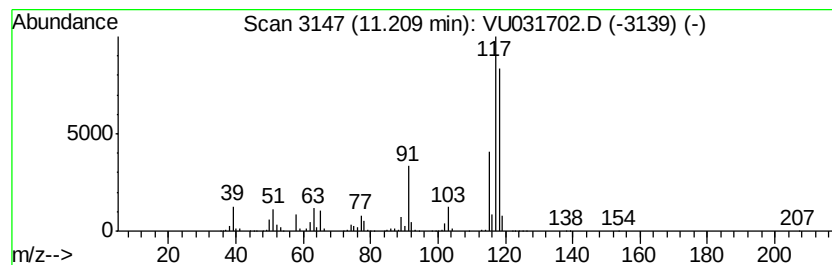
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Benzene, 1-ethenyl-3-methyl- Concentration Rank 16

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 11.21 | 78.50 ug/L | 4475600 | 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 | 95   |
| 5    |    | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
GAJ68DL

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

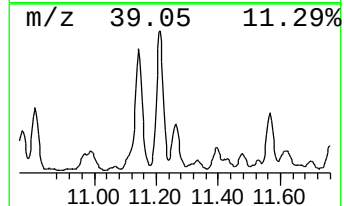
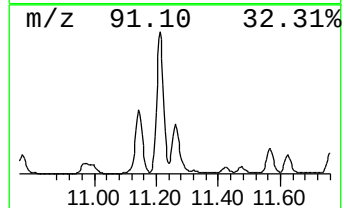
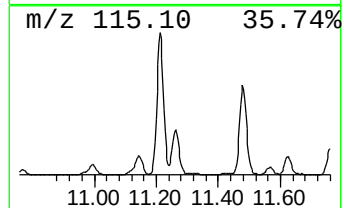
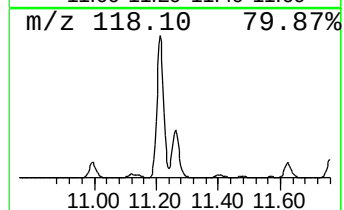
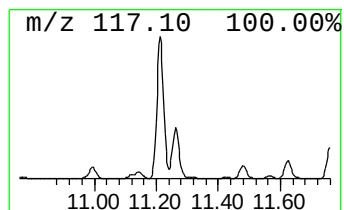
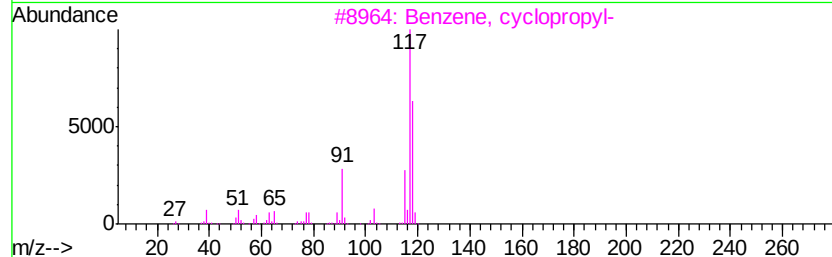
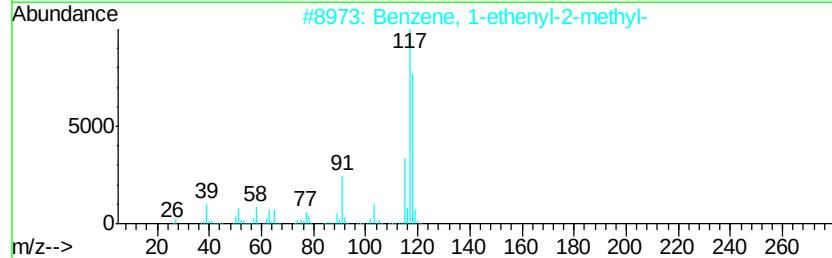
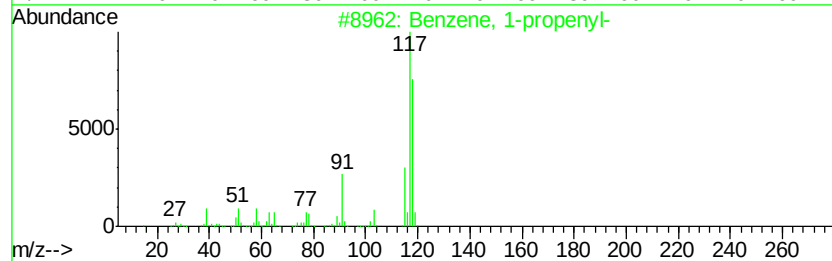
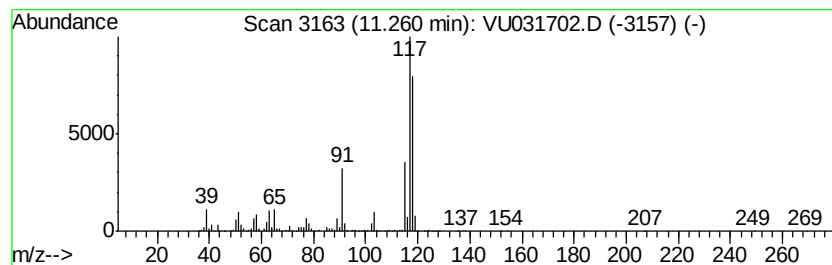
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Benzene, 1-propenyl- Concentration Rank 27

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 11.26 | 25.95 ug/L | 1479480 | 1,4-Dichlorobenzene-d4 | 11.48 |

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

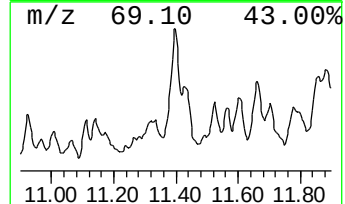
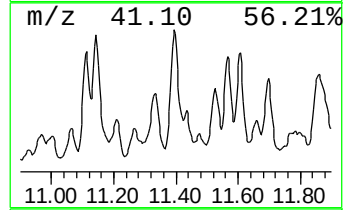
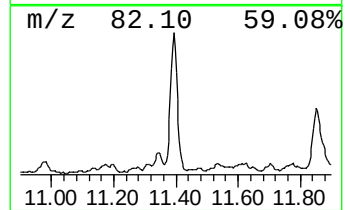
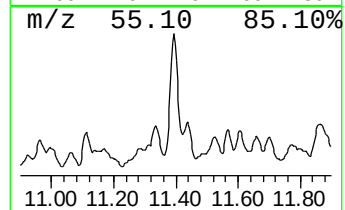
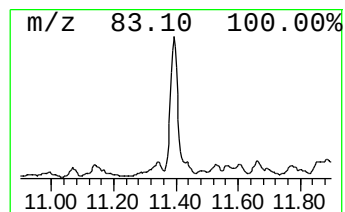
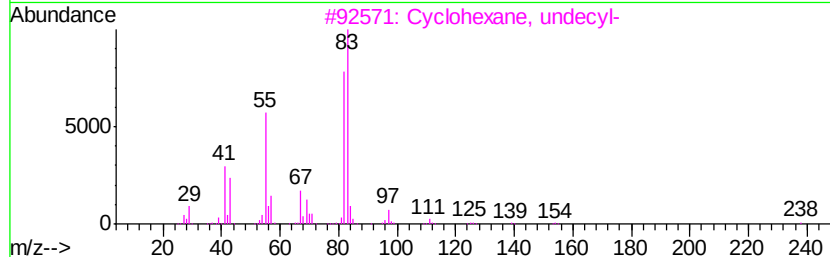
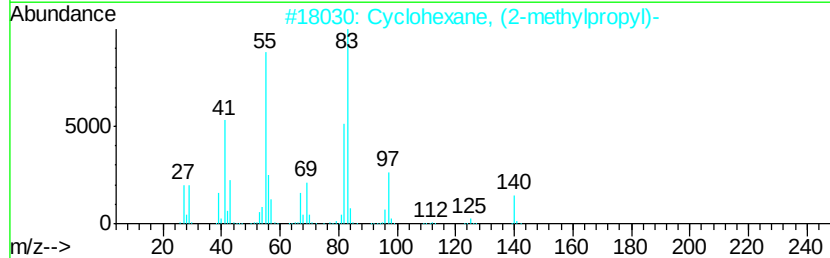
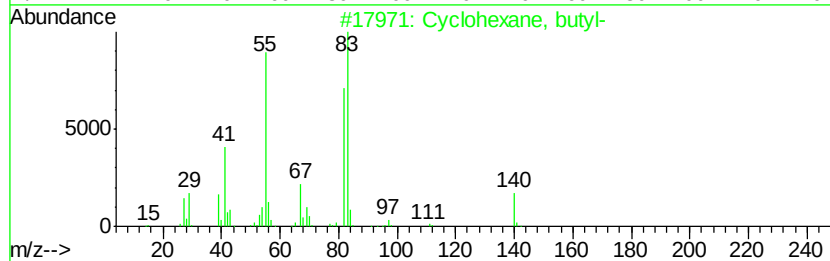
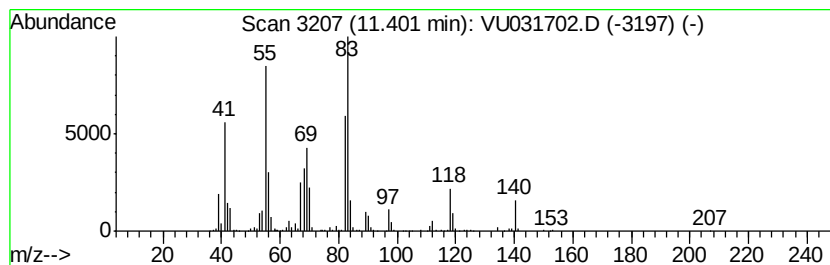
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

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Peak Number 13 (DEL) Alkane: Cyclic11.40 Concentration Rank 34

| R.T.      | EstConc                        | Area   | Relative to ISTD       | R.T.        |      |
|-----------|--------------------------------|--------|------------------------|-------------|------|
| 11.40     | 12.86 ug/L                     | 733369 | 1,4-Dichlorobenzene-d4 | 11.48       |      |
| Hit# of 5 | Tentative ID                   | MW     | MolForm                | CAS#        | Qual |
| 1         | Cyclohexane, butyl-            | 140    | C10H20                 | 001678-93-9 | 80   |
| 2         | Cyclohexane, (2-methylpropyl)- | 140    | C10H20                 | 001678-98-4 | 72   |
| 3         | Cyclohexane, undecyl-          | 238    | C17H34                 | 054105-66-7 | 64   |
| 4         | Cyclohexane, butyl-            | 140    | C10H20                 | 001678-93-9 | 64   |
| 5         | Cyclohexane, butyl-            | 140    | C10H20                 | 001678-93-9 | 64   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

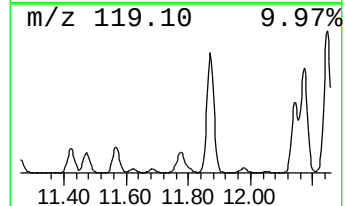
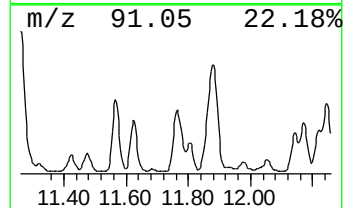
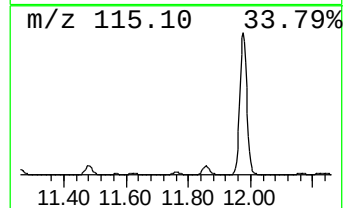
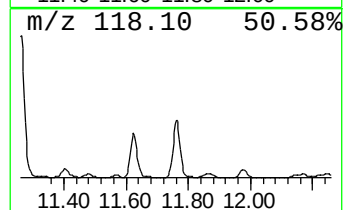
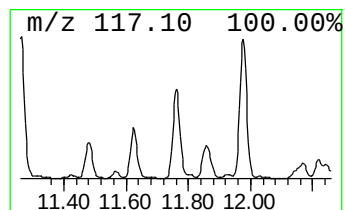
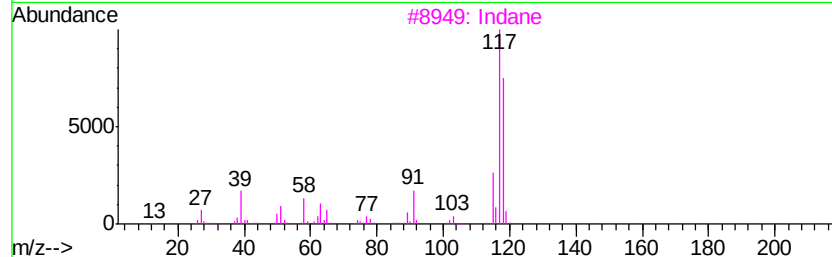
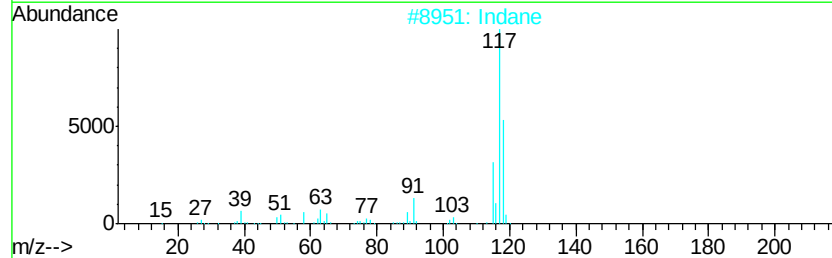
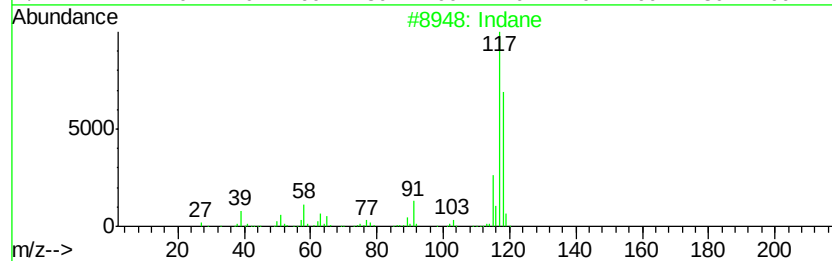
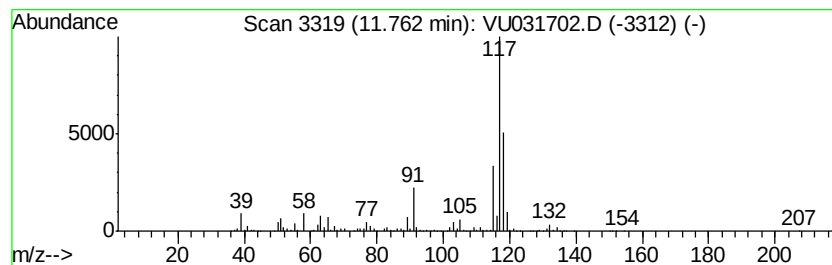
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

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Peak Number 15 Indane Concentration Rank 31

| R.T.    | EstConc                      | Area         | Relative to ISTD       | R.T.           |
|---------|------------------------------|--------------|------------------------|----------------|
| 11.76   | 16.97 ug/L                   | 967421       | 1,4-Dichlorobenzene-d4 | 11.48          |
| Hit# of | 5                            | Tentative ID | MW MolForm             | CAS# Qual      |
| 1       | Indane                       |              | 118 C9H10              | 000496-11-7 76 |
| 2       | Indane                       |              | 118 C9H10              | 000496-11-7 76 |
| 3       | Indane                       |              | 118 C9H10              | 000496-11-7 76 |
| 4       | Benzene, 1-ethenyl-2-methyl- |              | 118 C9H10              | 000611-15-4 70 |
| 5       | Benzene, cyclopropyl-        |              | 118 C9H10              | 000873-49-4 70 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

