

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
 Data File : VV007787.D  
 Acq On : 23 Sep 2018 14:43  
 Operator : SY/MD  
 Sample : J5014-02  
 Misc : 5.0 mL/MSVOA V/WATER  
 ALS Vial : 65 Sample Multiplier: 1

Instrument :  
 MSVOA\_V  
 ClientSampleId :  
 E8JB2

## 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\_V\METHOD\SOMVLM092318WMA.M  
 Title : VOC Analysis

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|--------|----------|------------|----------|-----------|-------|-------------|------------|--------------|------------|
| 1      | 1.322    | 64         | 72       | 91        | rVB   | 58041       | 63205      | 4.22%        | 0.418%     |
| 2      | 1.585    | 147        | 154      | 166       | rVB   | 61488       | 69854      | 4.66%        | 0.462%     |
| 3      | 2.139    | 316        | 326      | 338       | rBV   | 149511      | 212832     | 14.21%       | 1.406%     |
| 4      | 3.238    | 663        | 668      | 674       | rVB2  | 290         | 333        | 0.02%        | 0.002%     |
| 5      | 3.952    | 875        | 890      | 911       | rBV   | 36943       | 96067      | 6.41%        | 0.635%     |
| 6      | 4.167    | 955        | 957      | 962       | rVB   | 255         | 166        | 0.01%        | 0.001%     |
| 7      | 4.409    | 1015       | 1032     | 1054      | rBV   | 87265       | 210059     | 14.03%       | 1.388%     |
| 8      | 4.926    | 1189       | 1193     | 1199      | rVB   | 160         | 205        | 0.01%        | 0.001%     |
| 9      | 5.103    | 1228       | 1248     | 1274      | rBV2  | 217060      | 527364     | 35.21%       | 3.485%     |
| 10     | 5.363    | 1325       | 1329     | 1333      | rVB   | 216         | 200        | 0.01%        | 0.001%     |
| 11     | 5.408    | 1337       | 1343     | 1345      | rBV   | 227         | 167        | 0.01%        | 0.001%     |
| 12     | 5.527    | 1377       | 1380     | 1383      | rVB   | 216         | 136        | 0.01%        | 0.001%     |
| 13     | 5.672    | 1410       | 1425     | 1457      | rBV   | 193789      | 392017     | 26.18%       | 2.591%     |
| 14     | 5.894    | 1491       | 1494     | 1497      | rBV2  | 203         | 153        | 0.01%        | 0.001%     |
| 15     | 5.913    | 1497       | 1500     | 1503      | rVB2  | 204         | 138        | 0.01%        | 0.001%     |
| 16     | 5.968    | 1506       | 1517     | 1528      | rVV4  | 6269        | 11607      | 0.78%        | 0.077%     |
| 17     | 6.125    | 1552       | 1566     | 1592      | rBV   | 115932      | 236569     | 15.80%       | 1.563%     |
| 18     | 6.244    | 1602       | 1603     | 1610      | rVV3  | 351         | 304        | 0.02%        | 0.002%     |
| 19     | 6.273    | 1610       | 1612     | 1615      | rVB2  | 325         | 173        | 0.01%        | 0.001%     |
| 20     | 6.392    | 1647       | 1649     | 1653      | rVV2  | 162         | 155        | 0.01%        | 0.001%     |
| 21     | 7.035    | 1834       | 1849     | 1873      | rBV   | 61625       | 110273     | 7.36%        | 0.729%     |
| 22     | 7.306    | 1929       | 1933     | 1938      | rBV2  | 483         | 482        | 0.03%        | 0.003%     |
| 23     | 7.363    | 1938       | 1951     | 1970      | rBV   | 260024      | 465640     | 31.09%       | 3.077%     |
| 24     | 7.669    | 2035       | 2046     | 2066      | rBV   | 42428       | 76186      | 5.09%        | 0.503%     |
| 25     | 7.759    | 2072       | 2074     | 2078      | rBV2  | 262         | 188        | 0.01%        | 0.001%     |
| 26     | 7.778    | 2079       | 2080     | 2084      | rVB   | 369         | 187        | 0.01%        | 0.001%     |
| 27     | 7.836    | 2089       | 2098     | 2105      | rBV5  | 1285        | 1928       | 0.13%        | 0.013%     |
| 28     | 7.955    | 2132       | 2135     | 2137      | rBV   | 274         | 173        | 0.01%        | 0.001%     |
| 29     | 8.023    | 2151       | 2156     | 2161      | rVV7  | 874         | 978        | 0.07%        | 0.006%     |
| 30     | 8.138    | 2179       | 2192     | 2222      | rBV   | 124119      | 249202     | 16.64%       | 1.647%     |
| 31     | 8.338    | 2245       | 2254     | 2263      | rVB3  | 13860       | 22170      | 1.48%        | 0.147%     |
| 32     | 8.402    | 2264       | 2274     | 2282      | rBV   | 16923       | 29062      | 1.94%        | 0.192%     |
| 33     | 8.527    | 2310       | 2313     | 2314      | rBV2  | 323         | 182        | 0.01%        | 0.001%     |
| 34     | 8.550    | 2314       | 2320     | 2330      | rVB2  | 3305        | 4795       | 0.32%        | 0.032%     |

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 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\_V\METHOD\SOMVLM092318WMA.M  
 Title : VOC Analysis

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 8.640  | 2330 | 2348 | 2360 | rBV5 | 39143  | 85959   | 5.74%  | 0.568% |
| 36 | 8.704  | 2362 | 2368 | 2382 | rVB6 | 17765  | 31314   | 2.09%  | 0.207% |
| 37 | 8.772  | 2382 | 2389 | 2398 | rBV6 | 3459   | 5687    | 0.38%  | 0.038% |
| 38 | 8.833  | 2398 | 2408 | 2418 | rBV2 | 22133  | 38309   | 2.56%  | 0.253% |
| 39 | 8.900  | 2418 | 2429 | 2446 | rBV  | 299307 | 496230  | 33.13% | 3.279% |
| 40 | 9.048  | 2462 | 2475 | 2486 | rBV5 | 42452  | 99297   | 6.63%  | 0.656% |
| 41 | 9.109  | 2486 | 2494 | 2500 | rBV3 | 24790  | 38372   | 2.56%  | 0.254% |
| 42 | 9.154  | 2502 | 2508 | 2512 | rBV3 | 23595  | 34496   | 2.30%  | 0.228% |
| 43 | 9.367  | 2563 | 2574 | 2584 | rBV3 | 14753  | 25176   | 1.68%  | 0.166% |
| 44 | 9.453  | 2594 | 2601 | 2605 | rBV5 | 2273   | 3224    | 0.22%  | 0.021% |
| 45 | 9.521  | 2605 | 2622 | 2637 | rVV2 | 111121 | 283289  | 18.92% | 1.872% |
| 46 | 9.592  | 2639 | 2644 | 2665 | rVB3 | 47121  | 97390   | 6.50%  | 0.644% |
| 47 | 9.723  | 2676 | 2685 | 2688 | rBV2 | 33043  | 47383   | 3.16%  | 0.313% |
| 48 | 9.756  | 2690 | 2695 | 2704 | rVB4 | 75231  | 107960  | 7.21%  | 0.713% |
| 49 | 9.846  | 2715 | 2723 | 2734 | rBV3 | 91076  | 180808  | 12.07% | 1.195% |
| 50 | 9.958  | 2751 | 2758 | 2770 | rVB2 | 20251  | 39720   | 2.65%  | 0.262% |
| 51 | 10.180 | 2819 | 2827 | 2839 | rVB4 | 34838  | 58588   | 3.91%  | 0.387% |
| 52 | 10.267 | 2843 | 2854 | 2883 | rVB2 | 219026 | 496360  | 33.14% | 3.280% |
| 53 | 10.402 | 2889 | 2896 | 2901 | rBV  | 19608  | 29706   | 1.98%  | 0.196% |
| 54 | 10.498 | 2917 | 2926 | 2937 | rBV2 | 92965  | 202288  | 13.51% | 1.337% |
| 55 | 10.784 | 3005 | 3015 | 3033 | rVB3 | 109409 | 235270  | 15.71% | 1.555% |
| 56 | 10.926 | 3052 | 3059 | 3062 | rBV3 | 42368  | 60414   | 4.03%  | 0.399% |
| 57 | 10.961 | 3062 | 3070 | 3093 | rVB  | 494123 | 849023  | 56.69% | 5.611% |
| 58 | 11.202 | 3138 | 3145 | 3152 | rBV3 | 57355  | 81156   | 5.42%  | 0.536% |
| 59 | 11.299 | 3165 | 3175 | 3193 | rBV  | 369999 | 612028  | 40.87% | 4.044% |
| 60 | 11.386 | 3194 | 3202 | 3223 | rVB  | 267986 | 432141  | 28.86% | 2.856% |
| 61 | 11.588 | 3253 | 3265 | 3271 | rBV4 | 90222  | 201577  | 13.46% | 1.332% |
| 62 | 11.678 | 3282 | 3293 | 3313 | rVB4 | 569667 | 1302255 | 86.95% | 8.606% |
| 63 | 11.797 | 3323 | 3330 | 3343 | rVB4 | 34343  | 55705   | 3.72%  | 0.368% |
| 64 | 11.871 | 3345 | 3353 | 3370 | rVB  | 113903 | 189682  | 12.67% | 1.253% |
| 65 | 11.961 | 3371 | 3381 | 3385 | rBV  | 232740 | 344347  | 22.99% | 2.276% |
| 66 | 12.067 | 3405 | 3414 | 3435 | rVB  | 431965 | 726853  | 48.53% | 4.803% |
| 67 | 12.167 | 3436 | 3445 | 3450 | rBV  | 89441  | 144755  | 9.67%  | 0.957% |
| 68 | 12.309 | 3476 | 3489 | 3496 | rBV6 | 41883  | 93085   | 6.22%  | 0.615% |
| 69 | 12.363 | 3497 | 3506 | 3517 | rVB3 | 134361 | 235496  | 15.72% | 1.556% |
| 70 | 12.463 | 3518 | 3537 | 3546 | rBV  | 352502 | 620898  | 41.46% | 4.103% |
| 71 | 12.518 | 3546 | 3554 | 3569 | rVB  | 490623 | 754985  | 50.41% | 4.989% |

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Method : Z:\VOASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
 Title : VOC Analysis

|     |        |      |      |      |      |        |         |         |        |
|-----|--------|------|------|------|------|--------|---------|---------|--------|
| 72  | 12.601 | 3570 | 3580 | 3586 | rBV  | 67733  | 110698  | 7.39%   | 0.732% |
| 73  | 12.636 | 3587 | 3591 | 3596 | rVV  | 32429  | 31080   | 2.08%   | 0.205% |
| 74  | 12.778 | 3628 | 3635 | 3649 | rVB  | 175030 | 264455  | 17.66%  | 1.748% |
| 75  | 12.845 | 3649 | 3656 | 3663 | rVB2 | 53503  | 70434   | 4.70%   | 0.465% |
| 76  | 12.936 | 3664 | 3684 | 3700 | rBV2 | 761774 | 1497623 | 100.00% | 9.897% |
| 77  | 13.051 | 3713 | 3720 | 3729 | rBV  | 72860  | 100448  | 6.71%   | 0.664% |
| 78  | 13.109 | 3730 | 3738 | 3756 | rVB3 | 62184  | 121909  | 8.14%   | 0.806% |
| 79  | 13.218 | 3759 | 3772 | 3780 | rBV2 | 36072  | 69661   | 4.65%   | 0.460% |
| 80  | 13.321 | 3795 | 3804 | 3810 | rBV2 | 134565 | 214386  | 14.32%  | 1.417% |
| 81  | 13.521 | 3858 | 3866 | 3867 | rBV3 | 7719   | 9244    | 0.62%   | 0.061% |
| 82  | 13.601 | 3884 | 3891 | 3901 | rVB  | 24381  | 37938   | 2.53%   | 0.251% |
| 83  | 13.665 | 3902 | 3911 | 3918 | rBV5 | 10457  | 17848   | 1.19%   | 0.118% |
| 84  | 13.771 | 3935 | 3944 | 3954 | rBV5 | 22988  | 50414   | 3.37%   | 0.333% |
| 85  | 13.961 | 3994 | 4003 | 4019 | rBV2 | 34357  | 58488   | 3.91%   | 0.387% |
| 86  | 14.064 | 4028 | 4035 | 4039 | rBV7 | 2132   | 2793    | 0.19%   | 0.018% |
| 87  | 14.144 | 4051 | 4060 | 4074 | rVB4 | 23382  | 51991   | 3.47%   | 0.344% |
| 88  | 14.238 | 4084 | 4089 | 4090 | rBV2 | 739    | 526     | 0.04%   | 0.003% |
| 89  | 14.286 | 4095 | 4104 | 4115 | rVV  | 22996  | 35969   | 2.40%   | 0.238% |
| 90  | 14.347 | 4115 | 4123 | 4136 | rVB  | 15138  | 22094   | 1.48%   | 0.146% |
| 91  | 14.402 | 4136 | 4140 | 4144 | rBV2 | 404    | 295     | 0.02%   | 0.002% |
| 92  | 14.424 | 4144 | 4147 | 4148 | rBV2 | 253    | 162     | 0.01%   | 0.001% |
| 93  | 14.489 | 4159 | 4167 | 4174 | rVV8 | 3076   | 4943    | 0.33%   | 0.033% |
| 94  | 14.537 | 4177 | 4182 | 4190 | rVV6 | 1933   | 2642    | 0.18%   | 0.017% |
| 95  | 14.643 | 4212 | 4215 | 4218 | rBV2 | 546    | 442     | 0.03%   | 0.003% |
| 96  | 14.688 | 4218 | 4229 | 4246 | rVV  | 102354 | 156353  | 10.44%  | 1.033% |
| 97  | 14.768 | 4253 | 4254 | 4259 | rVB2 | 543    | 337     | 0.02%   | 0.002% |
| 98  | 14.826 | 4269 | 4272 | 4273 | rBV  | 303    | 144     | 0.01%   | 0.001% |
| 99  | 14.871 | 4275 | 4286 | 4299 | rVV  | 47775  | 74402   | 4.97%   | 0.492% |
| 100 | 15.032 | 4332 | 4336 | 4340 | rBV3 | 527    | 499     | 0.03%   | 0.003% |

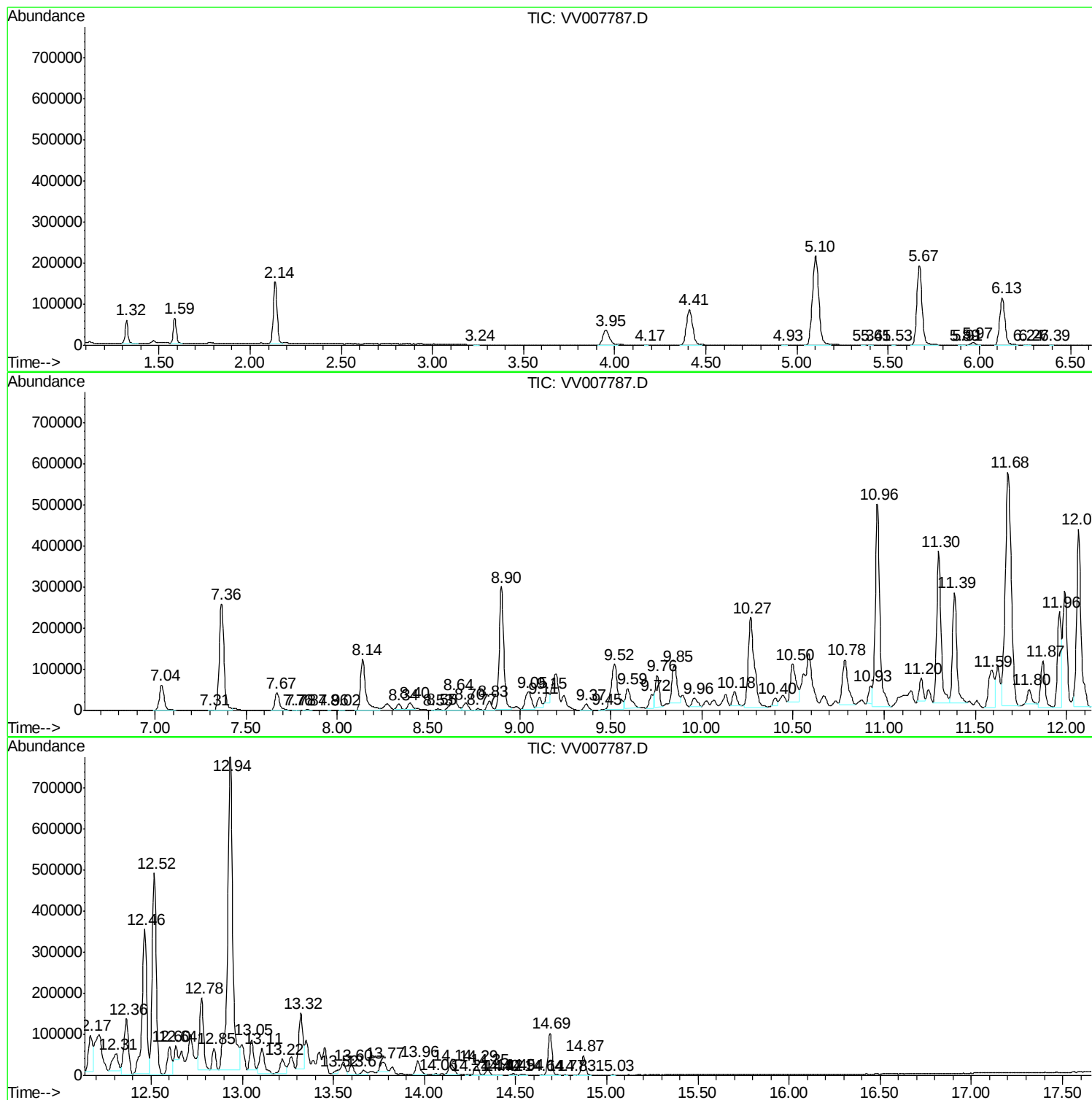
Sum of corrected areas: 15132594

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
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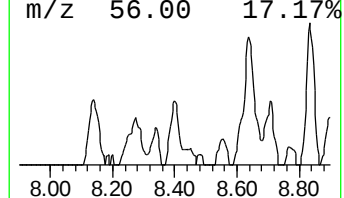
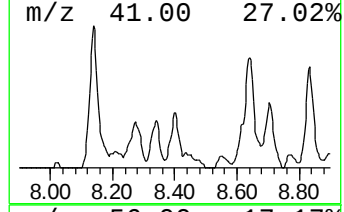
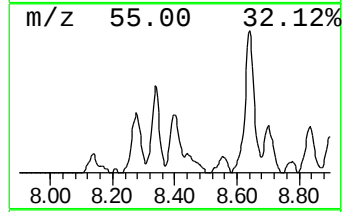
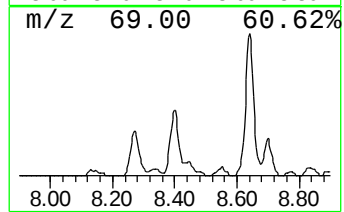
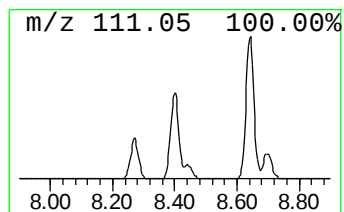
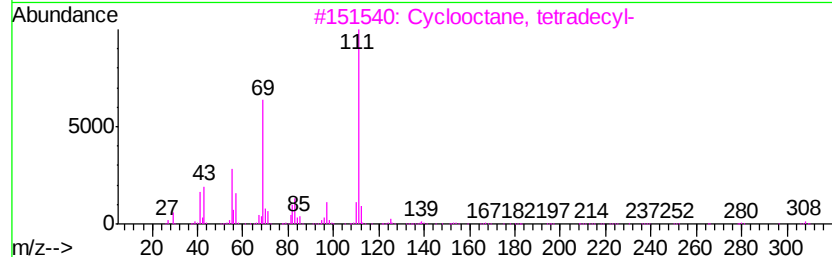
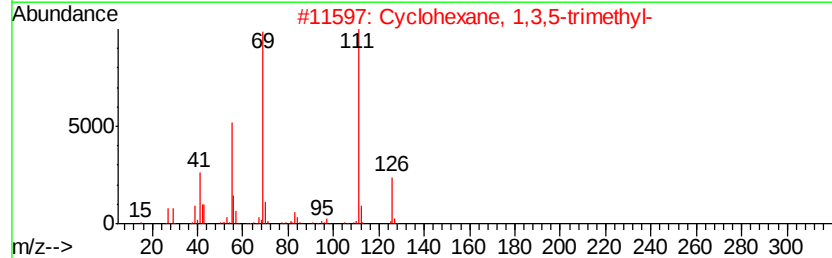
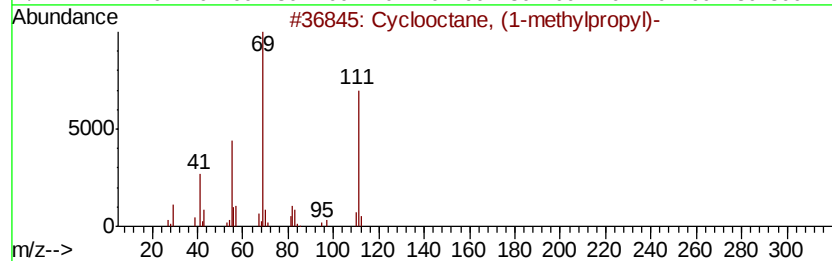
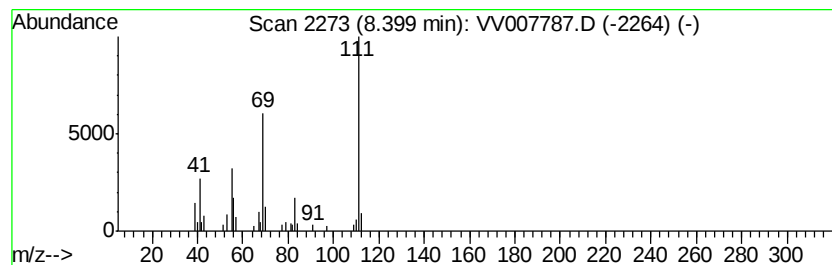
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Cyclic8.40 Concentration Rank 41

| R.T.      | EstConc                            | Area  | Relative to ISTD | R.T.        |      |
|-----------|------------------------------------|-------|------------------|-------------|------|
| 8.40      | 2.93 ug/L                          | 29062 | Chlorobenzene-d5 | 8.90        |      |
| Hit# of 5 | Tentative ID                       | MW    | MolForm          | CAS#        | Qual |
| 1         | Cvclooctane, (1-methylpropyl)-     | 168   | C12H24           | 016538-89-9 | 72   |
| 2         | Cyclohexane, 1,3,5-trimethyl-      | 126   | C9H18            | 001839-63-0 | 64   |
| 3         | Cvclooctane, tetradecyl-           | 308   | C22H44           | 149003-36-1 | 64   |
| 4         | Cyclohexane, 1,1,3-trimethyl-      | 126   | C9H18            | 003073-66-3 | 64   |
| 5         | Cyclohexane, 1-ethyl-2,4-dimethyl- | 140   | C10H20           | 061142-69-6 | 64   |