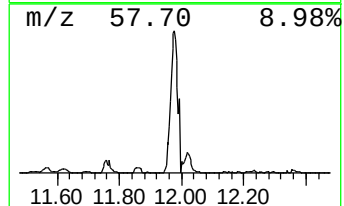
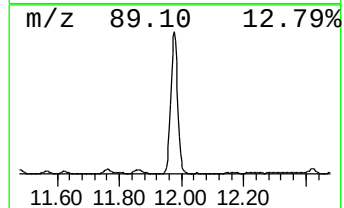
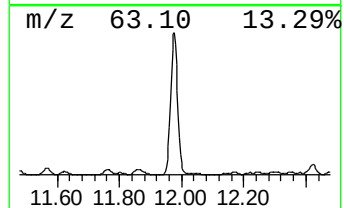
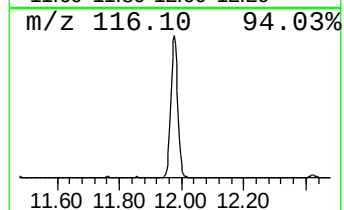
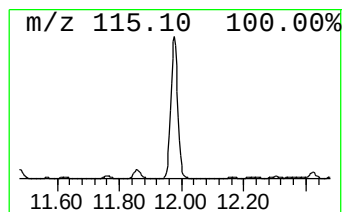
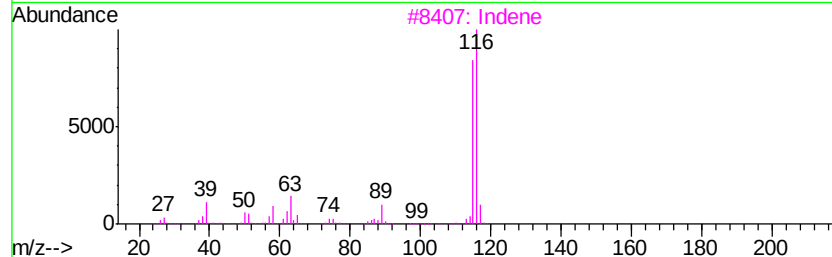
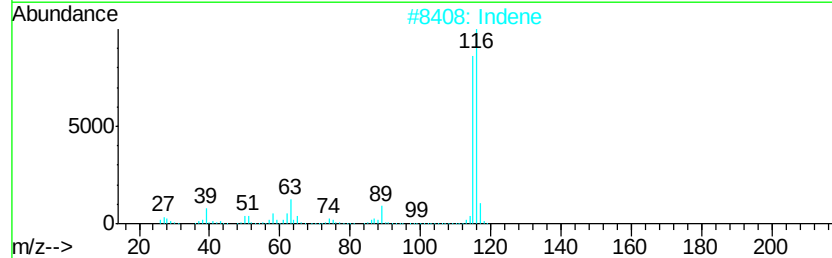
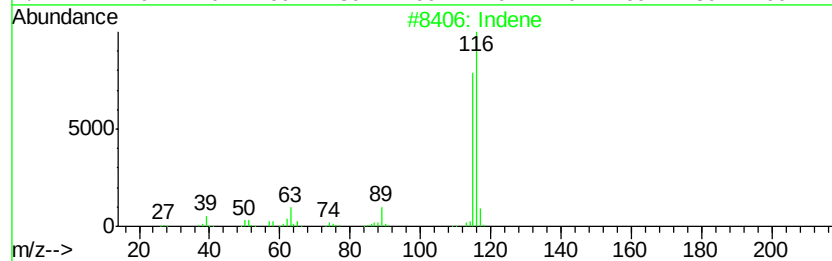
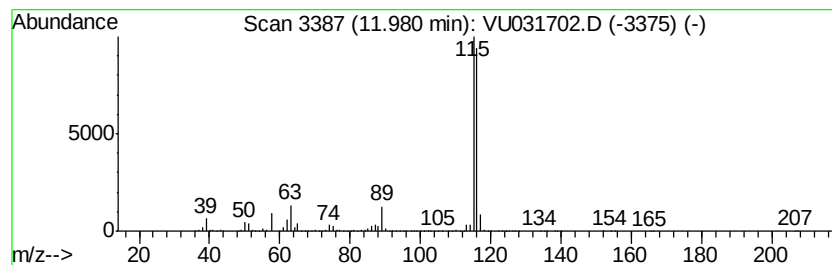
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Indene Concentration Rank 6

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 11.98 | 247.01 ug/L | 14082700 | 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 | 95   |
| 3    |    | Indene                  | 116 | C9H8    | 000095-13-6 | 94   |
| 4    |    | 3-Methylphenylacetylene | 116 | C9H8    | 000766-82-5 | 91   |
| 5    |    | Benzene, 1-propynyl-    | 116 | C9H8    | 000673-32-5 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
 Data File : VU031702.D  
 Acq On : 02 May 2019 11:15  
 Operator : JC/SP  
 Sample : K2603-16DL 10X  
 Misc : 5.10µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 GAJ68DL

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

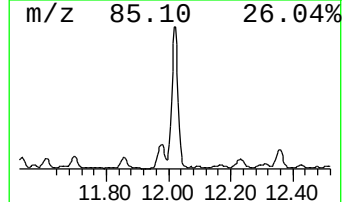
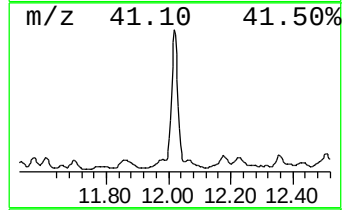
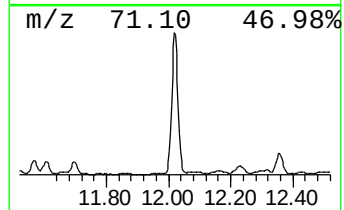
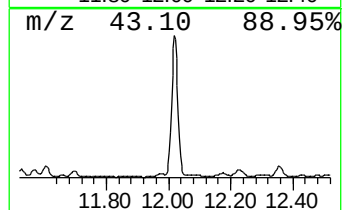
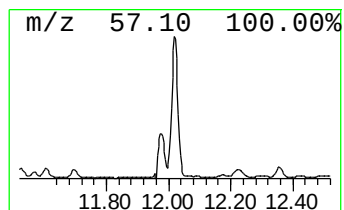
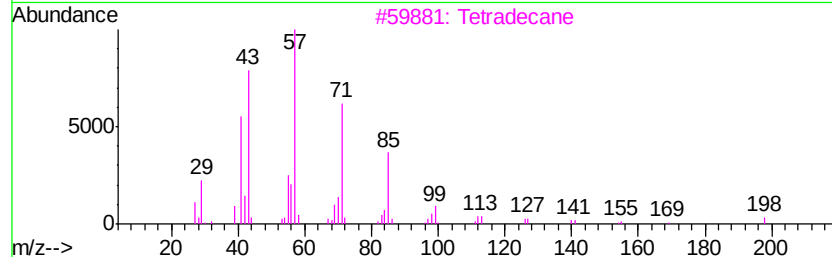
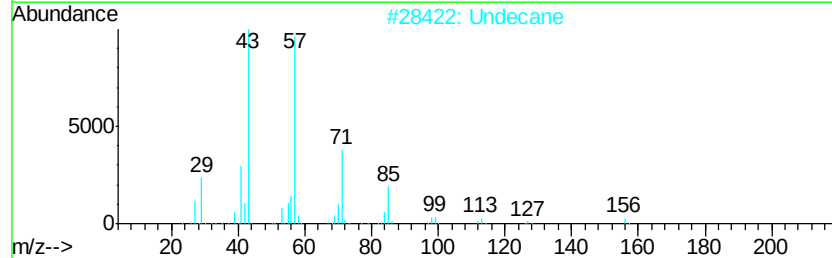
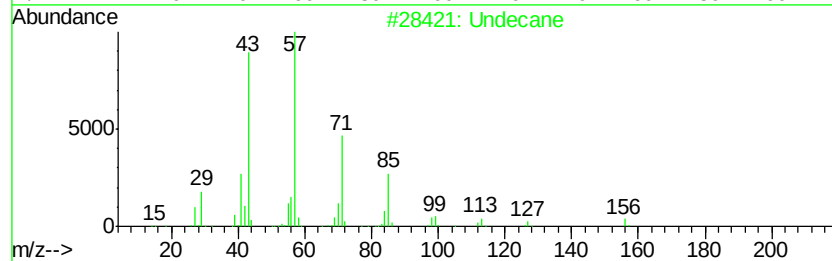
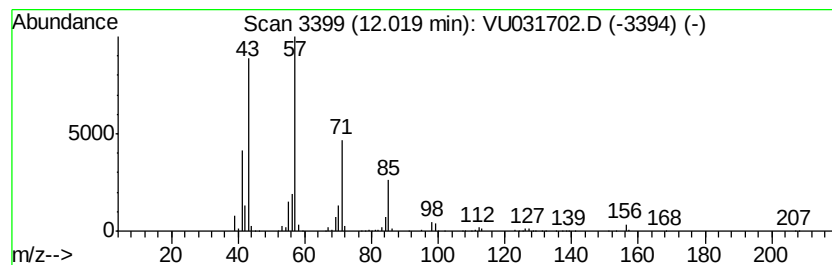
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.02 | 80.95 ug/L | 4615070 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Undecane     | 156 | C11H24  | 001120-21-4 | 91   |
| 2    |    | Undecane     | 156 | C11H24  | 001120-21-4 | 90   |
| 3    |    | Tetradecane  | 198 | C14H30  | 000629-59-4 | 90   |
| 4    |    | Undecane     | 156 | C11H24  | 001120-21-4 | 87   |
| 5    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
GAJ68DL

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

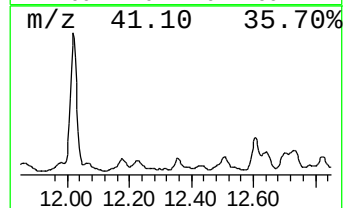
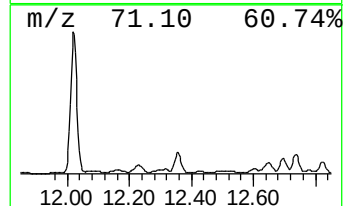
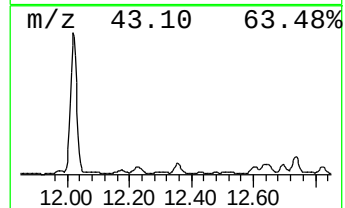
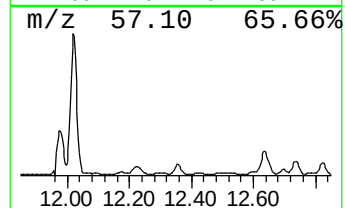
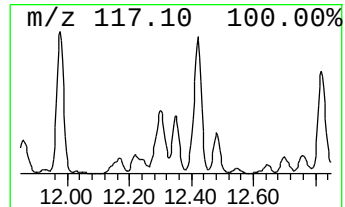
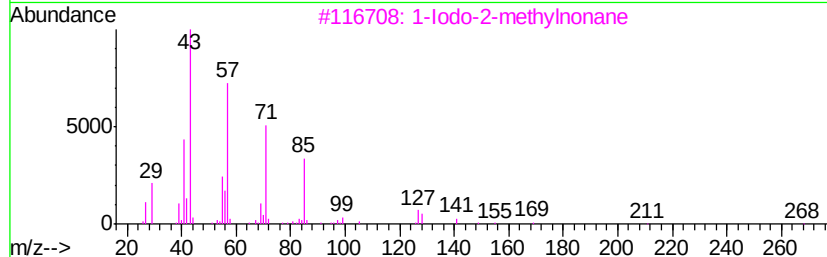
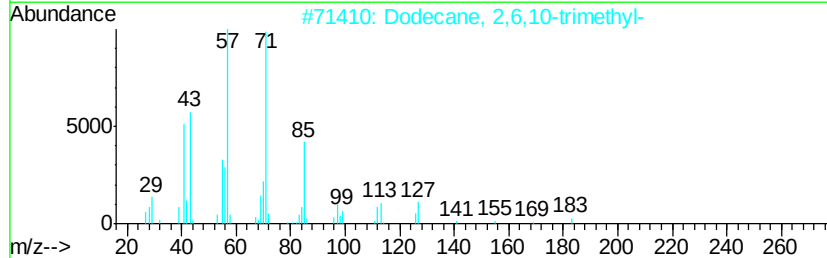
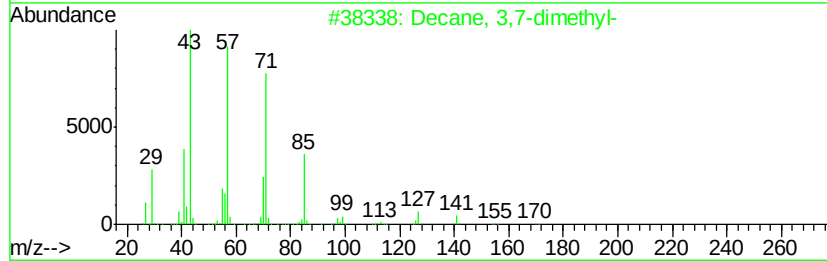
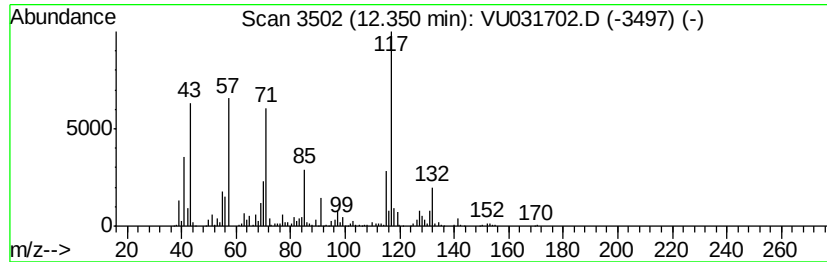
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 12.35 | 13.98 ug/L | 797191 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#         | Qual |
|------|----|-----------------------------|-----|---------|--------------|------|
| 1    | 5  | Decane, 3,7-dimethyl-       | 170 | C12H26  | 017312-54-8  | 50   |
| 2    |    | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3  | 43   |
| 3    |    | 1-Iodo-2-methylnonane       | 268 | C10H21I | 1000101-47-9 | 38   |
| 4    |    | Undecane, 4-methyl-         | 170 | C12H26  | 002980-69-0  | 38   |
| 5    |    | Indan, 1-methyl-            | 132 | C10H12  | 000767-58-8  | 38   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

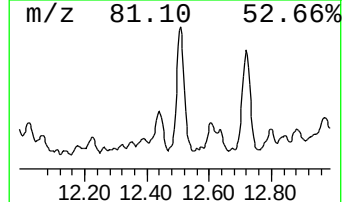
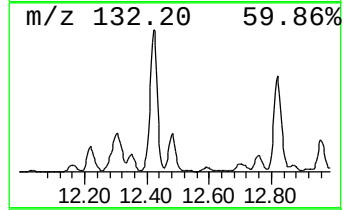
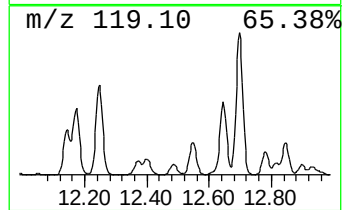
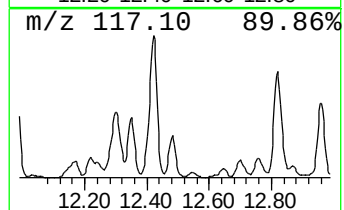
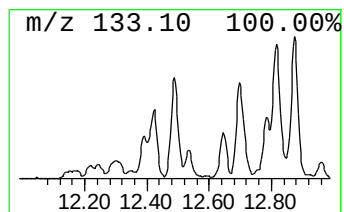
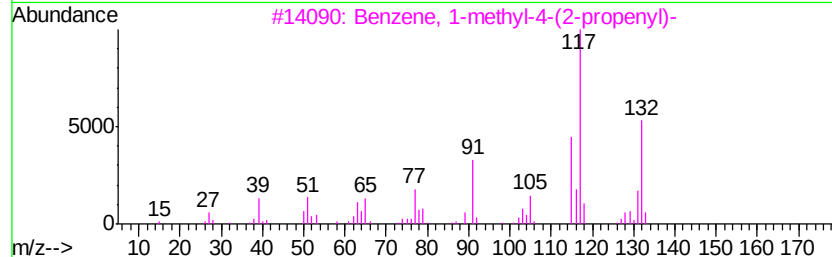
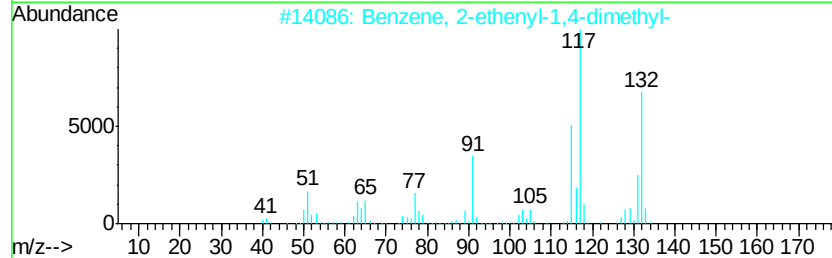
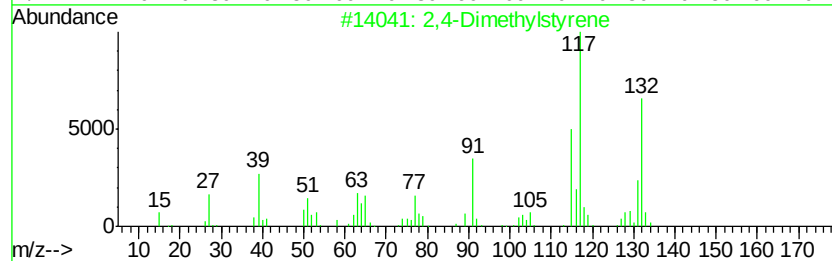
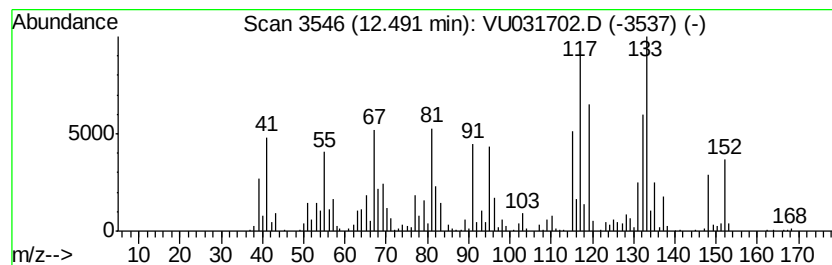
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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 2,4-Dimethylstyrene Concentration Rank 35