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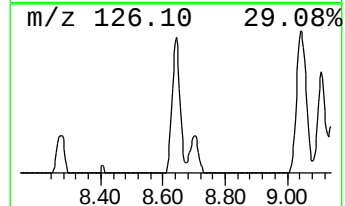
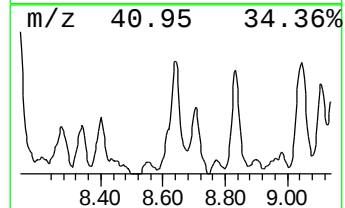
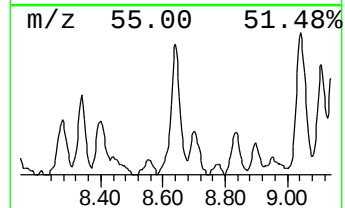
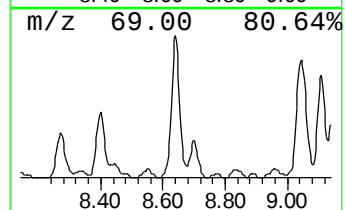
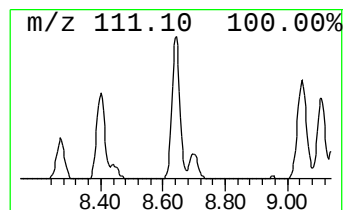
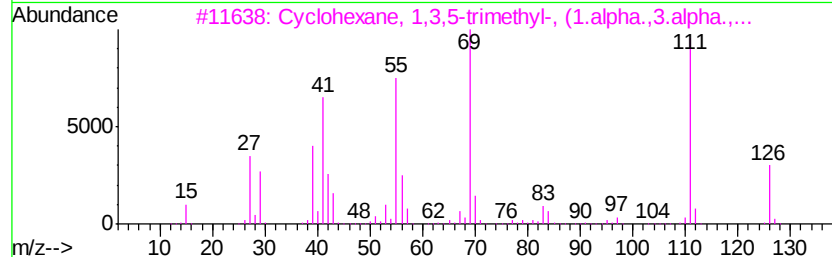
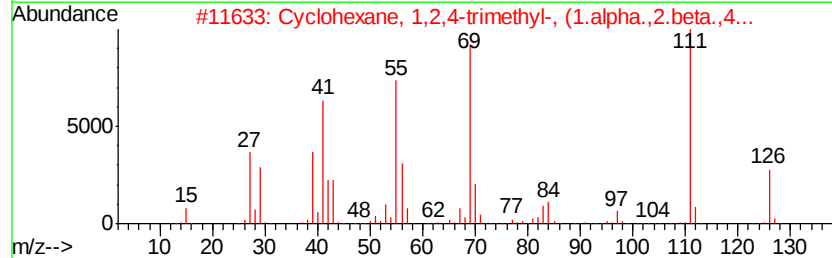
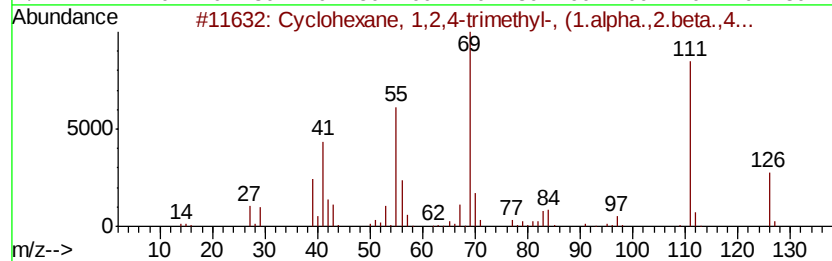
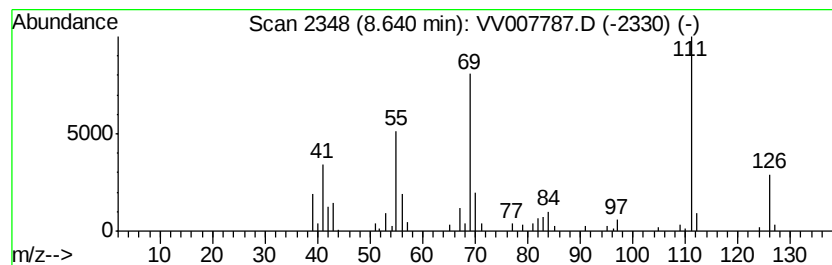
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Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Cyclic8.64 Concentration Rank 23

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.        |      |      |
|---------|-------------------------------------|--------------|------------------|-------------|------|------|
| 8.64    | 8.66 ug/L                           | 85959        | Chlorobenzene-d5 | 8.90        |      |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm     | CAS# | Qual |
| 1       | Cyclohexane, 1,2,4-trimethyl-, (... | 126          | C9H18            | 007667-60-9 | 97   |      |
| 2       | Cyclohexane, 1,2,4-trimethyl-, (... | 126          | C9H18            | 007667-60-9 | 94   |      |
| 3       | Cyclohexane, 1,3,5-trimethyl-, (... | 126          | C9H18            | 001795-26-2 | 90   |      |
| 4       | Cyclohexane, 1,3,5-trimethyl-, (... | 126          | C9H18            | 001795-27-3 | 87   |      |
| 5       | Cyclohexane, 1,3,5-trimethyl-       | 126          | C9H18            | 001839-63-0 | 83   |      |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

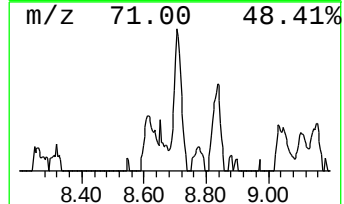
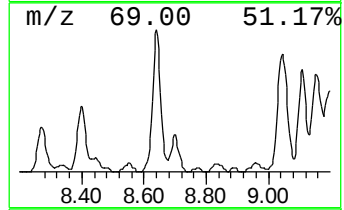
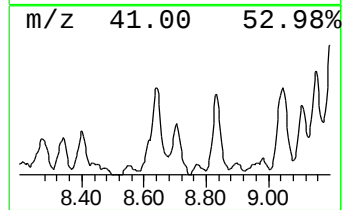
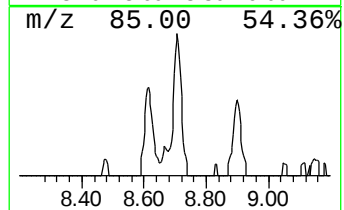
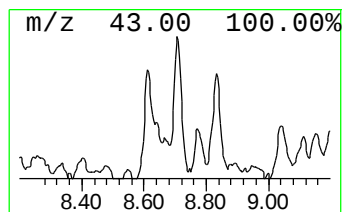
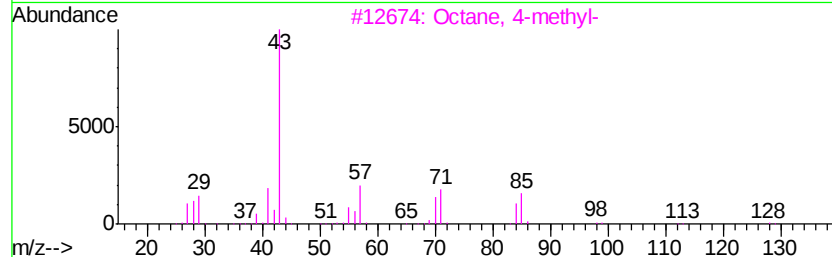
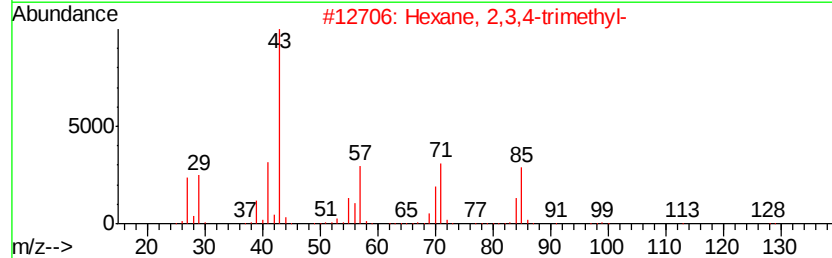
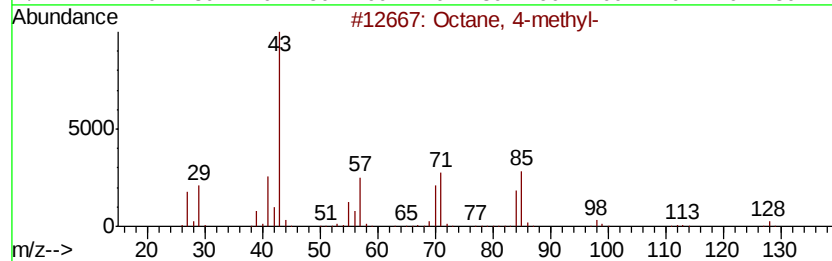
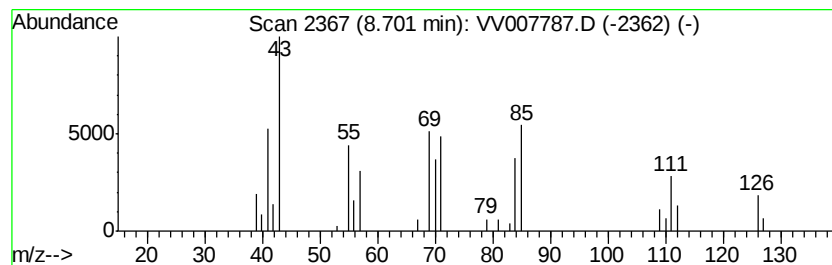
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 38

| R.T. | EstConc   | Area  | Relative to ISTD | R.T. |
|------|-----------|-------|------------------|------|
| 8.70 | 3.16 ug/L | 31314 | Chlorobenzene-d5 | 8.90 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 4-methyl-        | 128 | C9H20   | 002216-34-4 | 53   |
| 2    |    | Hexane, 2,3,4-trimethyl- | 128 | C9H20   | 000921-47-1 | 53   |
| 3    |    | Octane, 4-methyl-        | 128 | C9H20   | 002216-34-4 | 53   |
| 4    |    | Hexane, 2,3,4-trimethyl- | 128 | C9H20   | 000921-47-1 | 53   |
| 5    |    | Heptane, 2,4-dimethyl-   | 128 | C9H20   | 002213-23-2 | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampled :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

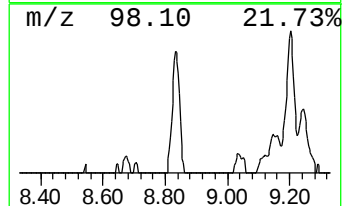
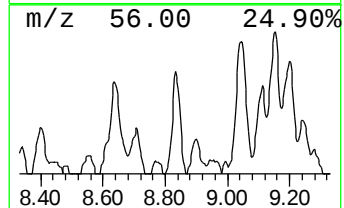
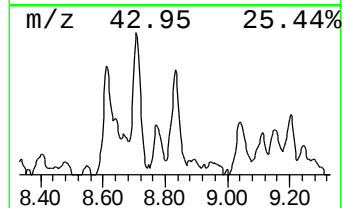
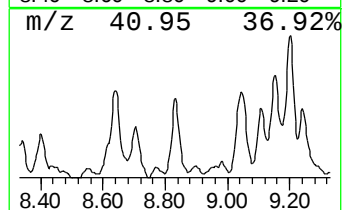
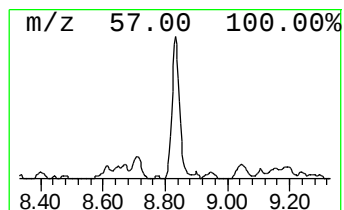
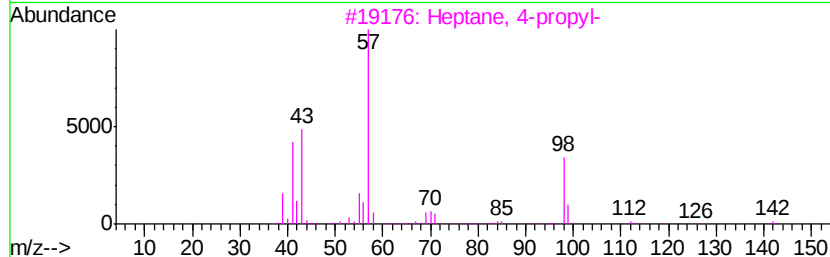
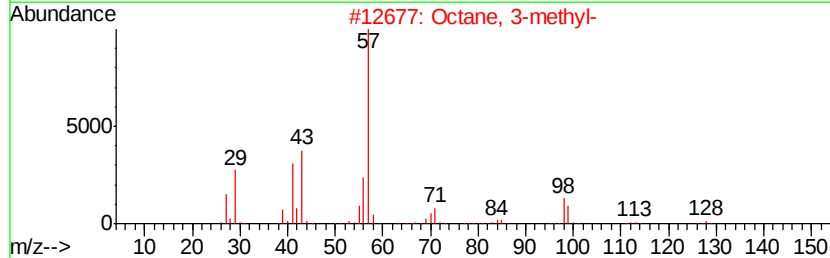
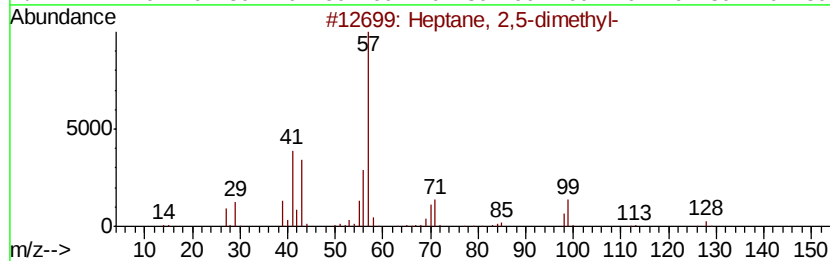
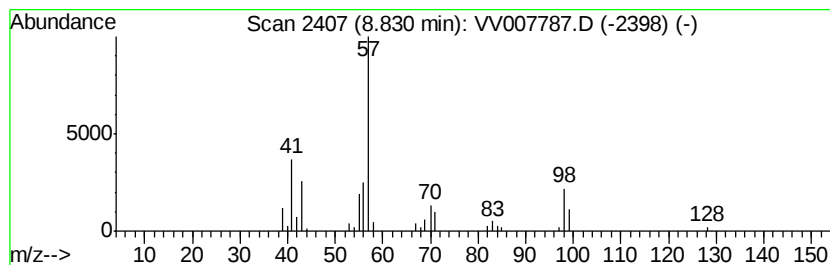
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 37

| R.T. | EstConc   | Area  | Relative to ISTD | R.T. |
|------|-----------|-------|------------------|------|
| 8.83 | 3.86 ug/L | 38309 | Chlorobenzene-d5 | 8.90 |

| Hit# of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|---------|---|------------------------|-----|---------|-------------|------|
| 1       |   | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 81   |
| 2       |   | Octane, 3-methyl-      | 128 | C9H20   | 002216-33-3 | 72   |
| 3       |   | Heptane, 4-propyl-     | 142 | C10H22  | 003178-29-8 | 72   |
| 4       |   | Octane, 3-methyl-      | 128 | C9H20   | 002216-33-3 | 72   |
| 5       |   | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

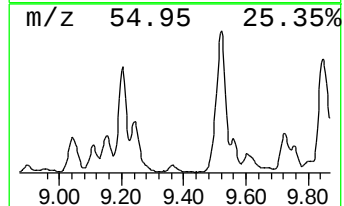
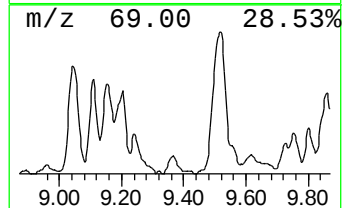
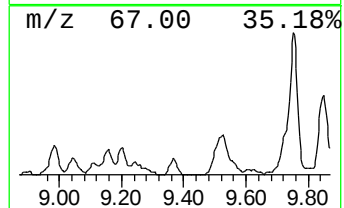
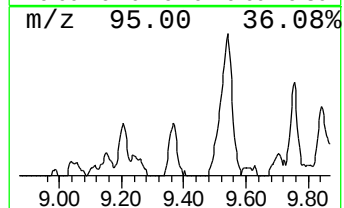
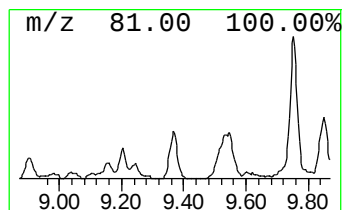
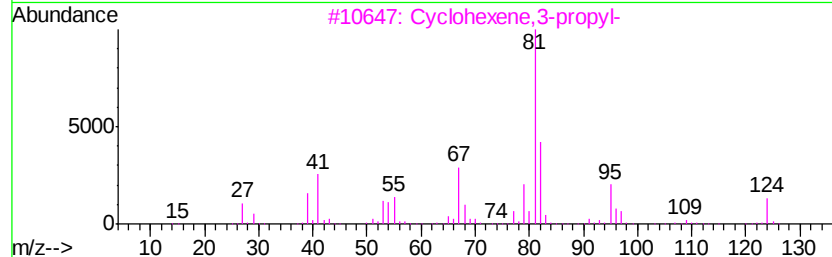
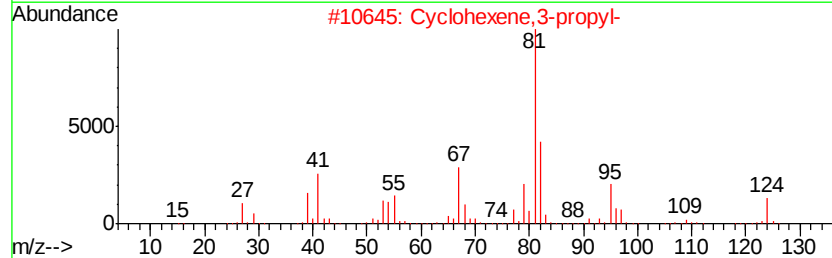
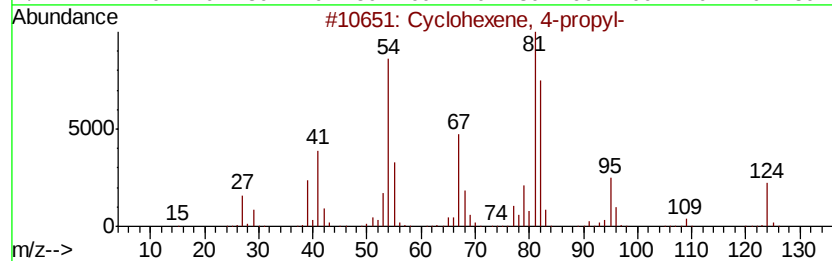
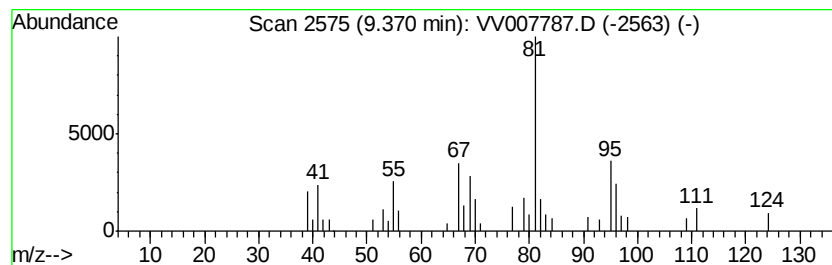
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Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 6 Cyclohexene, 4-propyl- Concentration Rank 43

| R.T. | EstConc   | Area  | Relative to ISTD | R.T. |
|------|-----------|-------|------------------|------|
| 9.37 | 2.54 ug/L | 25176 | Chlorobenzene-d5 | 8.90 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexene, 4-propyl-             | 124 | C9H16   | 013487-64-4 | 64   |
| 2    |    | Cyclohexene,3-propyl-              | 124 | C9H16   | 003983-06-0 | 58   |
| 3    |    | Cyclohexene,3-propyl-              | 124 | C9H16   | 003983-06-0 | 58   |
| 4    |    | Methyl ethyl cyclopentene          | 110 | C8H14   | 019780-56-4 | 53   |
| 5    |    | Cyclobutane, (1-methylethylidene)- | 96  | C7H12   | 001528-22-9 | 52   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

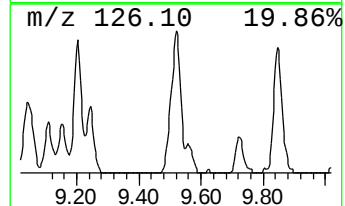
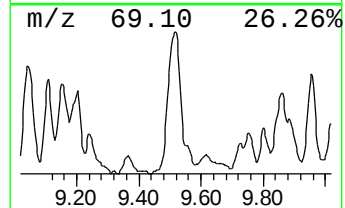
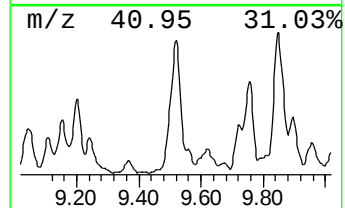
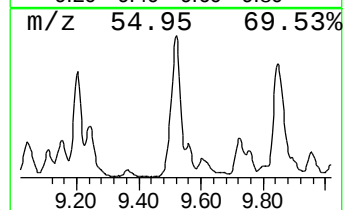
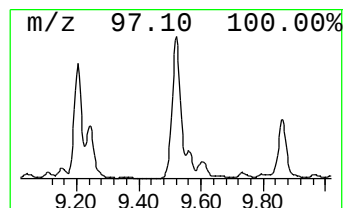
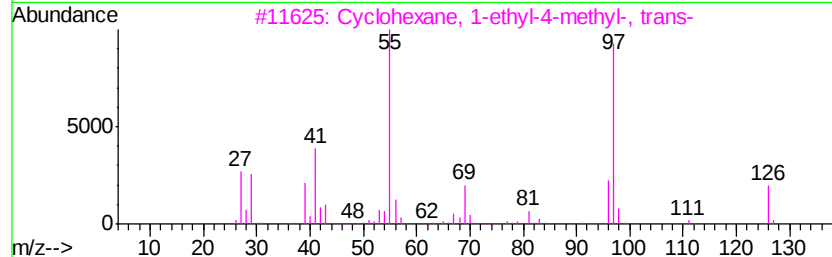
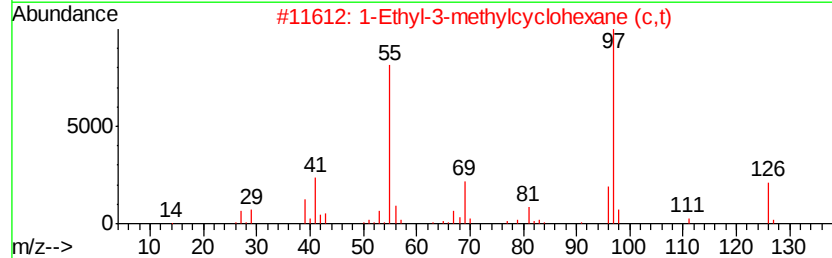
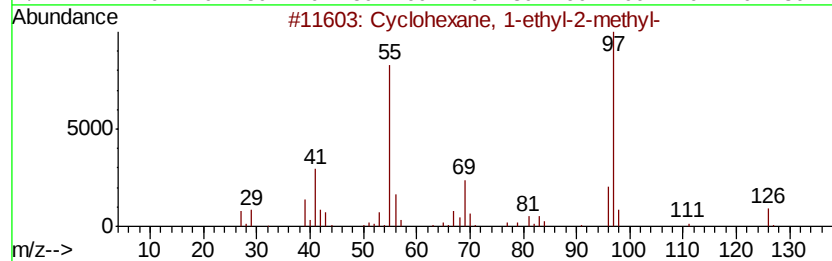
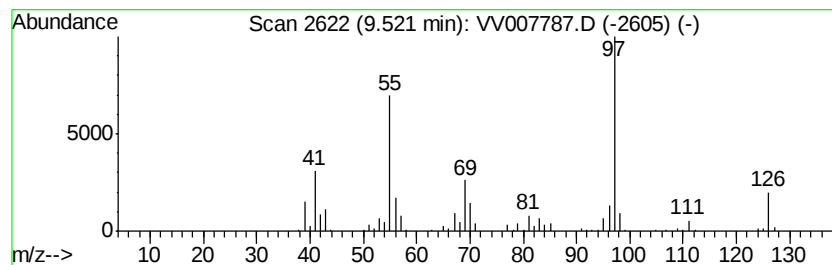
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 (DEL) Alkane: Cyclic9.52 Concentration Rank 7

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 9.52 | 28.54 ug/L | 283289 | Chlorobenzene-d5 | 8.90 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1-ethyl-2-methyl-      | 126 | C9H18   | 003728-54-9 | 81   |
| 2    |    | 1-Ethyl-3-methylcyclohexane (c,t)   | 126 | C9H18   | 003728-55-0 | 76   |
| 3    |    | Cyclohexane, 1-ethyl-4-methyl-, ... | 126 | C9H18   | 006236-88-0 | 74   |
| 4    |    | 3,5-Dimethyl-3-heptene              | 126 | C9H18   | 059643-68-4 | 72   |
| 5    |    | 1-Ethyl-4-methylcyclohexane         | 126 | C9H18   | 003728-56-1 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