| R.T.      | EstConc                           | Area   | Relative to ISTD       | R.T.        |      |
|-----------|-----------------------------------|--------|------------------------|-------------|------|
| 12.49     | 11.59 ug/L                        | 660567 | 1,4-Dichlorobenzene-d4 | 11.48       |      |
| Hit# of 5 | Tentative ID                      | MW     | MolForm                | CAS#        | Qual |
| 1         | 2,4-Dimethylstyrene               | 132    | C10H12                 | 002234-20-0 | 97   |
| 2         | Benzene, 2-ethenyl-1,4-dimethyl-  | 132    | C10H12                 | 002039-89-6 | 96   |
| 3         | Benzene, 1-methyl-4-(2-propenyl)- | 132    | C10H12                 | 003333-13-9 | 93   |
| 4         | Benzene, (2-methyl-1-propenyl)-   | 132    | C10H12                 | 000768-49-0 | 92   |
| 5         | Benzene, 4-ethenyl-1,2-dimethyl-  | 132    | C10H12                 | 027831-13-6 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

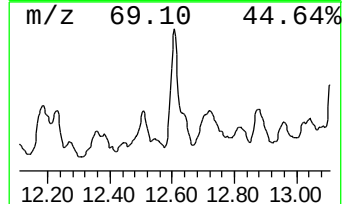
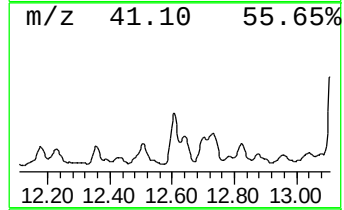
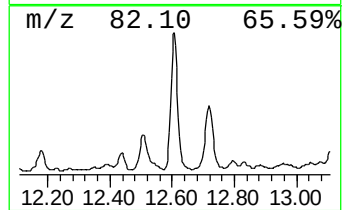
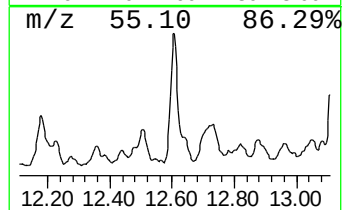
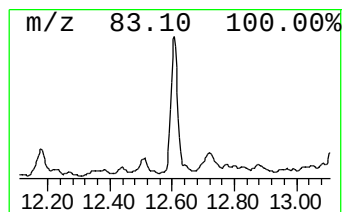
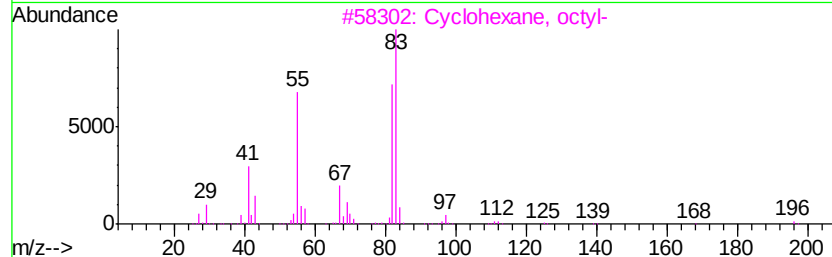
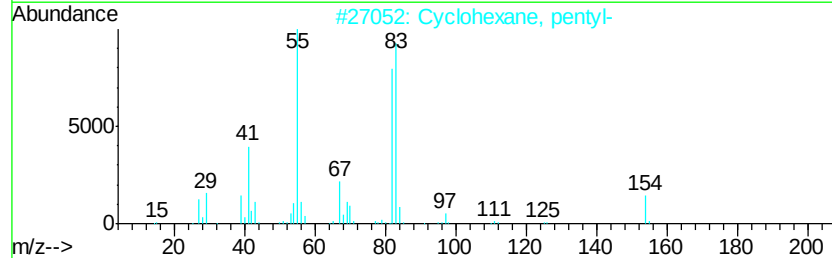
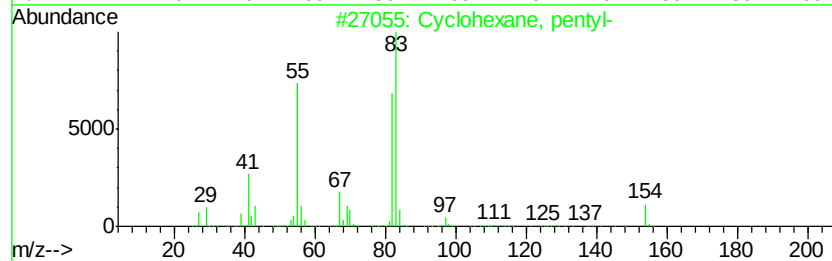
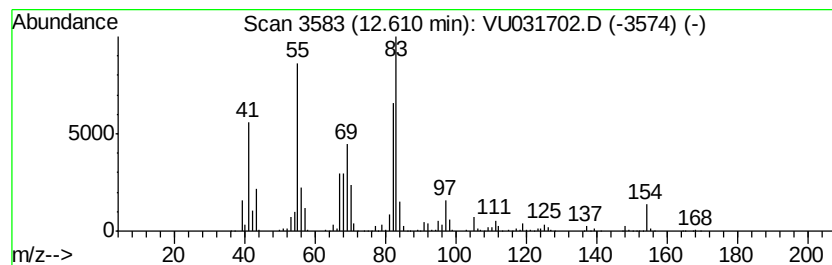
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 (DEL) Alkane: Cyclic12.61 Concentration Rank 26

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.61 | 26.40 ug/L | 1505340 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 64   |
| 2    |    | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 64   |
| 3    |    | Cyclohexane, octyl-            | 196 | C14H28  | 001795-15-9 | 59   |
| 4    |    | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 58   |
| 5    |    | Cyclohexane, (4-methylpentyl)- | 168 | C12H24  | 061142-20-9 | 53   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

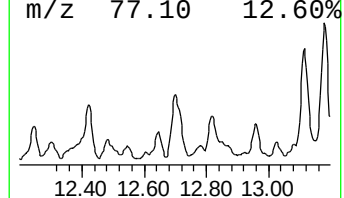
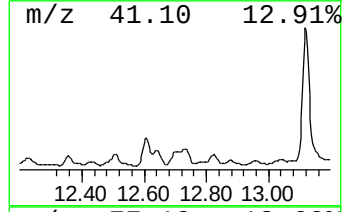
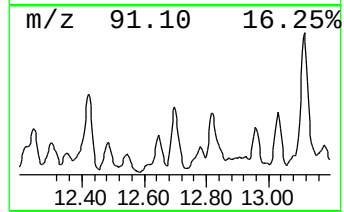
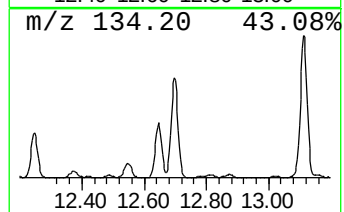
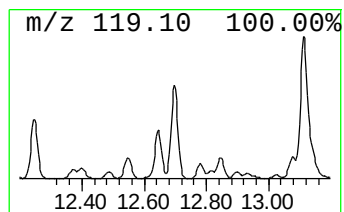
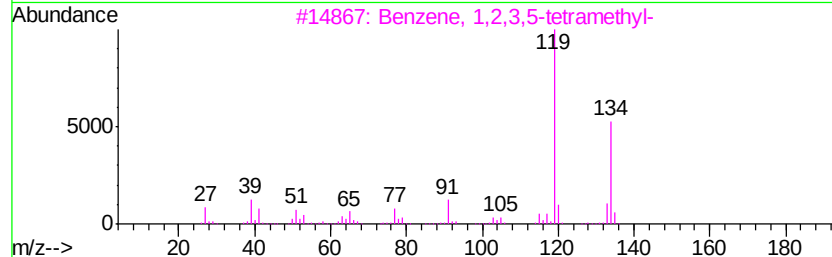
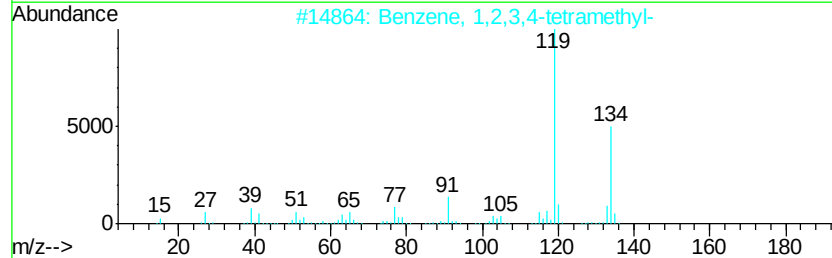
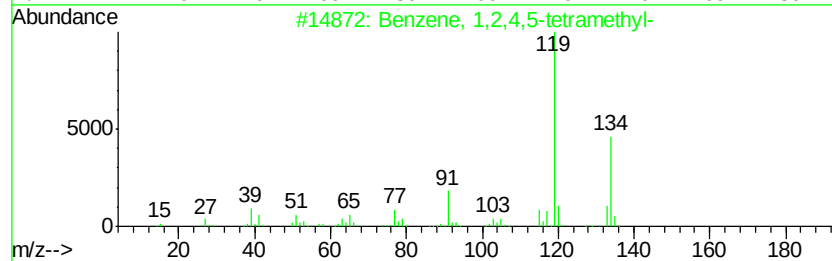
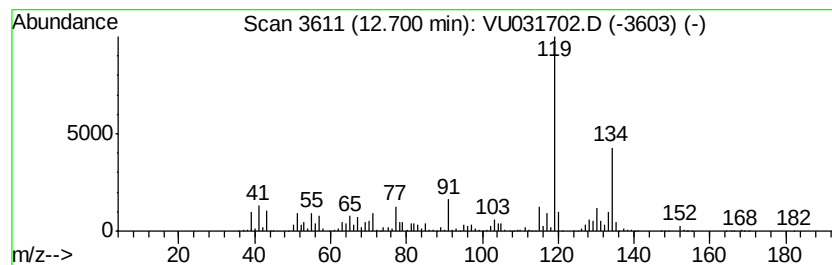
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 19

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.70 | 65.82 ug/L | 3752490 | 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 | 93   |
| 3    |    |   | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 93   |
| 4    |    |   | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 93   |
| 5    |    |   | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

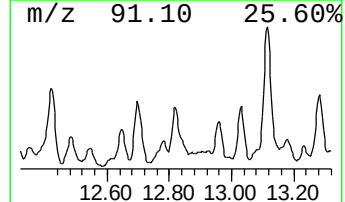
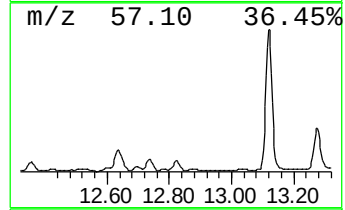
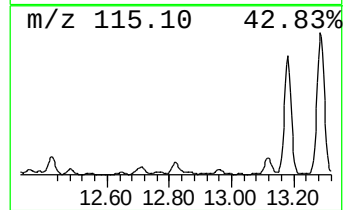
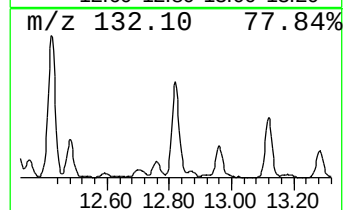
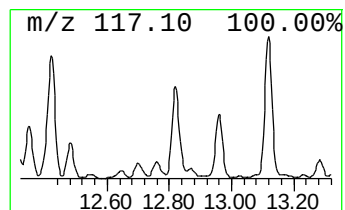
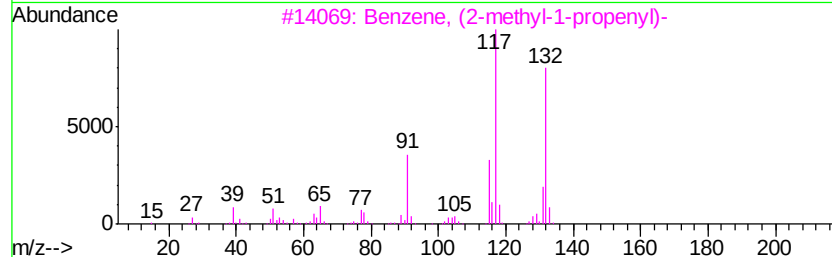
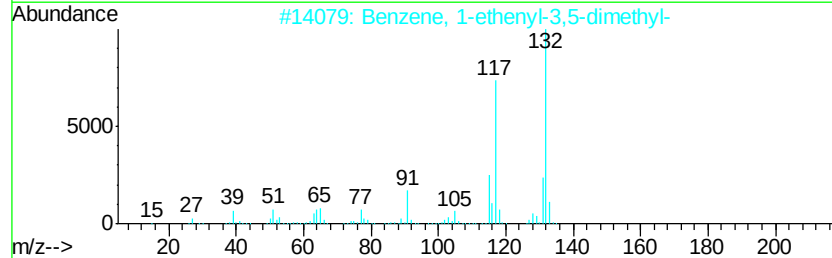
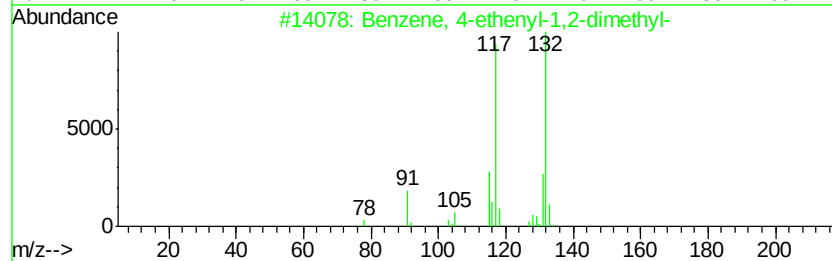
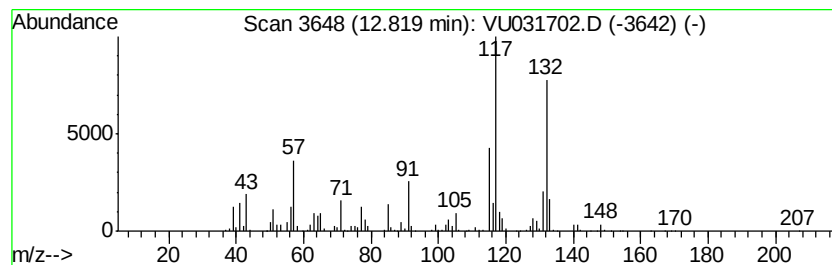
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 Benzene, 4-ethenyl-1,2-dime... Concentration Rank 28

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.82 | 24.93 ug/L | 1421600 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-ethenyl-1,2-dimethyl-    | 132 | C10H12  | 027831-13-6 | 97   |
| 2    |    | Benzene, 1-ethenyl-3,5-dimethyl-    | 132 | C10H12  | 005379-20-4 | 96   |
| 3    |    | Benzene, (2-methyl-1-propenyl)-     | 132 | C10H12  | 000768-49-0 | 94   |
| 4    |    | Benzene, 2-butenyl-                 | 132 | C10H12  | 001560-06-1 | 93   |
| 5    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 93   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