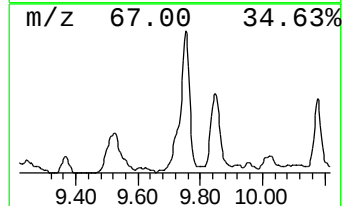
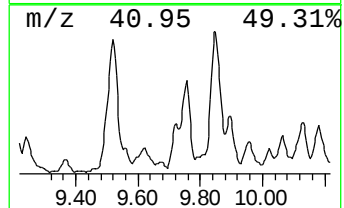
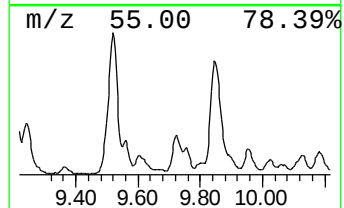
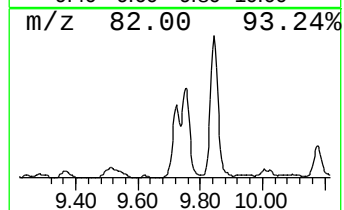
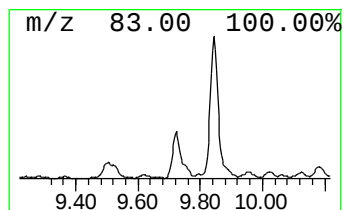
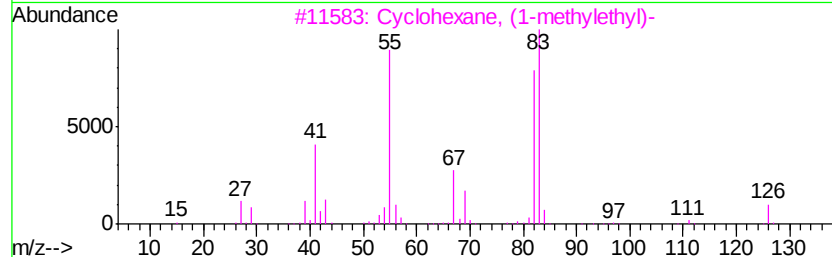
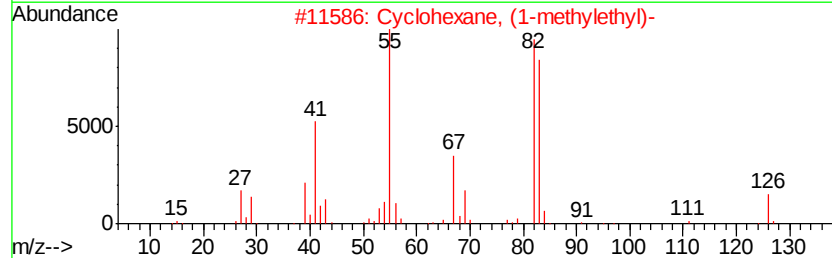
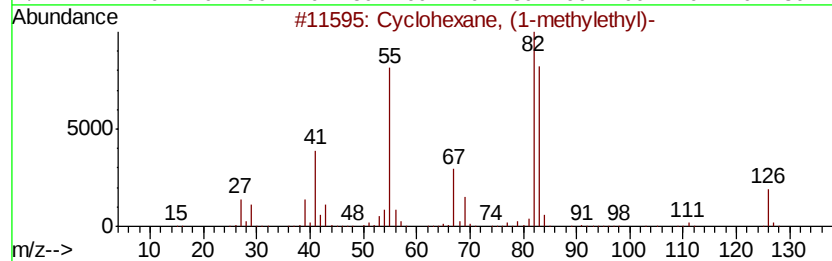
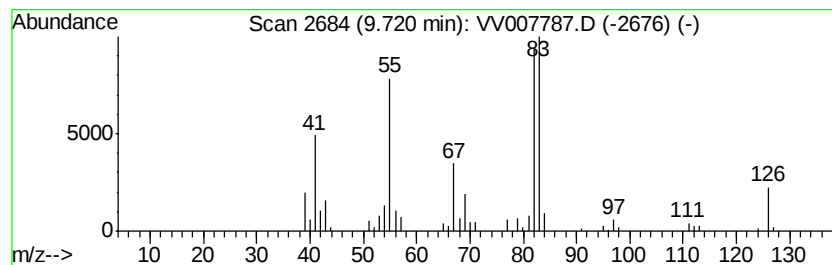
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 (DEL) Alkane: Cyclic9.72 Concentration Rank 33

| R.T. | EstConc   | Area  | Relative to ISTD | R.T. |
|------|-----------|-------|------------------|------|
| 9.72 | 4.77 ug/L | 47383 | Chlorobenzene-d5 | 8.90 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 95   |
| 2    |    | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 93   |
| 3    |    | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 91   |
| 4    |    | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 91   |
| 5    |    | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

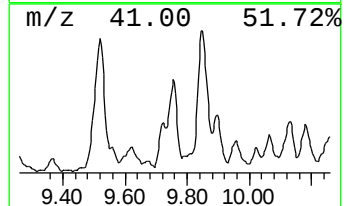
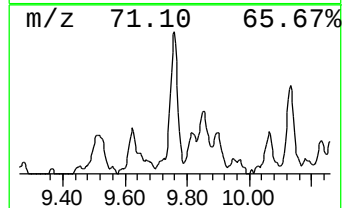
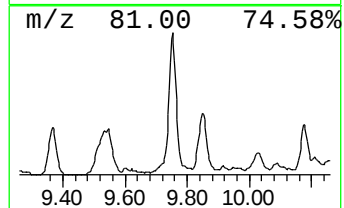
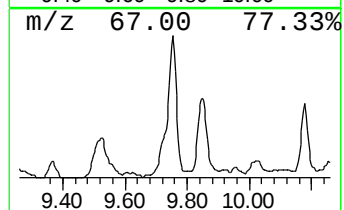
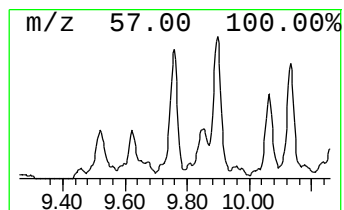
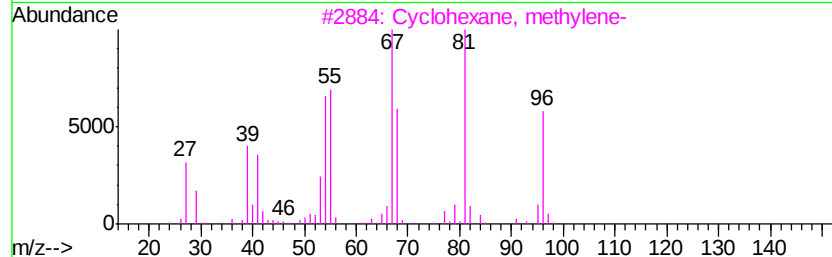
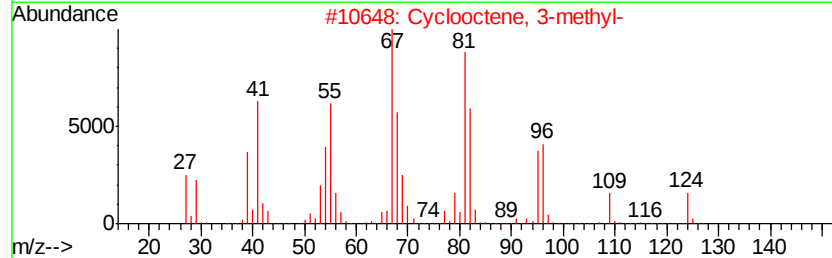
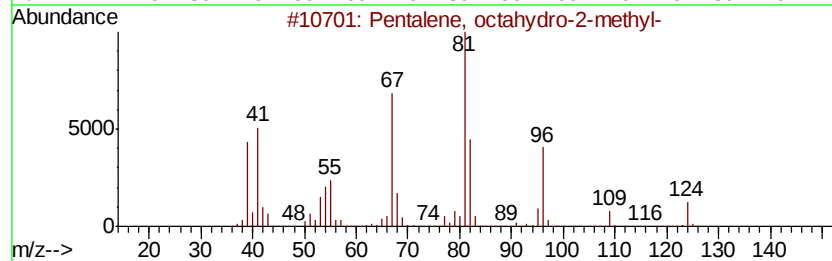
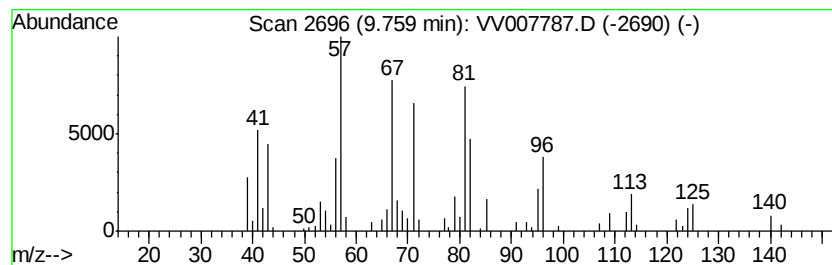
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Pentalene, octahydro-2-methyl- Concentration Rank 20

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 9.76 | 10.88 ug/L | 107960 | Chlorobenzene-d5 | 8.90 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentalene, octahydro-2-methyl-   | 124 | C9H16   | 003868-64-2 | 76   |
| 2    |    | Cyclooctene, 3-methyl-           | 124 | C9H16   | 013152-05-1 | 58   |
| 3    |    | Cyclohexane, methylene-          | 96  | C7H12   | 001192-37-6 | 49   |
| 4    |    | Cyclohexene, 1-propyl-           | 124 | C9H16   | 002539-75-5 | 46   |
| 5    |    | Cyclohexanol, 2-(1-methylethyl)- | 142 | C9H18O  | 000096-07-1 | 43   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

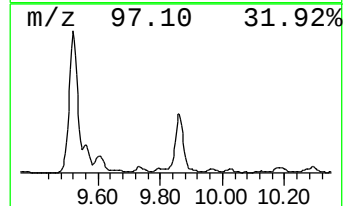
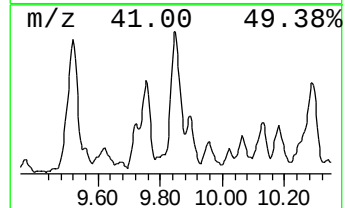
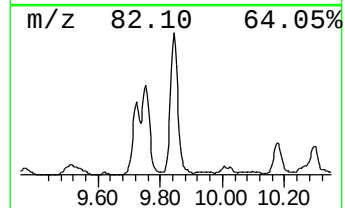
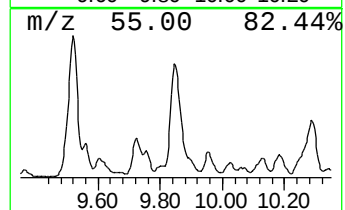
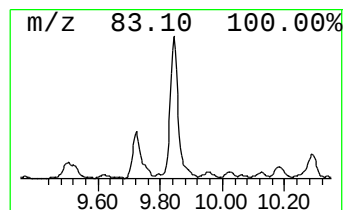
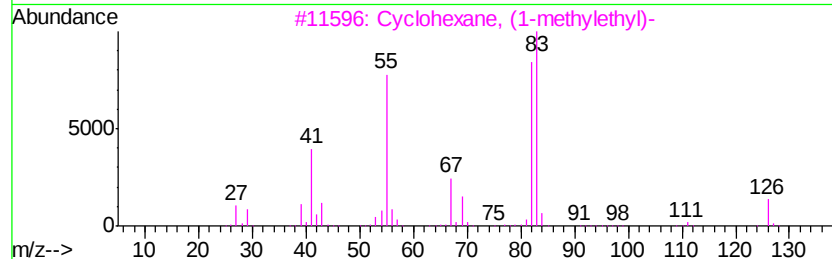
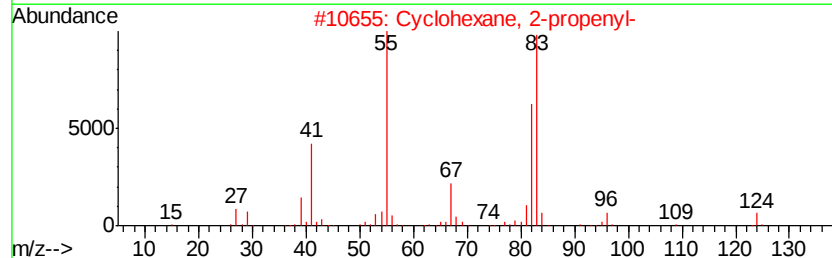
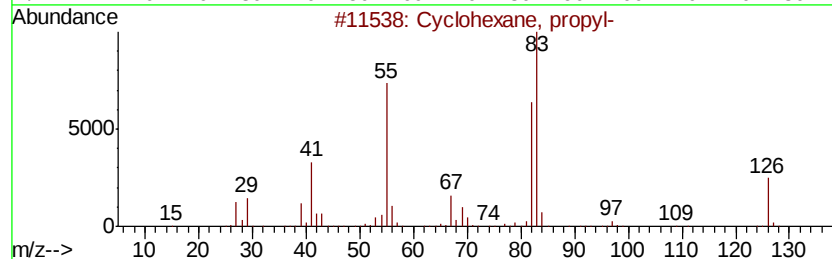
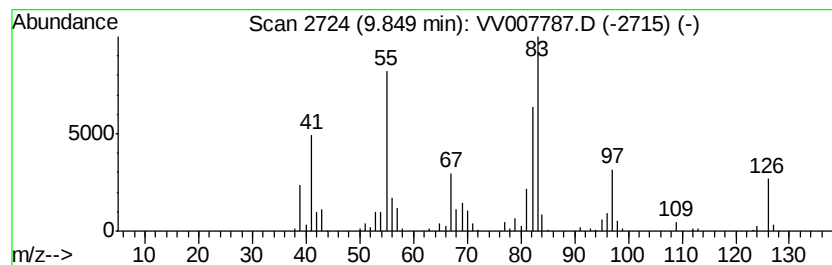
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Cyclic9.85 Concentration Rank 12