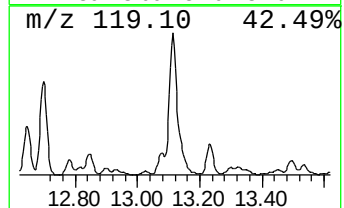
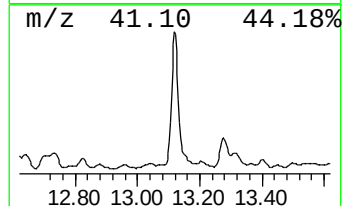
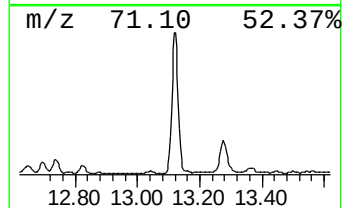
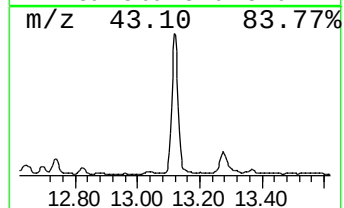
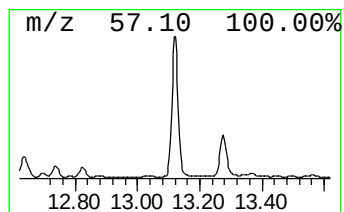
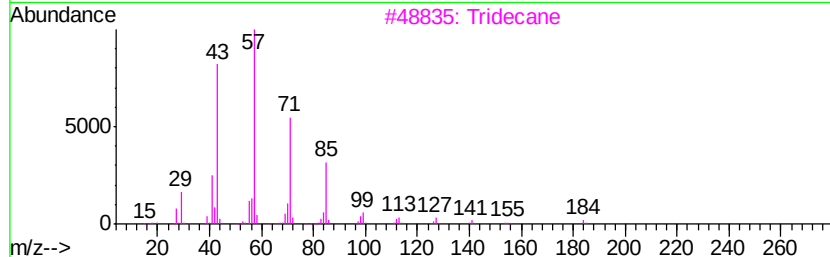
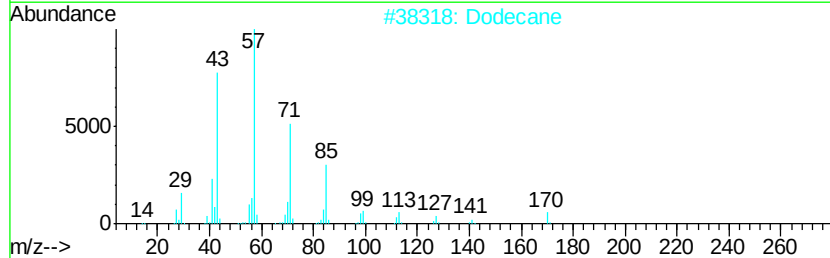
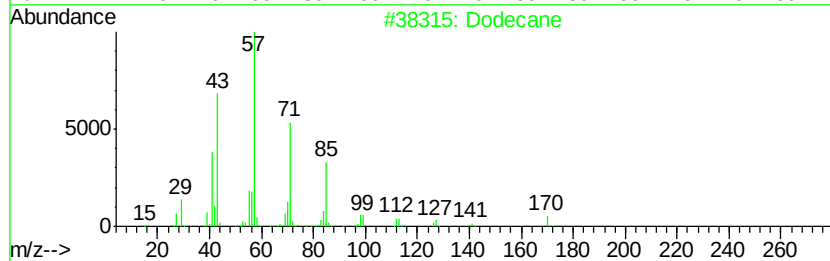
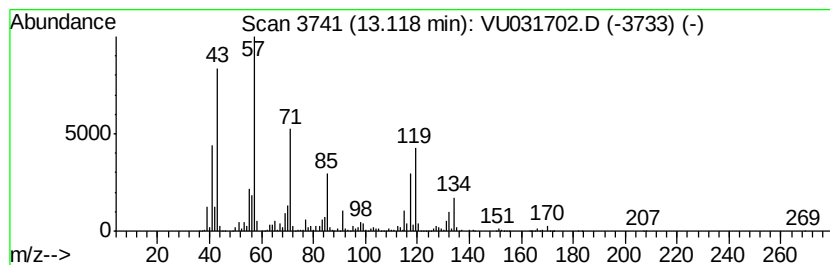
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 13.12 | 161.48 ug/L | 9206370 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane     | 170 | C12H26  | 000112-40-3 | 86   |
| 2    |    | Dodecane     | 170 | C12H26  | 000112-40-3 | 46   |
| 3    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 38   |
| 4    |    | Undecane     | 156 | C11H24  | 001120-21-4 | 38   |
| 5    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 38   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

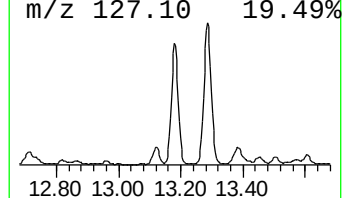
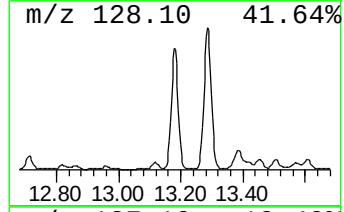
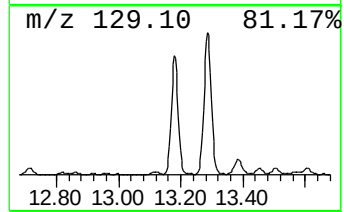
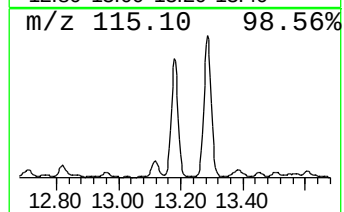
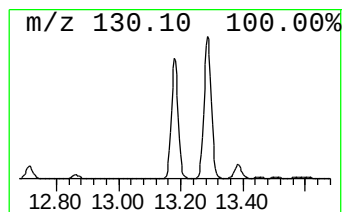
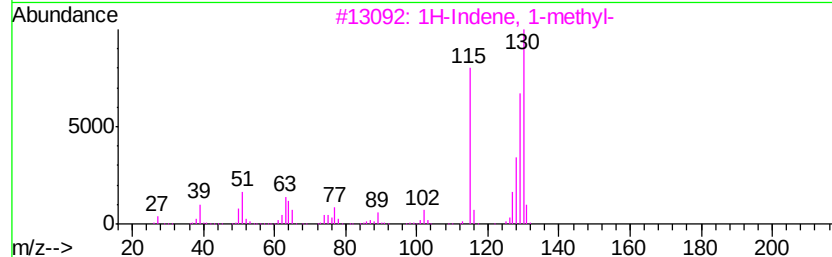
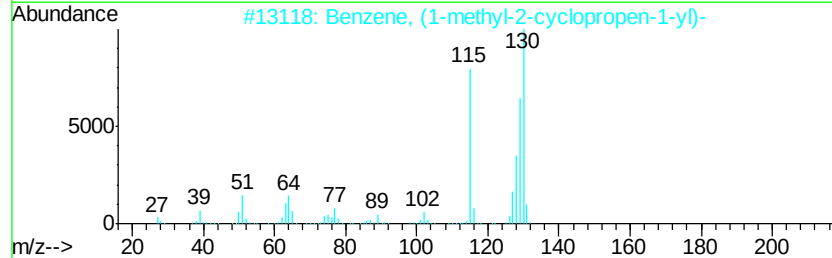
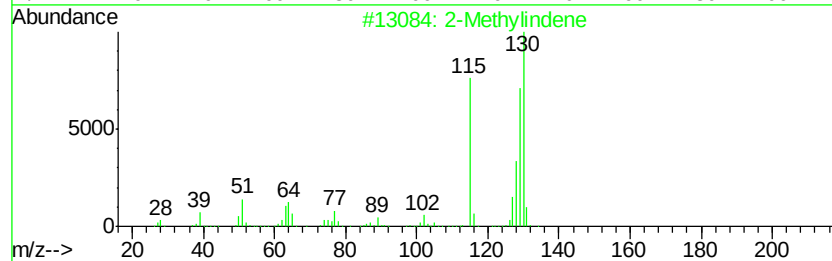
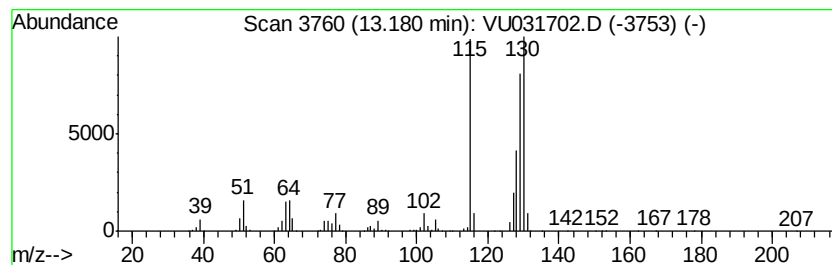
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 2-Methylindene Concentration Rank 11

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 13.18 | 113.24 ug/L | 6456040 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Methylindene                      | 130 | C10H10  | 002177-47-1 | 97   |
| 2    |    | Benzene, (1-methyl-2-cyclopropen... | 130 | C10H10  | 065051-83-4 | 96   |
| 3    |    | 1H-Indene, 1-methyl-                | 130 | C10H10  | 000767-59-9 | 96   |
| 4    |    | Benzene,1-methyl-1,2-propadienyl-   | 130 | C10H10  | 022433-39-2 | 95   |
| 5    |    | 1H-Indene, 3-methyl-                | 130 | C10H10  | 000767-60-2 | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

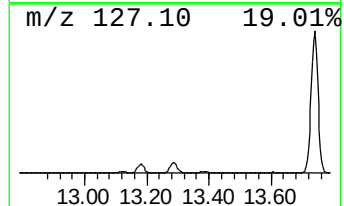
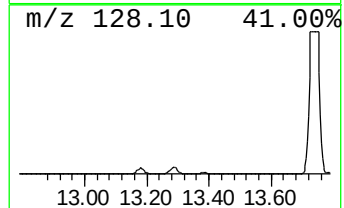
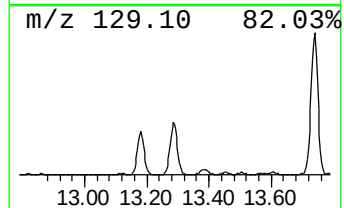
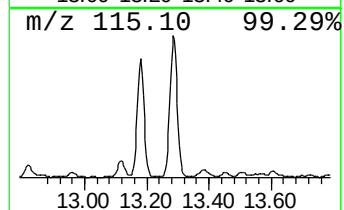
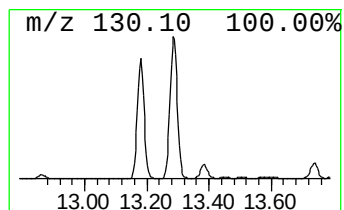
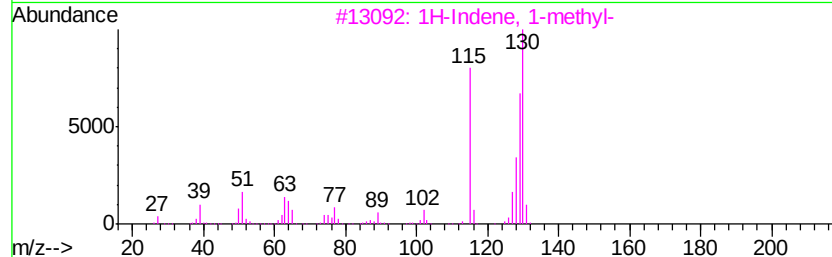
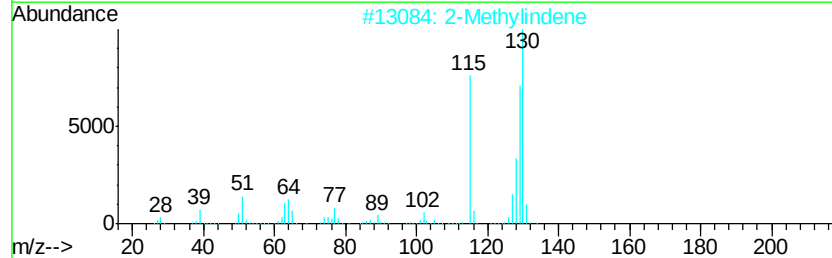
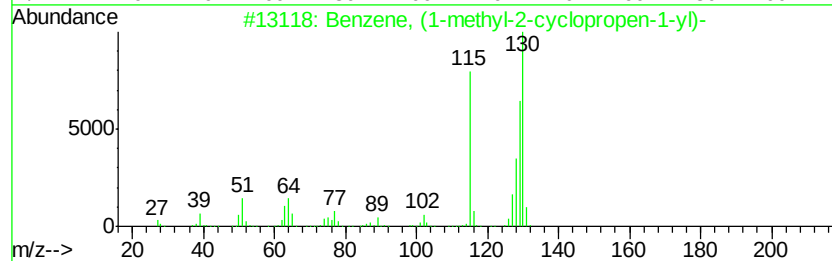
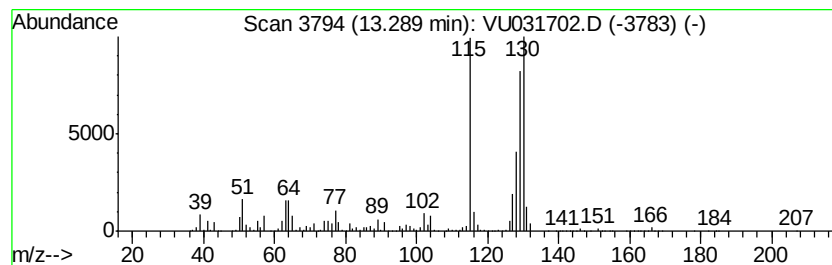
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 25 Benzene, (1-methyl-2-cyclop... Concentration Rank 7

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 13.29 | 196.71 ug/L | 11214700 | 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 | 96   |
| 2    |    | 2-Methylindene                      | 130 | C10H10  | 002177-47-1 | 96   |
| 3    |    | 1H-Indene, 1-methyl-                | 130 | C10H10  | 000767-59-9 | 96   |
| 4    |    | 1H-Indene, 1-methyl-                | 130 | C10H10  | 000767-59-9 | 96   |
| 5    |    | Benzene, 1-butylnyl-                | 130 | C10H10  | 000622-76-4 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

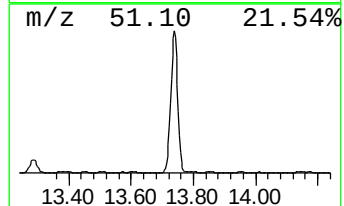
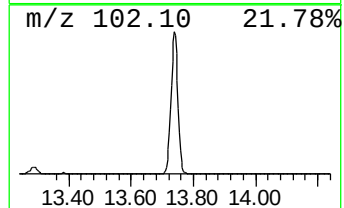
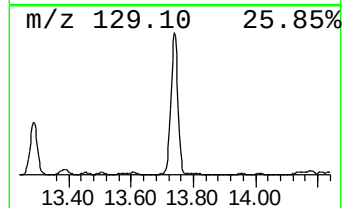
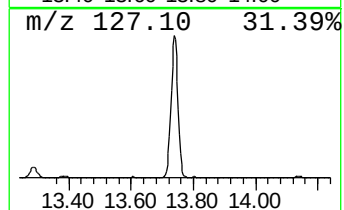
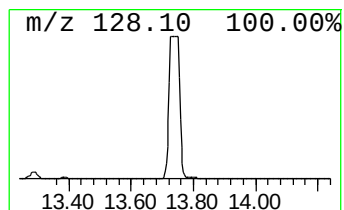
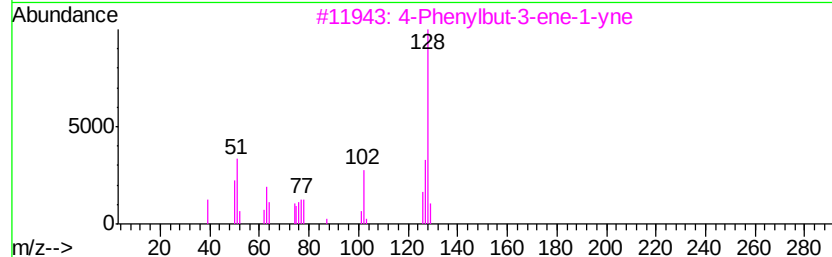
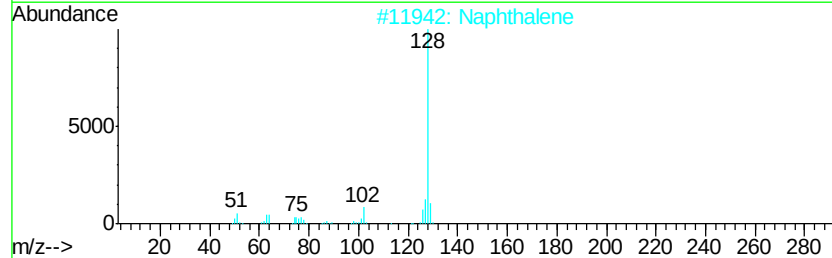
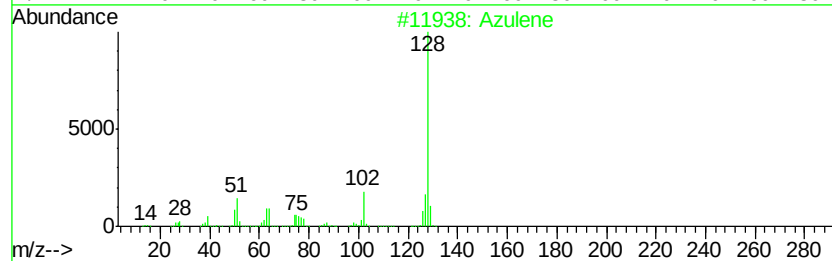
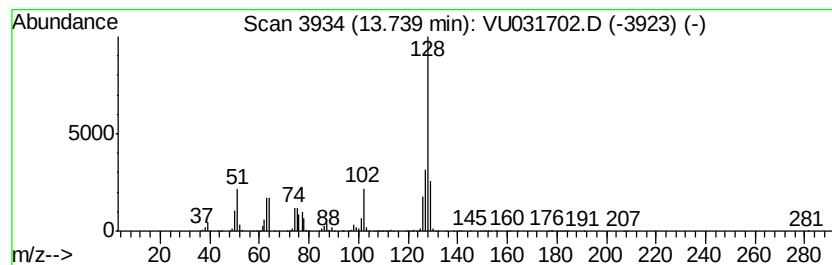
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 Azulene Concentration Rank 2

| R.T.  | EstConc      | Area     | Relative to ISTD       | R.T.  |
|-------|--------------|----------|------------------------|-------|
| 13.74 | 1073.09 ug/L | 61179700 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------|-----|---------|--------------|------|
| 1    | 5  | Azulene                 | 128 | C10H8   | 000275-51-4  | 87   |
| 2    |    | Naphthalene             | 128 | C10H8   | 000091-20-3  | 87   |
| 3    |    | 4-Phenylbut-3-ene-1-yne | 128 | C10H8   | 1000222-12-0 | 86   |
| 4    |    | 1H-Indene, 1-methylene- | 128 | C10H8   | 002471-84-3  | 83   |
| 5    |    | Azulene                 | 128 | C10H8   | 000275-51-4  | 81   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

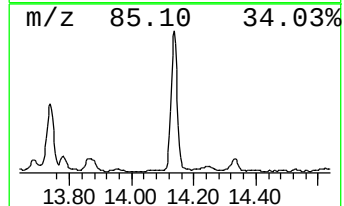
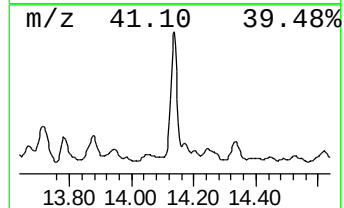
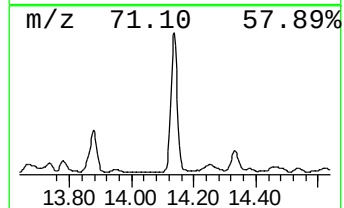
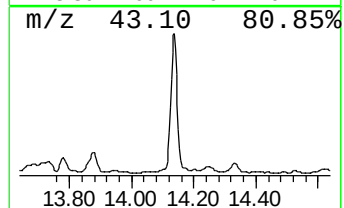
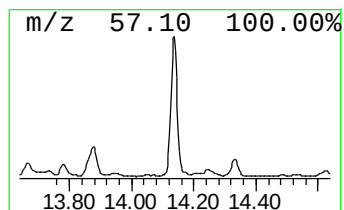
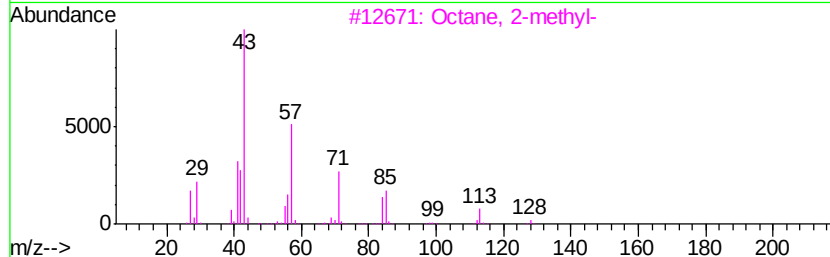
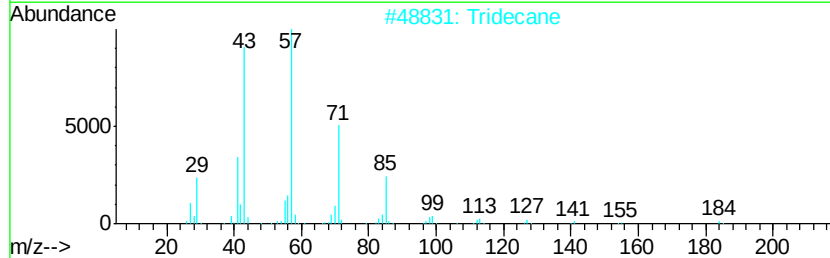
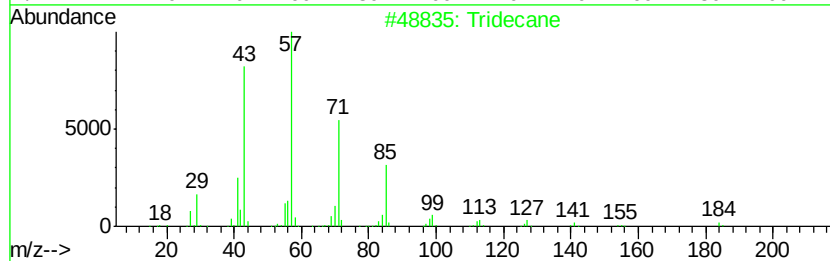
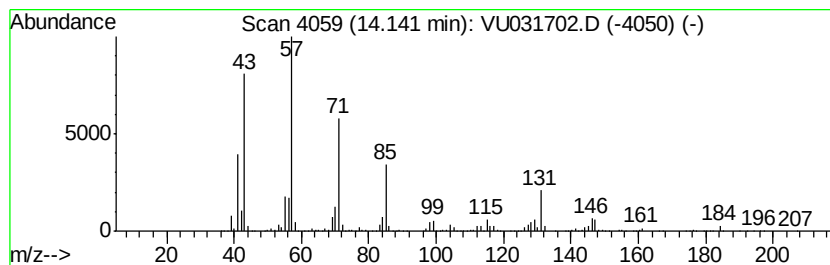
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 14.14 | 80.94 ug/L | 4614800 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane         | 184 | C13H28  | 000629-50-5 | 94   |
| 2    |    | Tridecane         | 184 | C13H28  | 000629-50-5 | 94   |
| 3    |    | Octane, 2-methyl- | 128 | C9H20   | 003221-61-2 | 58   |
| 4    |    | Tetradecane       | 198 | C14H30  | 000629-59-4 | 53   |
| 5    |    | Heptadecane       | 240 | C17H36  | 000629-78-7 | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

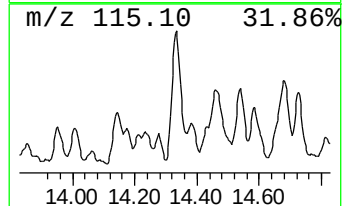
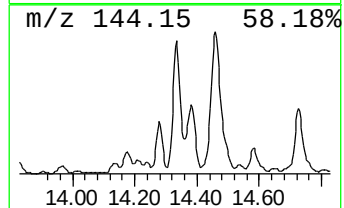
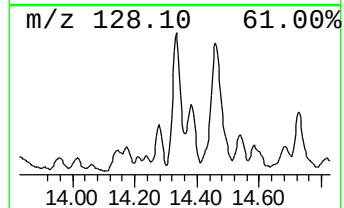
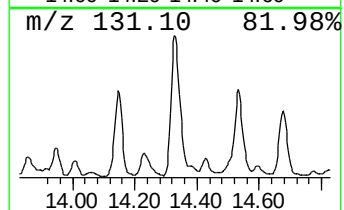
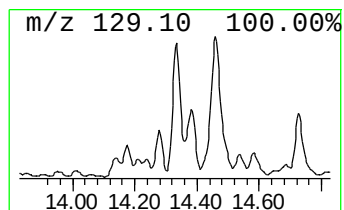
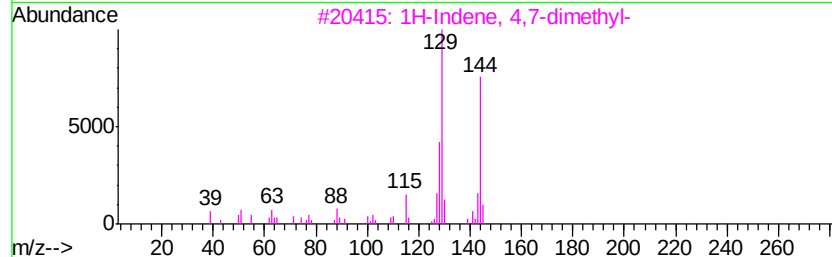
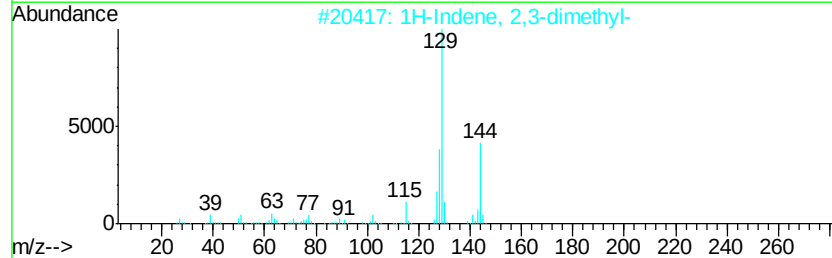
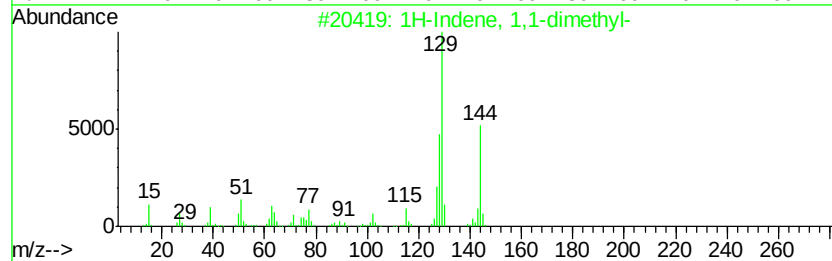
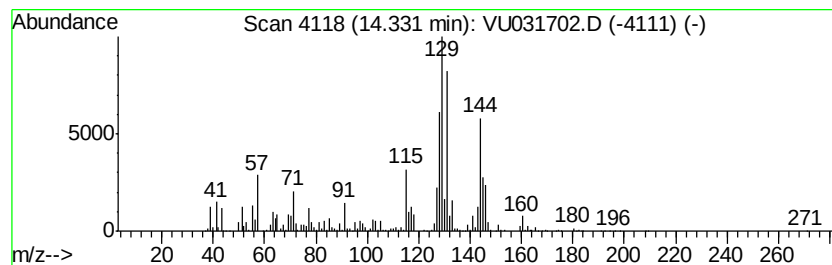
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 28 1H-Indene, 1,1-dimethyl- Concentration Rank 20

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 14.33 | 64.50 ug/L | 3677100 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 1,1-dimethyl-       | 144 | C11H12  | 018636-55-0 | 91   |
| 2    |    |   | 1H-Indene, 2,3-dimethyl-       | 144 | C11H12  | 004773-82-4 | 60   |
| 3    |    |   | 1H-Indene, 4,7-dimethyl-       | 144 | C11H12  | 006974-97-6 | 60   |
| 4    |    |   | 1H-Indene, 1,3-dimethyl-       | 144 | C11H12  | 002177-48-2 | 60   |
| 5    |    |   | Benzene, (1-methyl-2-butynyl)- | 144 | C11H12  | 087712-69-4 | 55   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

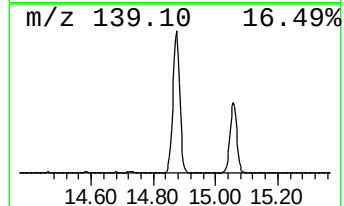
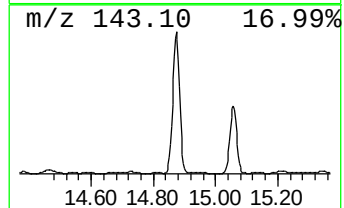
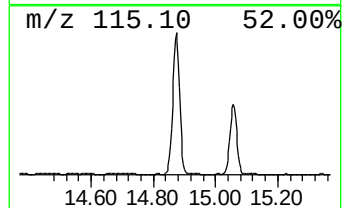
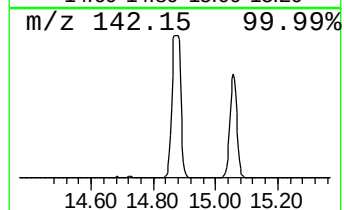
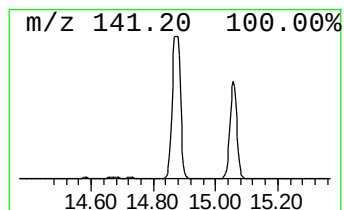
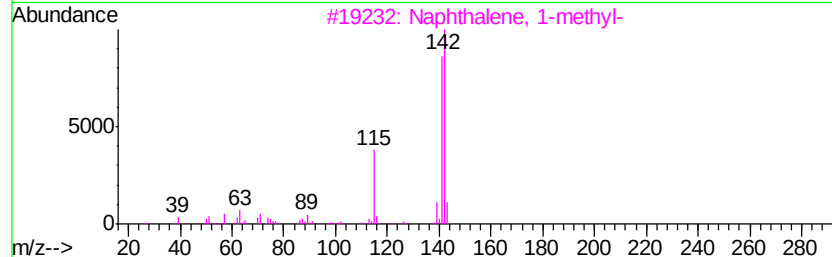
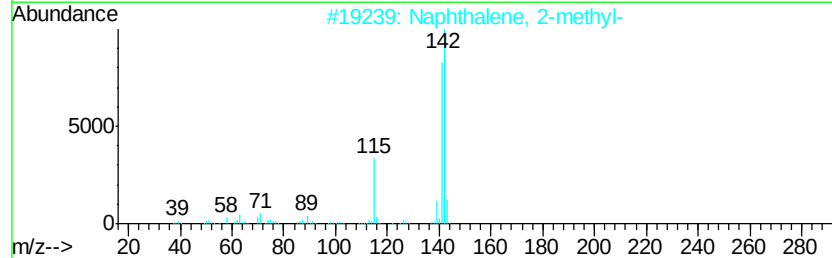
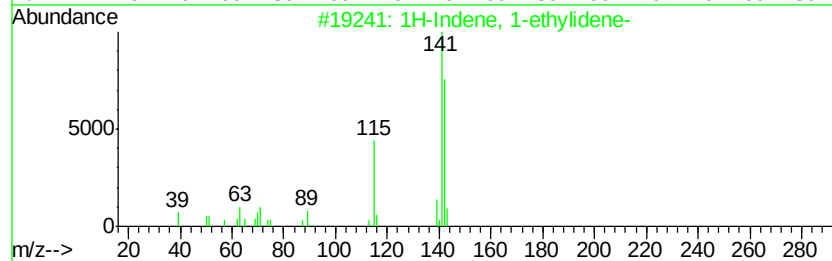
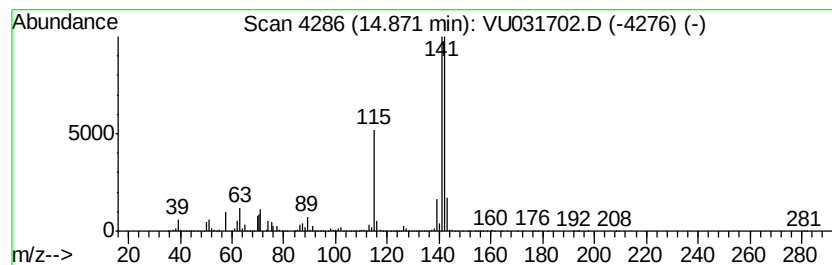
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 1H-Indene, 1-ethylidene- Concentration Rank 1

| R.T.  | EstConc      | Area     | Relative to ISTD       | R.T.  |
|-------|--------------|----------|------------------------|-------|
| 14.87 | 1209.13 ug/L | 68935800 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 1-ethylidene- | 142 | C11H10  | 002471-83-2 | 97   |
| 2    |    | Naphthalene, 2-methyl-   | 142 | C11H10  | 000091-57-6 | 95   |
| 3    |    | Naphthalene, 1-methyl-   | 142 | C11H10  | 000090-12-0 | 94   |
| 4    |    | Naphthalene, 1-methyl-   | 142 | C11H10  | 000090-12-0 | 94   |
| 5    |    | Naphthalene, 2-methyl-   | 142 | C11H10  | 000091-57-6 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