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 9.85 | 18.22 ug/L | 180808 | Chlorobenzene-d5 | 8.90 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 93   |
| 2    |    | Cyclohexane, 2-propenyl-      | 124 | C9H16   | 002114-42-3 | 76   |
| 3    |    | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 76   |
| 4    |    | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 76   |
| 5    |    | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

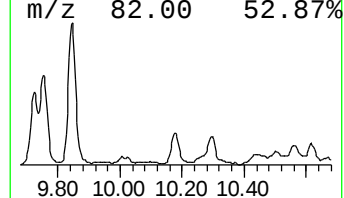
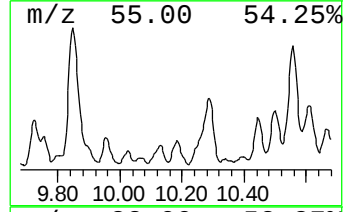
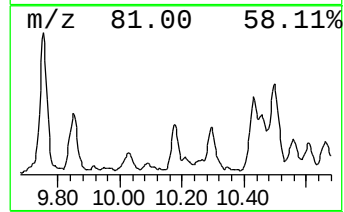
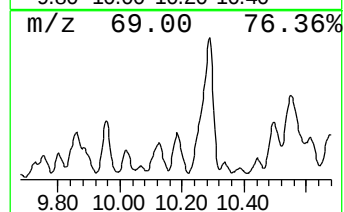
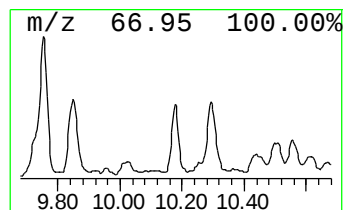
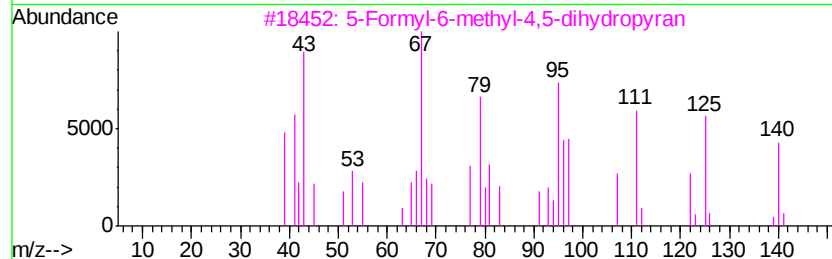
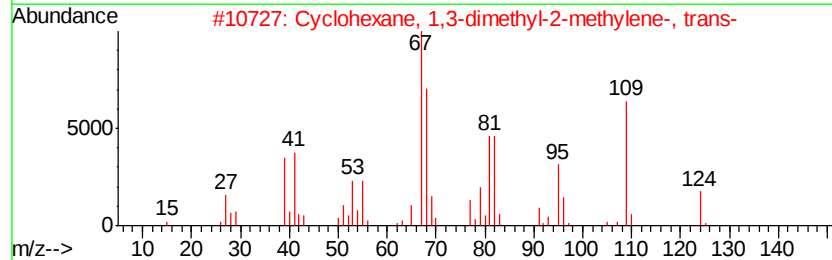
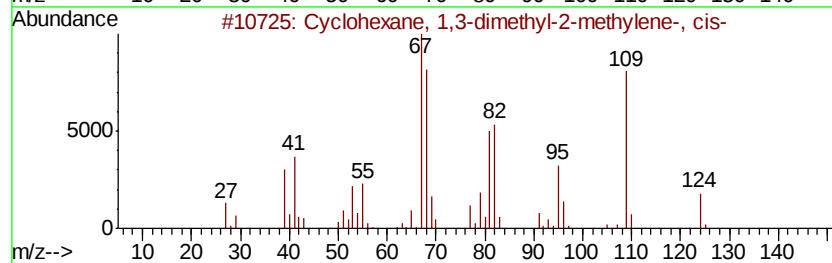
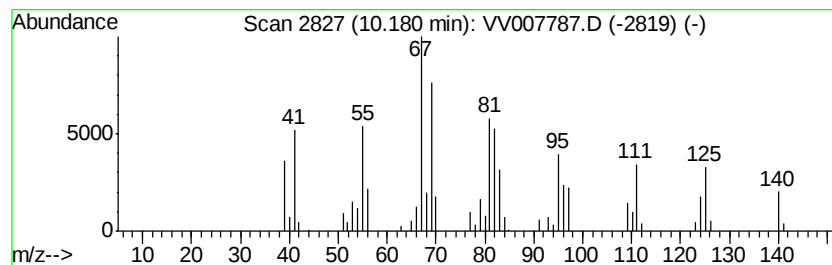
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 unknown-01 Concentration Rank 31

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 10.18 | 4.79 ug/L | 58588 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Cyclohexane, 1,3-dimethyl-2-meth... | 124 | C9H16   | 019781-47-6  | 46   |
| 2    |    |   | Cyclohexane, 1,3-dimethyl-2-meth... | 124 | C9H16   | 020348-74-7  | 46   |
| 3    |    |   | 5-Formyl-6-methyl-4,5-dihydropran   | 140 | C8H12O2 | 1000288-90-6 | 43   |
| 4    |    |   | 4-Nonyne                            | 124 | C9H16   | 020184-91-2  | 38   |
| 5    |    |   | Bicyclo[2.2.1]heptane, 2-methyl-... | 110 | C8H14   | 000872-78-6  | 38   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampled :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

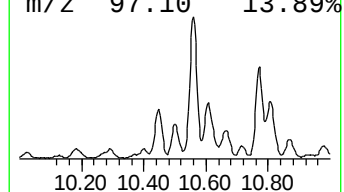
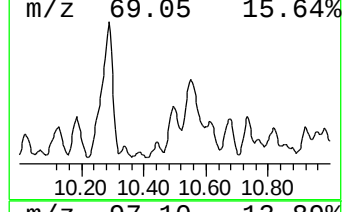
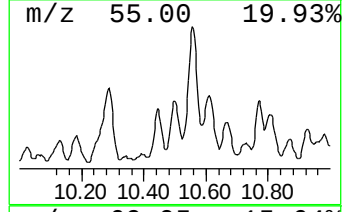
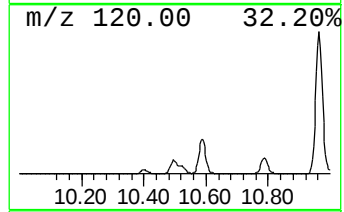
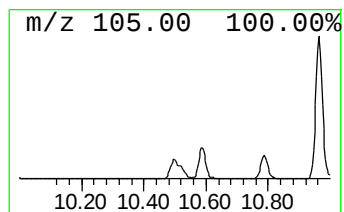
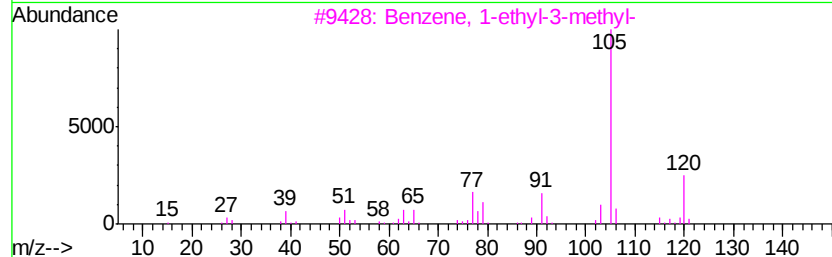
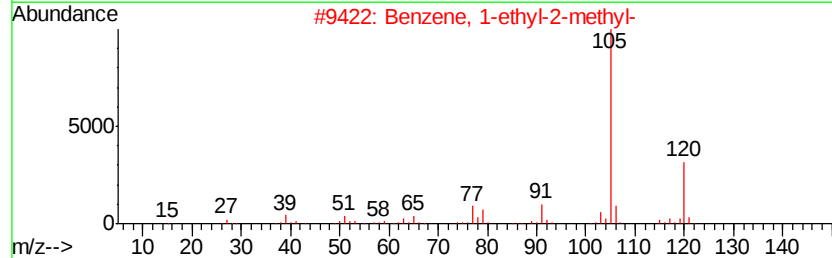
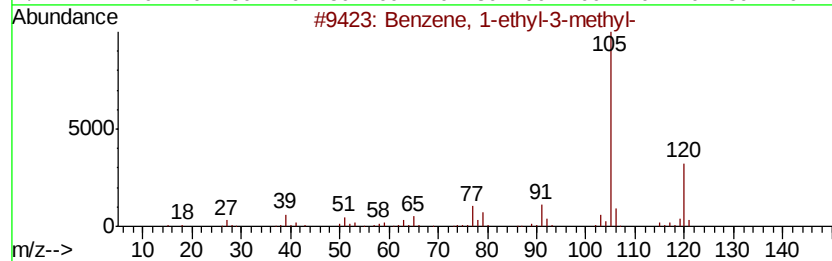
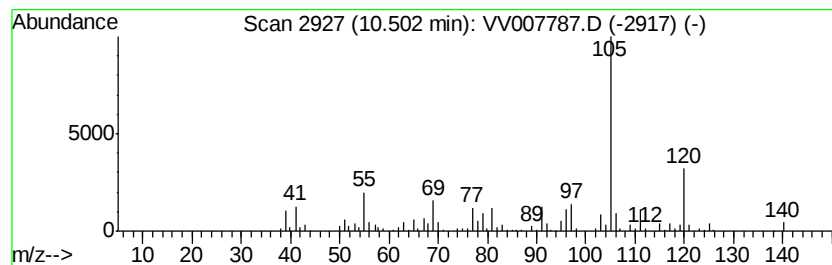
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Benzene, 1-ethyl-3-methyl- Concentration Rank 14

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 10.50 | 16.53 ug/L | 202288 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 93   |
| 2    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 92   |
| 3    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 4    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 90   |
| 5    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 90   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