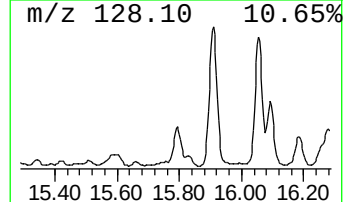
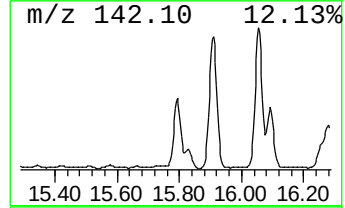
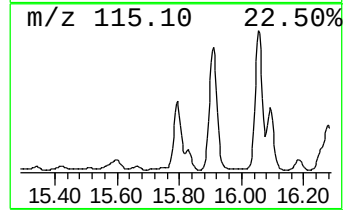
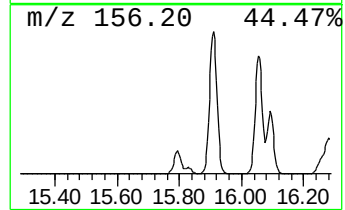
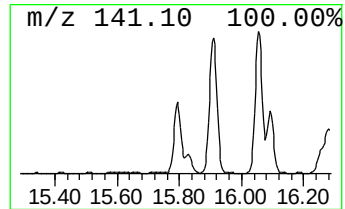
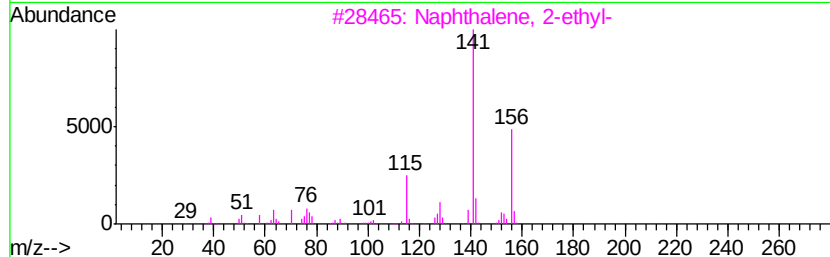
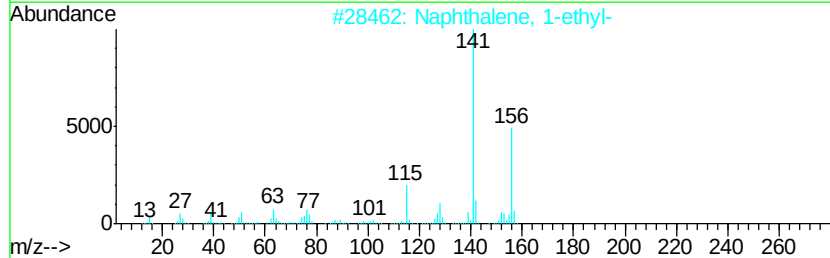
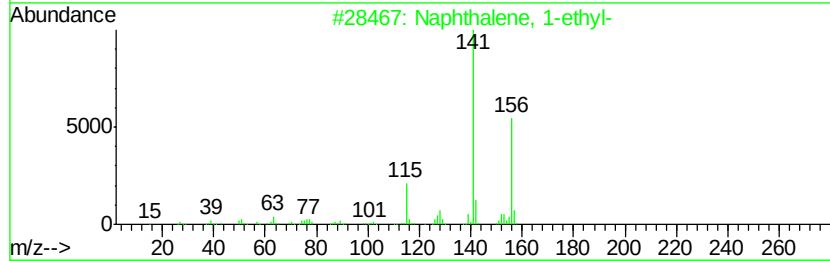
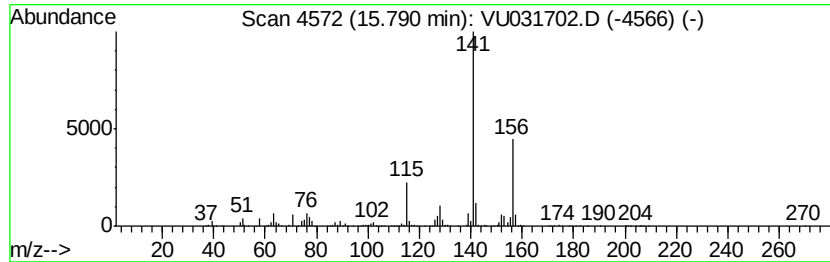
Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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 Naphthalene, 1-ethyl- Concentration Rank 13

| R.T.      | EstConc               | Area    | Relative to ISTD       | R.T.        |      |
|-----------|-----------------------|---------|------------------------|-------------|------|
| 15.79     | 82.01 ug/L            | 4675620 | 1,4-Dichlorobenzene-d4 | 11.48       |      |
| Hit# of 5 | Tentative ID          | MW      | MolForm                | CAS#        | Qual |
| 1         | Naphthalene, 1-ethyl- | 156     | C12H12                 | 001127-76-0 | 97   |
| 2         | Naphthalene, 1-ethyl- | 156     | C12H12                 | 001127-76-0 | 96   |
| 3         | Naphthalene, 2-ethyl- | 156     | C12H12                 | 000939-27-5 | 96   |
| 4         | Naphthalene, 2-ethyl- | 156     | C12H12                 | 000939-27-5 | 95   |
| 5         | Naphthalene, 1-ethyl- | 156     | C12H12                 | 001127-76-0 | 94   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

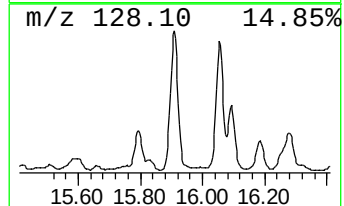
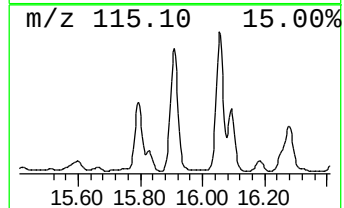
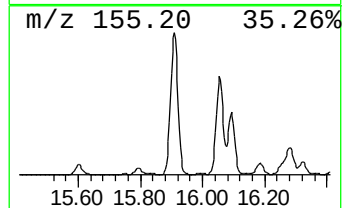
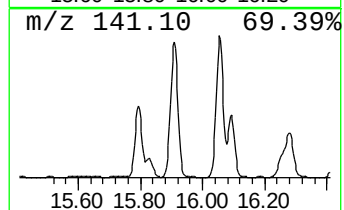
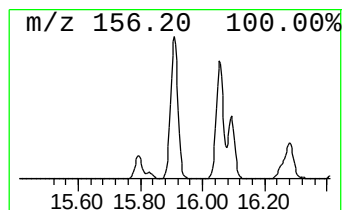
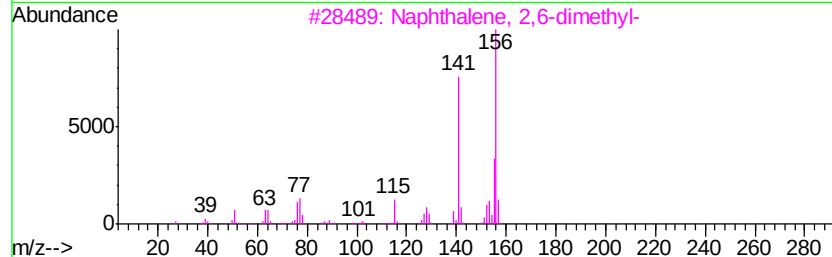
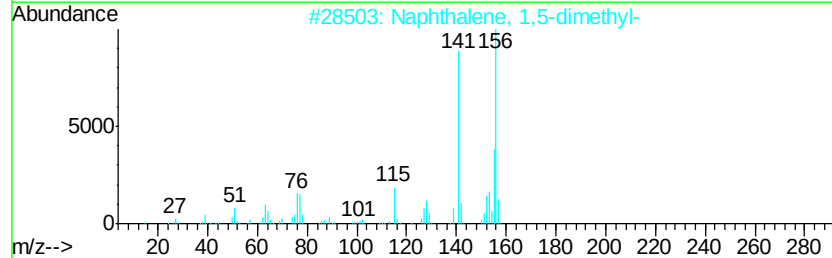
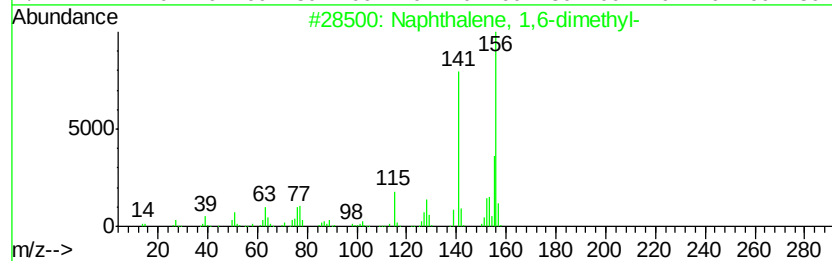
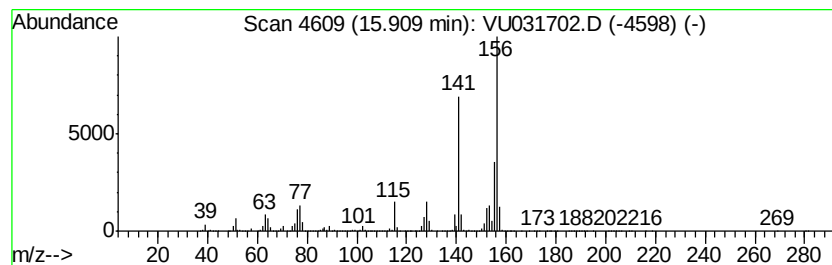
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 Naphthalene, 1,6-dimethyl- Concentration Rank 4

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 15.91 | 382.62 ug/L | 21814200 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 2    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 98   |
| 3    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 4    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 98   |
| 5    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

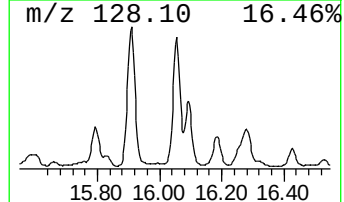
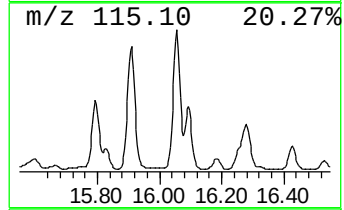
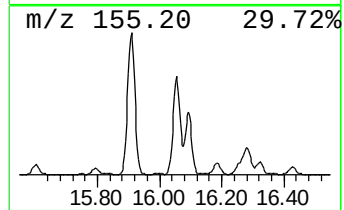
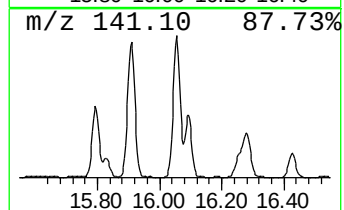
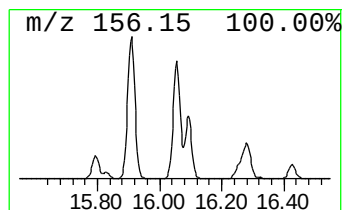
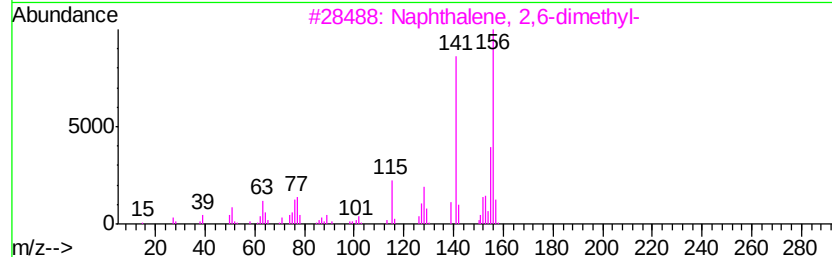
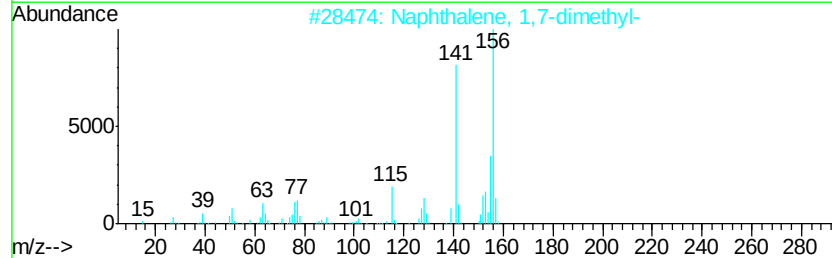
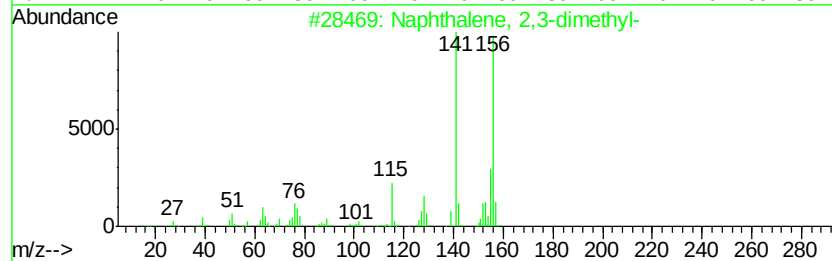
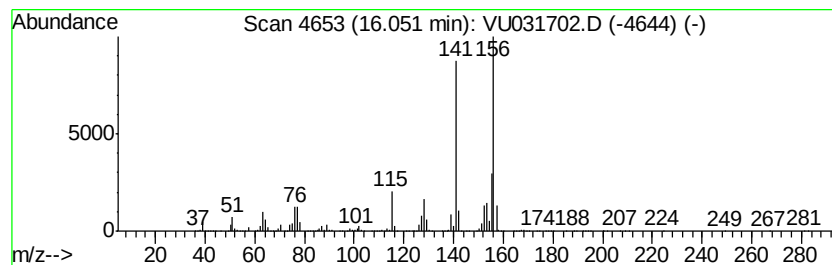
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 Naphthalene, 2,3-dimethyl- Concentration Rank 5

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 16.05 | 316.52 ug/L | 18045700 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 98   |
| 2    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 98   |
| 3    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 4    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 5    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

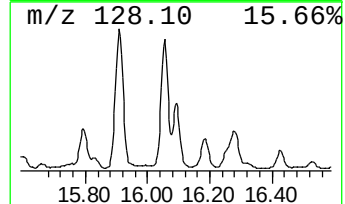
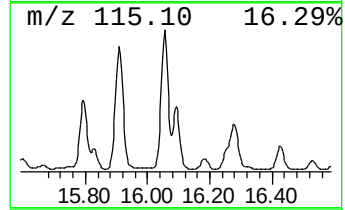
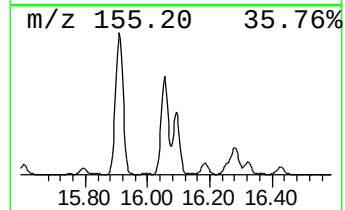
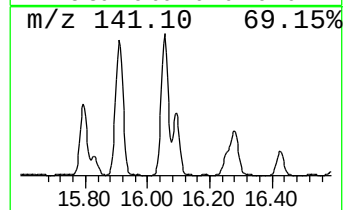
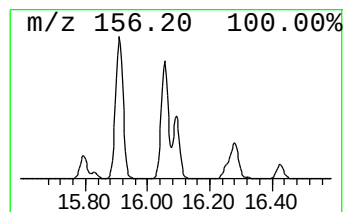
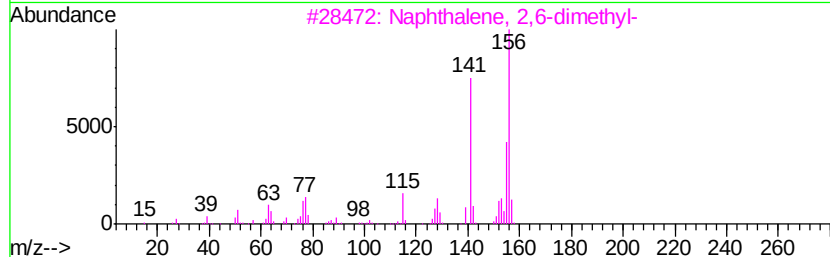
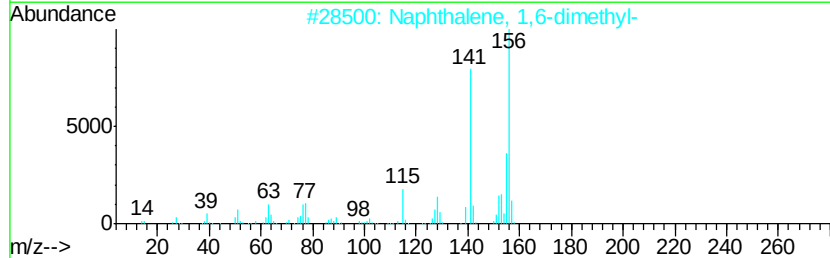
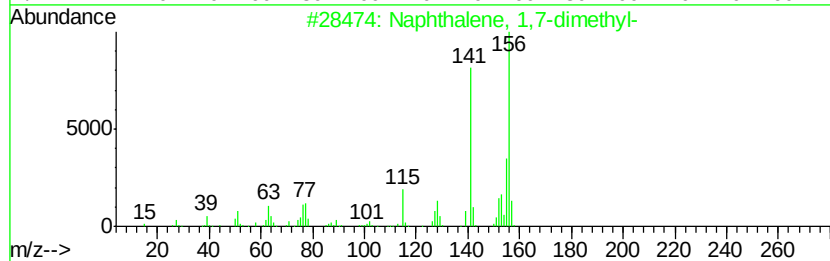
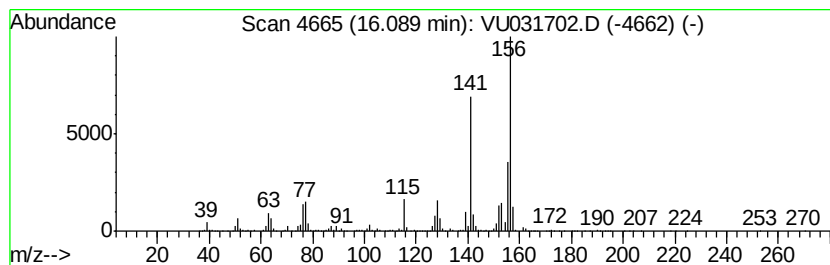
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 Naphthalene, 1,7-dimethyl- Concentration Rank 9

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 16.09 | 146.39 ug/L | 8346160 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 98   |
| 2    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 3    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 97   |
| 4    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 97   |
| 5    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

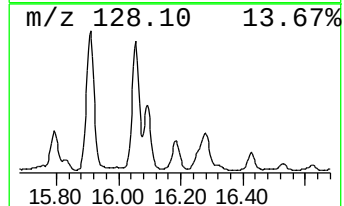
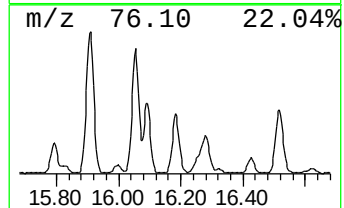
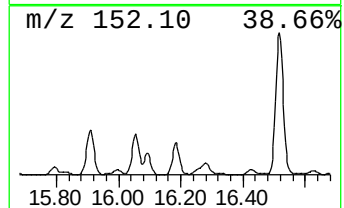
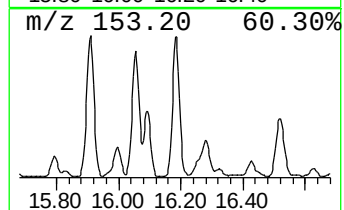
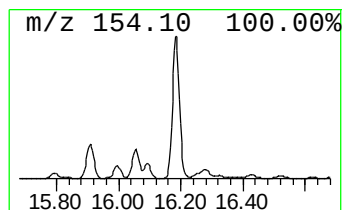
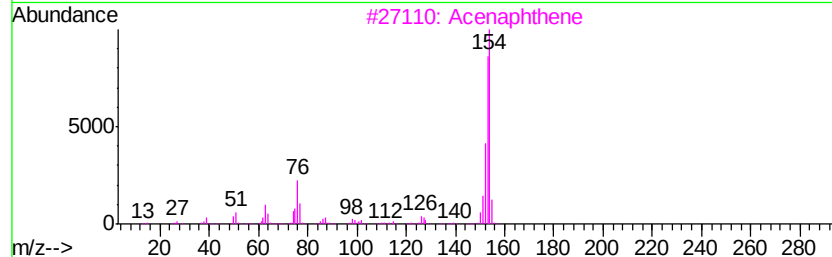
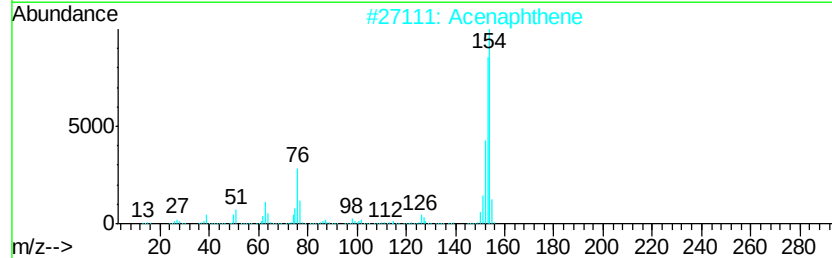
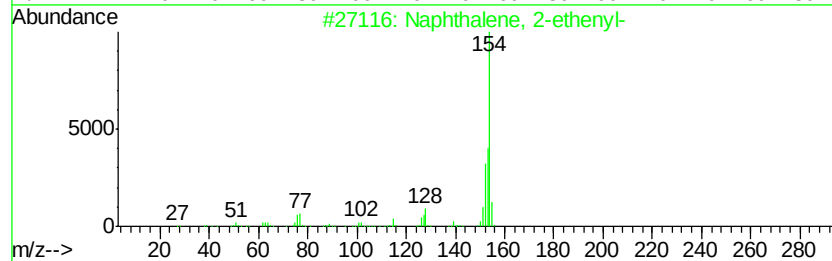
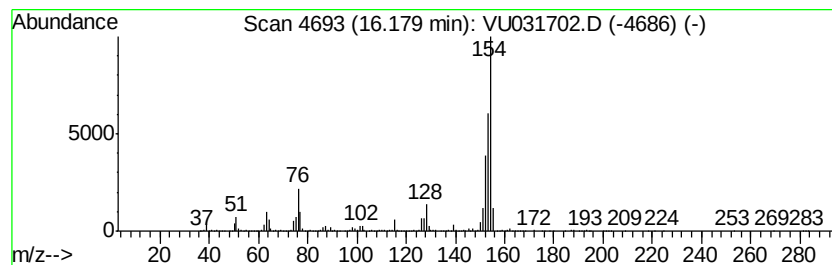
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 Naphthalene, 2-ethenyl- Concentration Rank 18

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 16.18 | 73.71 ug/L | 4202350 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-ethenyl- | 154 | C12H10  | 000827-54-3 | 93   |
| 2    |    | Acenaphthene            | 154 | C12H10  | 000083-32-9 | 93   |
| 3    |    | Acenaphthene            | 154 | C12H10  | 000083-32-9 | 93   |
| 4    |    | Biphenyl                | 154 | C12H10  | 000092-52-4 | 90   |
| 5    |    | Naphthalene, 2-ethenyl- | 154 | C12H10  | 000827-54-3 | 76   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

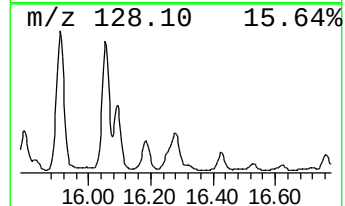
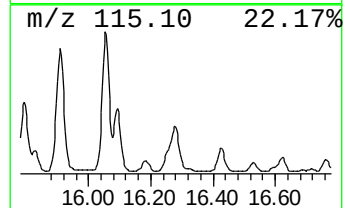
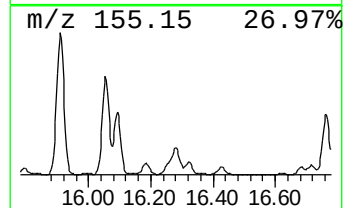
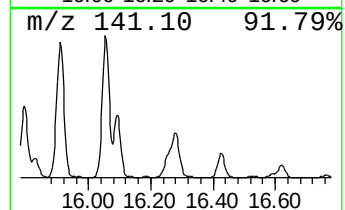
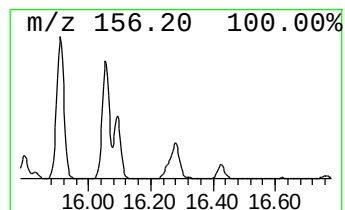
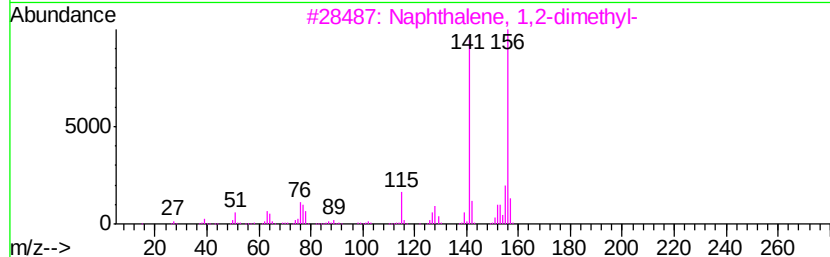
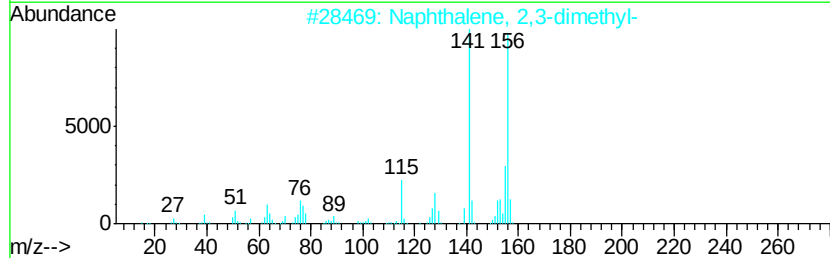
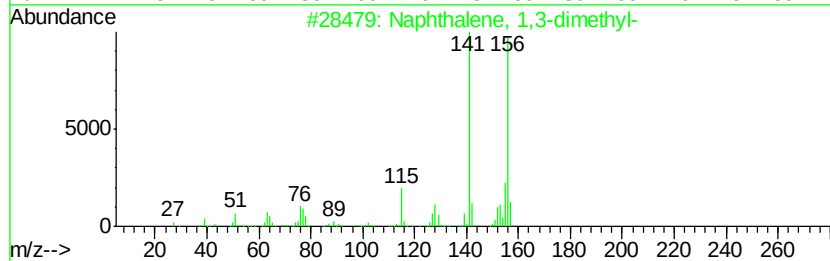
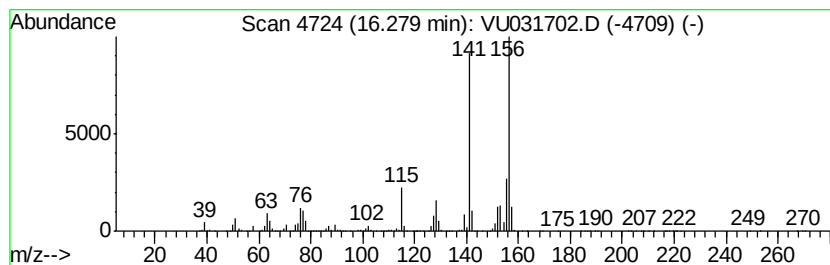
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 36 Naphthalene, 1,3-dimethyl- Concentration Rank 10

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 16.28 | 129.42 ug/L | 7378620 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 98   |
| 2    |    | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 98   |
| 3    |    | Naphthalene, 1,2-dimethyl- | 156 | C12H12  | 000573-98-8 | 98   |
| 4    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 98   |
| 5    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

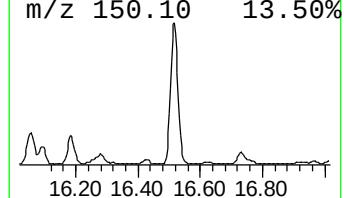
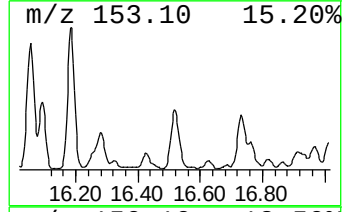
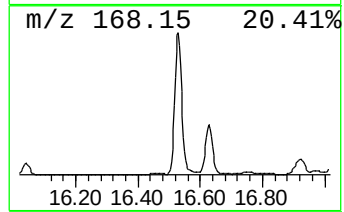
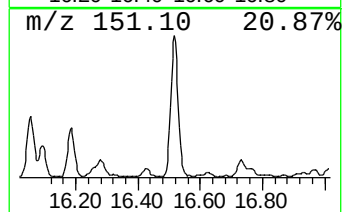
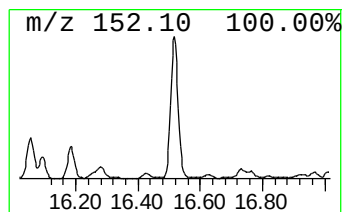
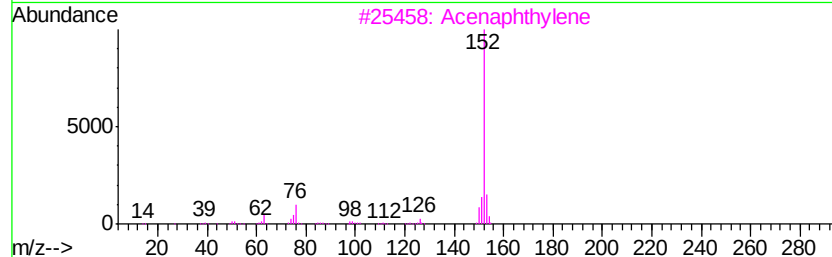
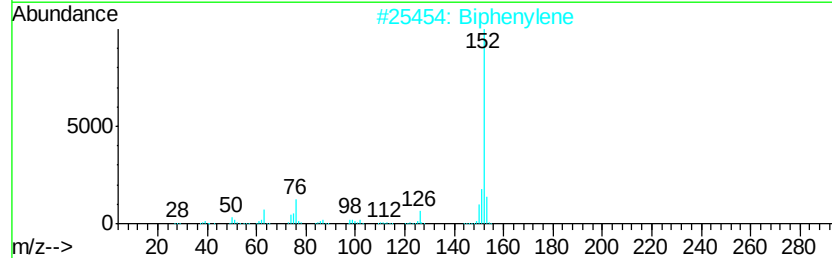
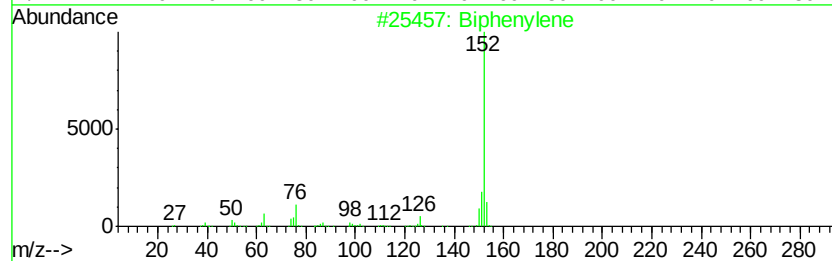
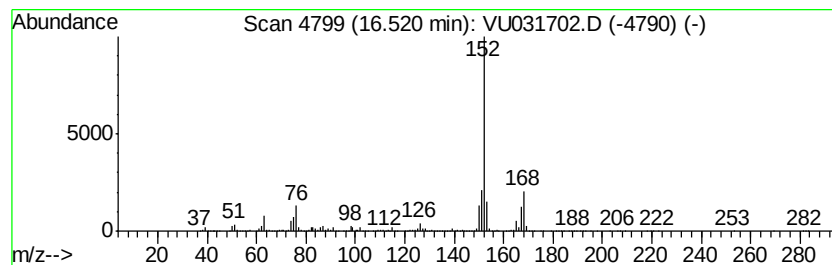
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 38 Biphenylene Concentration Rank 12