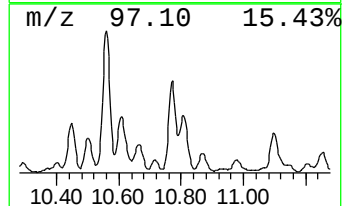
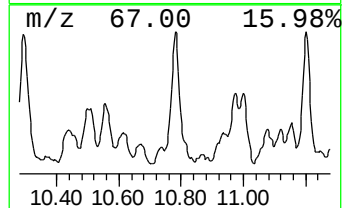
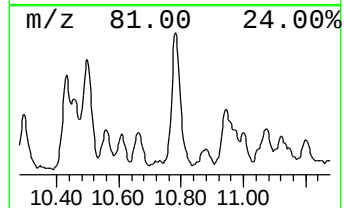
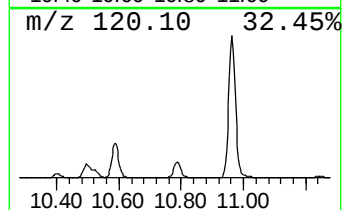
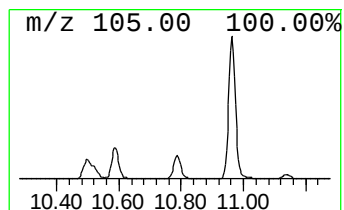
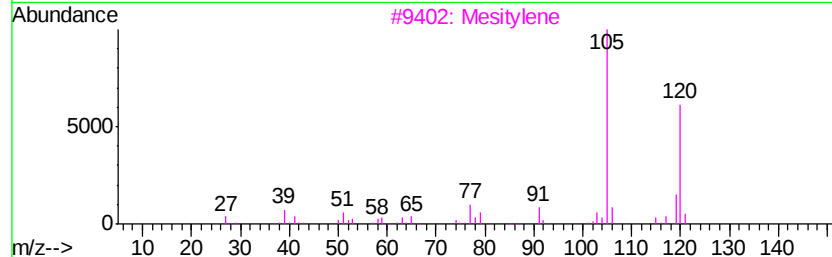
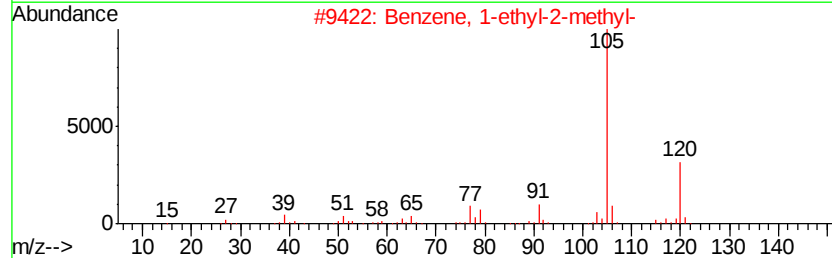
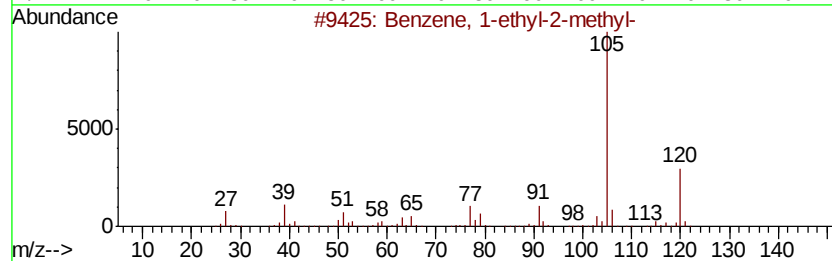
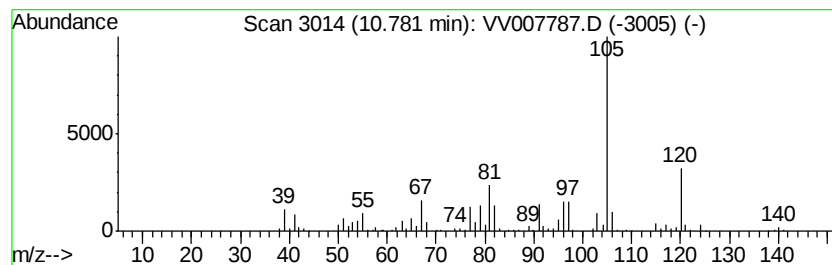
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 10.78 | 19.22 ug/L | 235270 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 92   |
| 2    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 91   |
| 3    |    | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 64   |
| 4    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 64   |
| 5    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

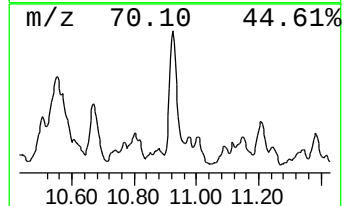
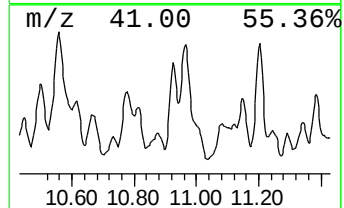
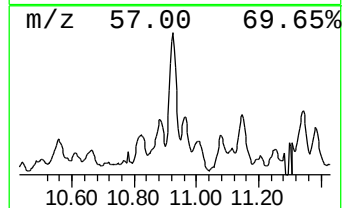
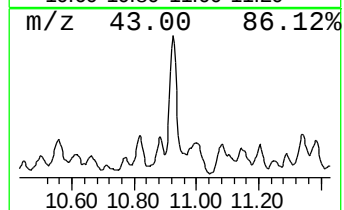
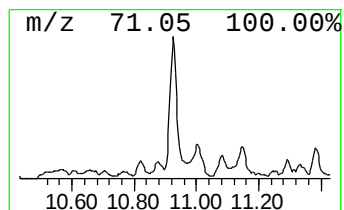
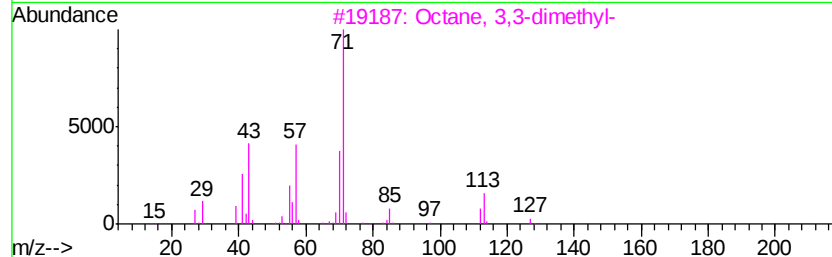
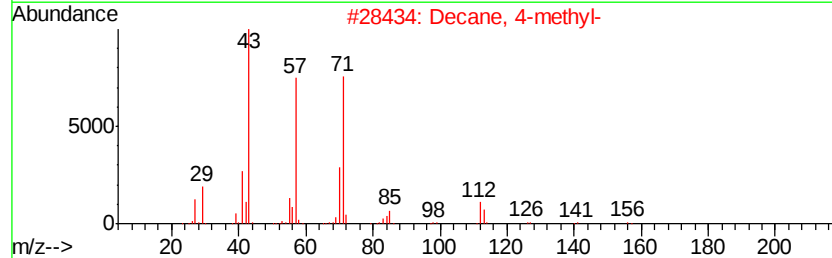
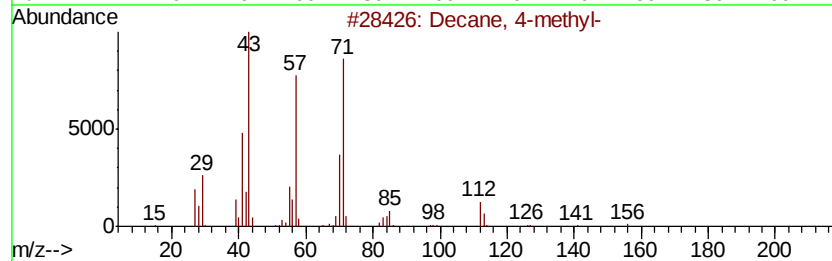
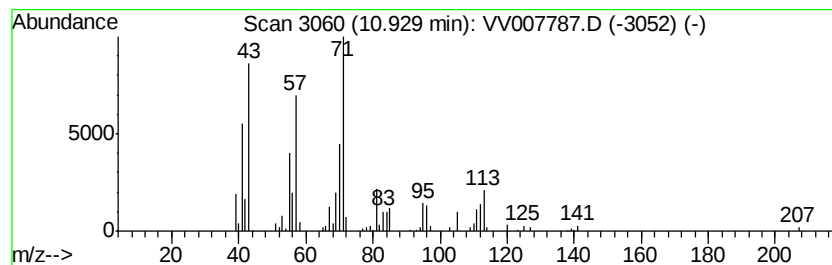
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 30

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 10.93 | 4.94 ug/L | 60414 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 4-methyl-     | 156 | C11H24  | 002847-72-5 | 76   |
| 2    |    |   | Decane, 4-methyl-     | 156 | C11H24  | 002847-72-5 | 64   |
| 3    |    |   | Octane, 3,3-dimethyl- | 142 | C10H22  | 004110-44-5 | 58   |
| 4    |    |   | Octane, 3,3-dimethyl- | 142 | C10H22  | 004110-44-5 | 53   |
| 5    |    |   | Nonane, 2,6-dimethyl- | 156 | C11H24  | 017302-28-2 | 52   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

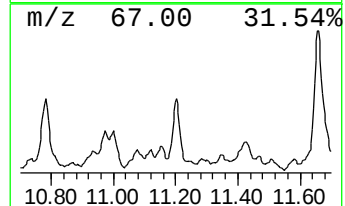
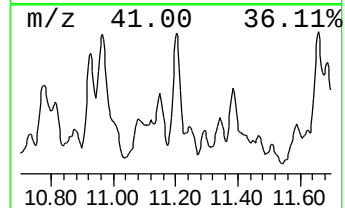
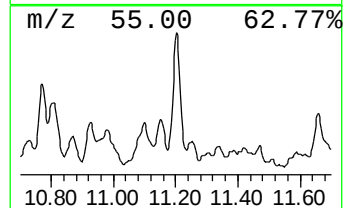
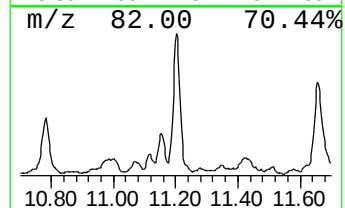
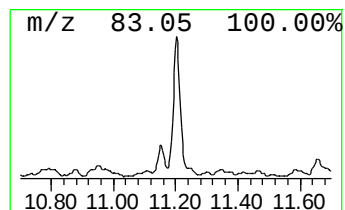
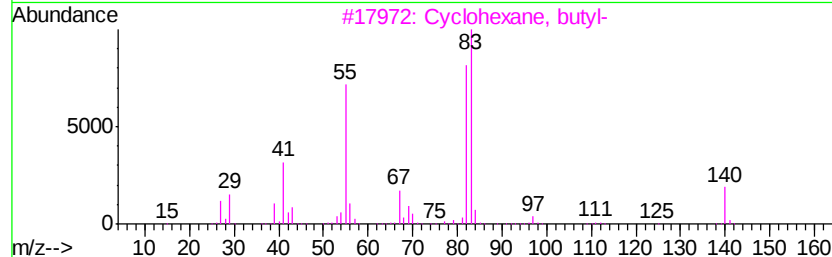
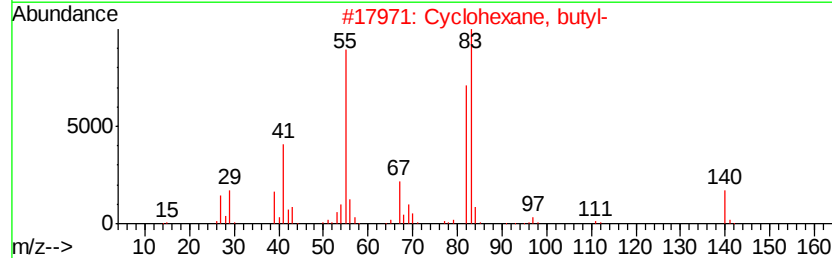
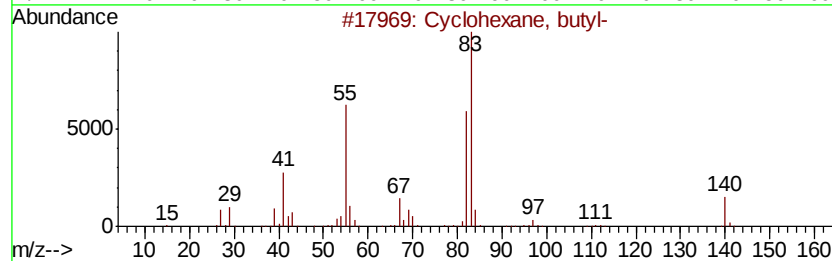
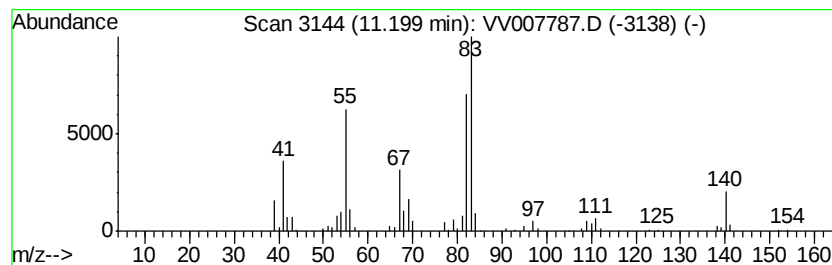
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Cyclic11.20 Concentration Rank 26

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 11.20 | 6.63 ug/L | 81156 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, butyl-   | 140 | C10H20  | 001678-93-9 | 93   |
| 2    |    |   | Cyclohexane, butyl-   | 140 | C10H20  | 001678-93-9 | 90   |
| 3    |    |   | Cyclohexane, butyl-   | 140 | C10H20  | 001678-93-9 | 87   |
| 4    |    |   | Cyclohexane, octyl-   | 196 | C14H28  | 001795-15-9 | 64   |
| 5    |    |   | Cyclohexane, undecyl- | 238 | C17H34  | 054105-66-7 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