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 16.52 | 105.37 ug/L | 6007660 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Biphenylene                         | 152 | C12H8     | 000259-79-0  | 93   |
| 2    |    | Biphenylene                         | 152 | C12H8     | 000259-79-0  | 89   |
| 3    |    | Acenaphthylene                      | 152 | C12H8     | 000208-96-8  | 70   |
| 4    |    | Acenaphthylene                      | 152 | C12H8     | 000208-96-8  | 64   |
| 5    |    | 1-acenaphthenol, trifluoroacetat... | 266 | C14H9F3O2 | 1000376-45-2 | 62   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

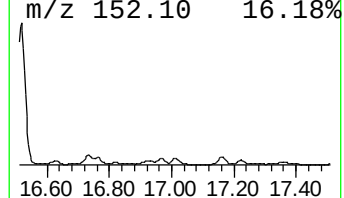
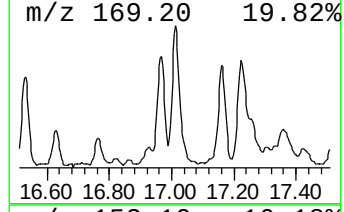
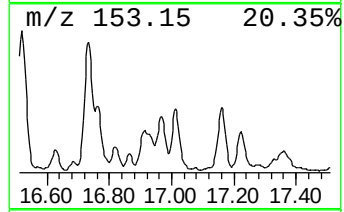
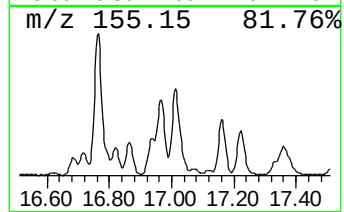
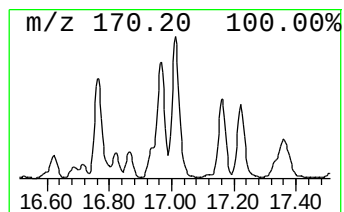
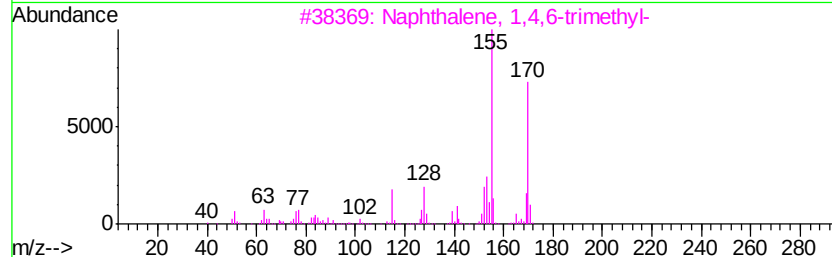
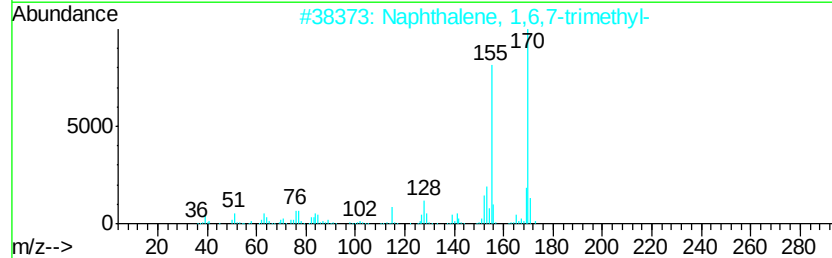
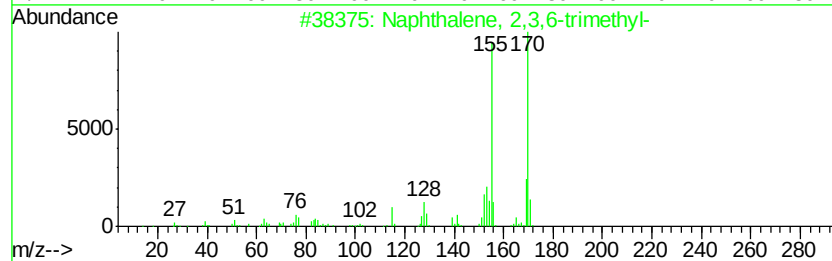
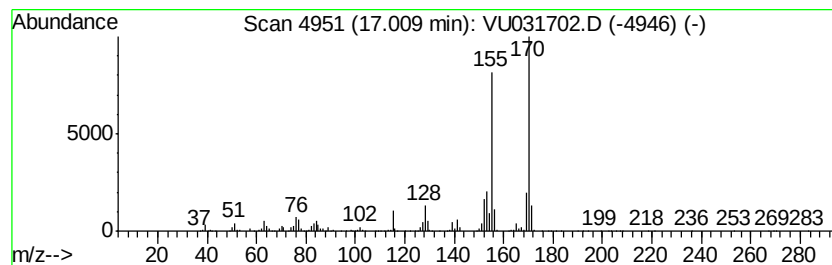
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 Naphthalene, 2,3,6-trimethyl- Concentration Rank 21

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 17.01 | 47.05 ug/L | 2682520 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 98   |
| 2    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 98   |
| 3    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 96   |
| 4    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 95   |
| 5    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU050219\  
Data File : VU031702.D  
Acq On : 02 May 2019 11:15  
Operator : JC/SP  
Sample : K2603-16DL 10X  
Misc : 5.10µ/5mL/100µL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
GAJ68DL

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

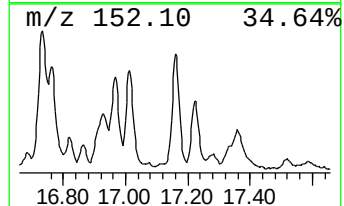
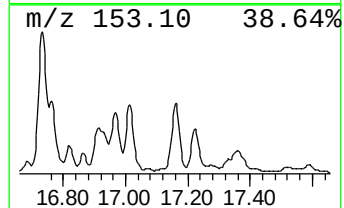
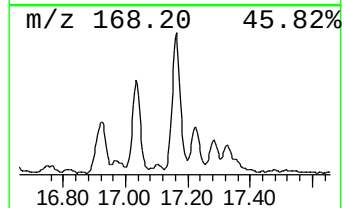
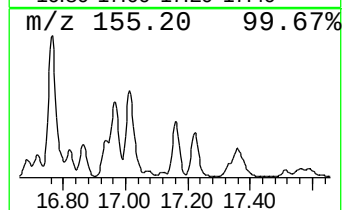
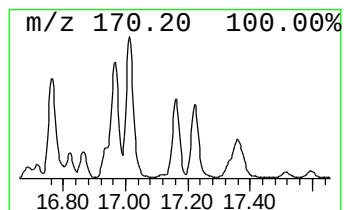
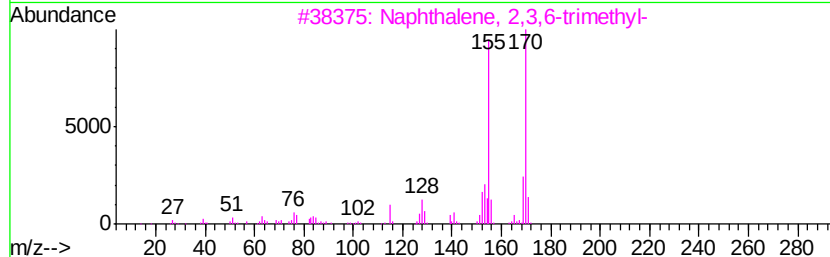
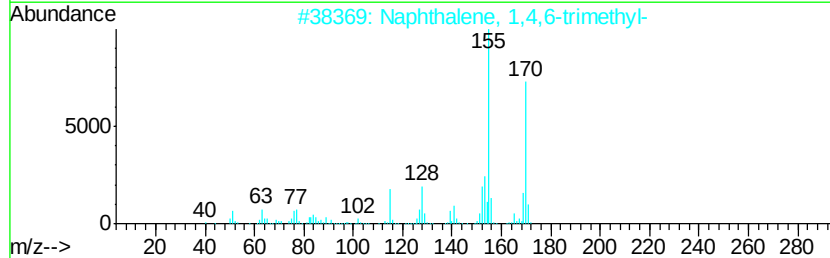
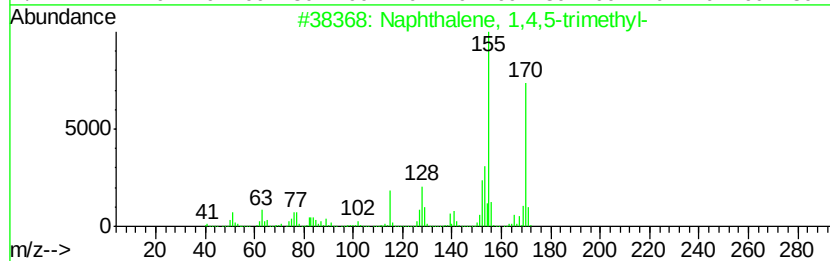
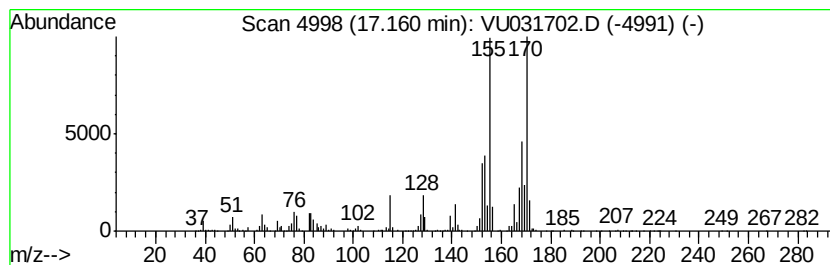
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 40 Naphthalene, 1,4,5-trimethyl- Concentration Rank 24

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 17.16 | 31.43 ug/L | 1792080 | 1,4-Dichlorobenzene-d4 | 11.48 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 95   |
| 2    |    | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 93   |
| 3    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 92   |
| 4    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 91   |
| 5    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 91   |





Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU050219\  
 Data File : VU031702.D  
 Acq On : 02 May 2019 11:15  
 Operator : JC/SP  
 Sample : K2603-16DL 10X  
 Misc : 5.10g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 GAJ68DL

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 |
| (DEL) Alkane: Str... | 9.44  | 10.3    | ug/L  | 625945   | 2                     | 9.09  | 3054340 | 50.0 |
| (DEL) Alkane: Str... | 9.95  | 5.1     | ug/L  | 312920   | 2                     | 9.09  | 3054340 | 50.0 |
| (DEL) Alkane: Cyc... | 10.04 | 3.9     | ug/L  | 237108   | 2                     | 9.09  | 3054340 | 50.0 |
| Benzene, propyl-     | 10.58 | 4.2     | ug/L  | 239773   | 3                     | 11.48 | 2850630 | 50.0 |
| Benzene, 1-ethyl-... | 10.70 | 21.8    | ug/L  | 1240760  | 3                     | 11.48 | 2850630 | 50.0 |
| (DEL) Alkane: Str... | 10.81 | 42.2    | ug/L  | 2404670  | 3                     | 11.48 | 2850630 | 50.0 |
| (DEL) Alkane: Str... | 11.11 | 9.6     | ug/L  | 546190   | 3                     | 11.48 | 2850630 | 50.0 |
| Benzene, 1,2,3-tr... | 11.14 | 78.1    | ug/L  | 4450260  | 3                     | 11.48 | 2850630 | 50.0 |
| Benzene, 1-etheny... | 11.21 | 78.5    | ug/L  | 4475600  | 3                     | 11.48 | 2850630 | 50.0 |
| Benzene, 1-propenyl- | 11.26 | 25.9    | ug/L  | 1479480  | 3                     | 11.48 | 2850630 | 50.0 |
| (DEL) Alkane: Cyc... | 11.40 | 12.9    | ug/L  | 733369   | 3                     | 11.48 | 2850630 | 50.0 |
| Indane               | 11.76 | 17.0    | ug/L  | 967421   | 3                     | 11.48 | 2850630 | 50.0 |
| Indene               | 11.98 | 247.0   | ug/L  | 14082700 | 3                     | 11.48 | 2850630 | 50.0 |
| (DEL) Alkane: Str... | 12.02 | 81.0    | ug/L  | 4615070  | 3                     | 11.48 | 2850630 | 50.0 |
| (DEL) Alkane: Str... | 12.35 | 14.0    | ug/L  | 797191   | 3                     | 11.48 | 2850630 | 50.0 |
| 2,4-Dimethylstyrene  | 12.49 | 11.6    | ug/L  | 660567   | 3                     | 11.48 | 2850630 | 50.0 |
| (DEL) Alkane: Cyc... | 12.61 | 26.4    | ug/L  | 1505340  | 3                     | 11.48 | 2850630 | 50.0 |
| Benzene, 1,2,4,5-... | 12.70 | 65.8    | ug/L  | 3752490  | 3                     | 11.48 | 2850630 | 50.0 |
| Benzene, 4-etheny... | 12.82 | 24.9    | ug/L  | 1421600  | 3                     | 11.48 | 2850630 | 50.0 |
| (DEL) Alkane: Str... | 13.12 | 161.5   | ug/L  | 9206370  | 3                     | 11.48 | 2850630 | 50.0 |
| 2-Methylindene       | 13.18 | 113.2   | ug/L  | 6456040  | 3                     | 11.48 | 2850630 | 50.0 |
| Benzene, (1-methy... | 13.29 | 196.7   | ug/L  | 11214700 | 3                     | 11.48 | 2850630 | 50.0 |
| Azulene              | 13.74 | 1073.1  | ug/L  | 61179700 | 3                     | 11.48 | 2850630 | 50.0 |
| (DEL) Alkane: Str... | 14.14 | 80.9    | ug/L  | 4614800  | 3                     | 11.48 | 2850630 | 50.0 |
| 1H-Indene, 1,1-di... | 14.33 | 64.5    | ug/L  | 3677100  | 3                     | 11.48 | 2850630 | 50.0 |
| 1H-Indene, 1-ethy... | 14.87 | 1209.1  | ug/L  | 68935800 | 3                     | 11.48 | 2850630 | 50.0 |
| Naphthalene, 1-et... | 15.79 | 82.0    | ug/L  | 4675620  | 3                     | 11.48 | 2850630 | 50.0 |
| Naphthalene, 1,6-... | 15.91 | 382.6   | ug/L  | 21814200 | 3                     | 11.48 | 2850630 | 50.0 |
| Naphthalene, 2,3-... | 16.05 | 316.5   | ug/L  | 18045700 | 3                     | 11.48 | 2850630 | 50.0 |
| Naphthalene, 1,7-... | 16.09 | 146.4   | ug/L  | 8346160  | 3                     | 11.48 | 2850630 | 50.0 |
| Naphthalene, 2-et... | 16.18 | 73.7    | ug/L  | 4202350  | 3                     | 11.48 | 2850630 | 50.0 |
| Naphthalene, 1,3-... | 16.28 | 129.4   | ug/L  | 7378620  | 3                     | 11.48 | 2850630 | 50.0 |
| Biphenylene          | 16.52 | 105.4   | ug/L  | 6007660  | 3                     | 11.48 | 2850630 | 50.0 |
| Naphthalene, 2,3,... | 17.01 | 47.0    | ug/L  | 2682520  | 3                     | 11.48 | 2850630 | 50.0 |
| Naphthalene, 1,4,... | 17.16 | 31.4    | ug/L  | 1792080  | 3                     | 11.48 | 2850630 | 50.0 |