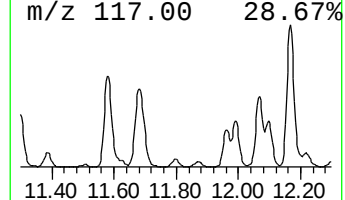
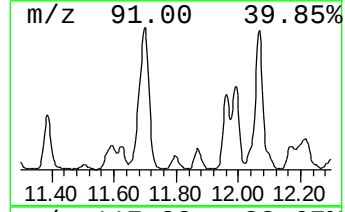
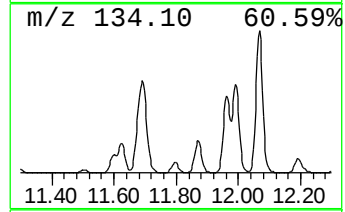
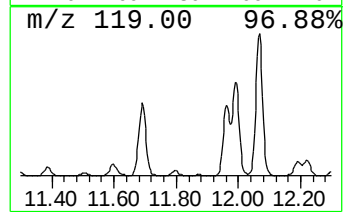
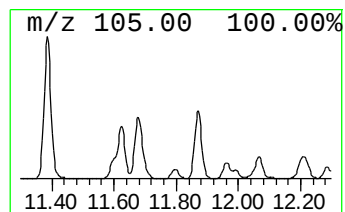
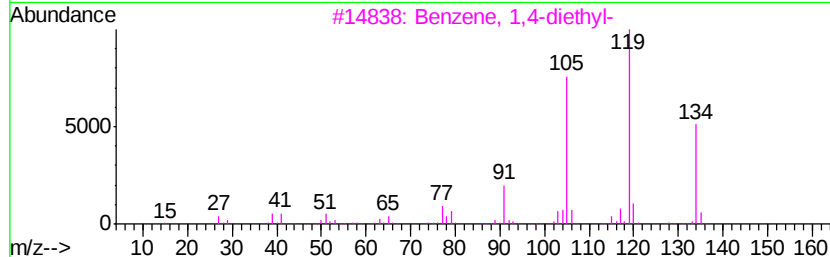
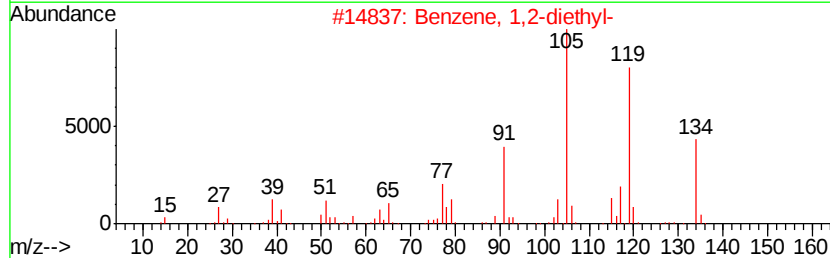
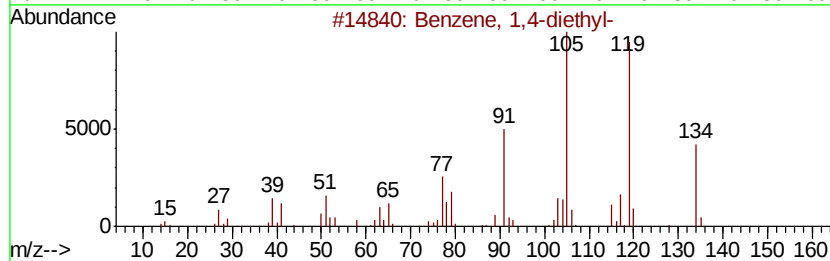
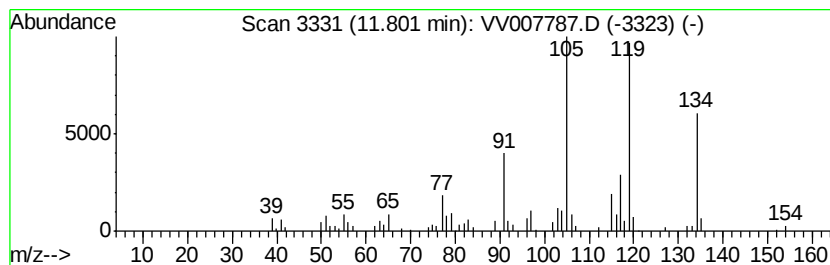
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Benzene, 1,4-diethyl- Concentration Rank 34

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 11.80 | 4.55 ug/L | 55705 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,4-diethyl- | 134 | C10H14  | 000105-05-5 | 93   |
| 2    |    |   | Benzene, 1,2-diethyl- | 134 | C10H14  | 000135-01-3 | 93   |
| 3    |    |   | Benzene, 1,4-diethyl- | 134 | C10H14  | 000105-05-5 | 87   |
| 4    |    |   | Benzene, 1,2-diethyl- | 134 | C10H14  | 000135-01-3 | 87   |
| 5    |    |   | Benzene, 1,4-diethyl- | 134 | C10H14  | 000105-05-5 | 87   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

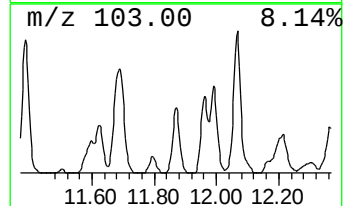
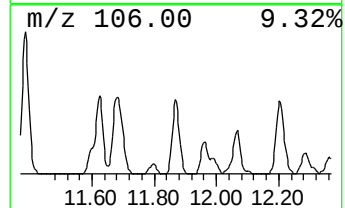
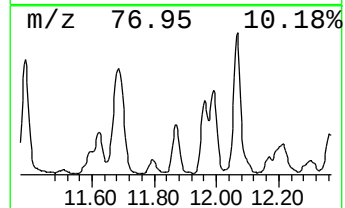
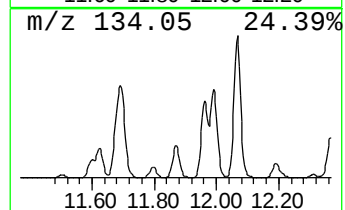
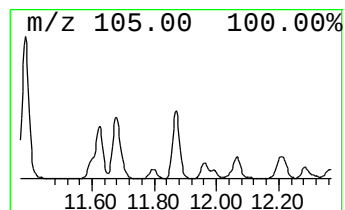
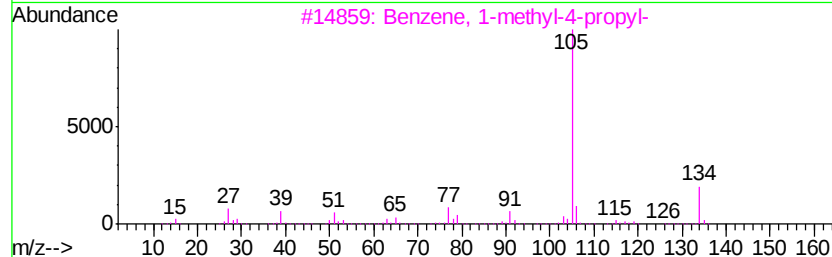
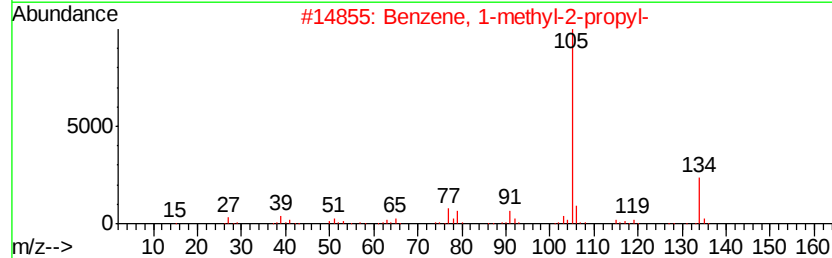
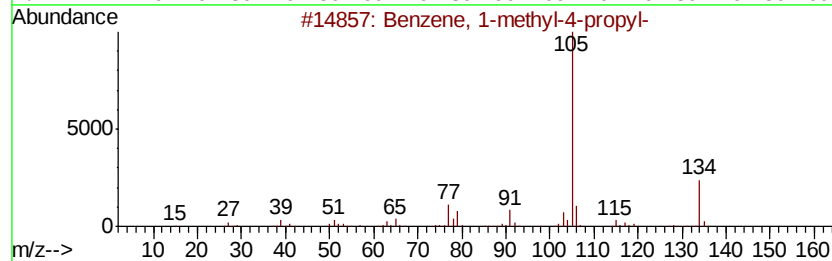
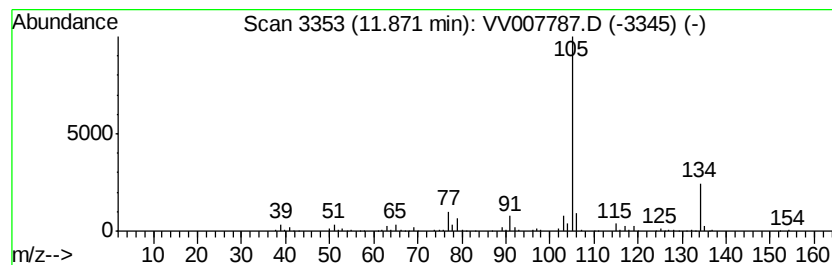
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 Benzene, 1-methyl-4-propyl- Concentration Rank 16

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 11.87 | 15.50 ug/L | 189682 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 94   |
| 2    |    | Benzene, 1-methyl-2-propyl- | 134 | C10H14  | 001074-17-5 | 91   |
| 3    |    | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 91   |
| 4    |    | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 91   |
| 5    |    | Benzene, (1-methylpropyl)-  | 134 | C10H14  | 000135-98-8 | 90   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

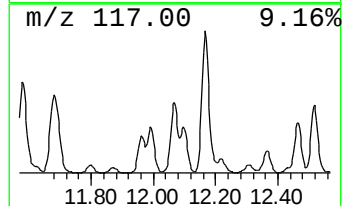
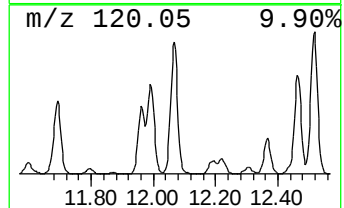
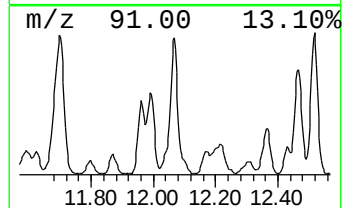
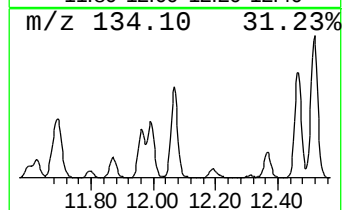
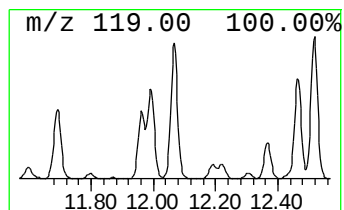
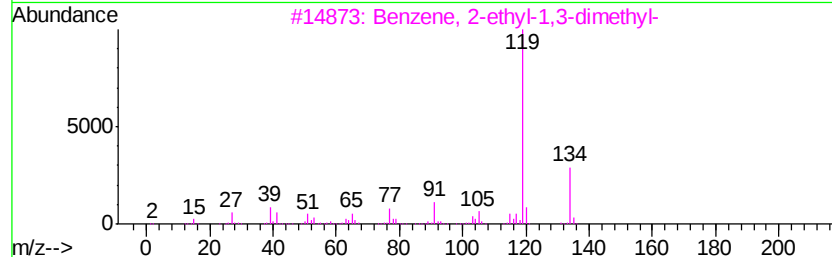
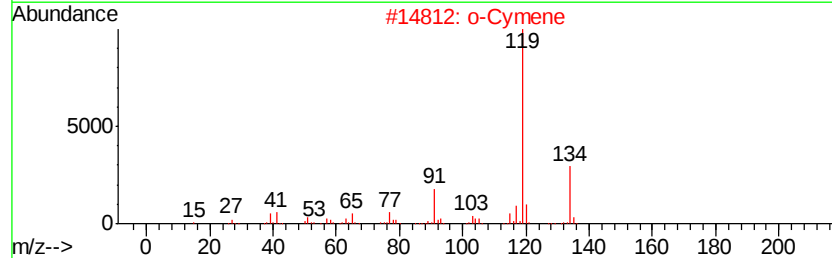
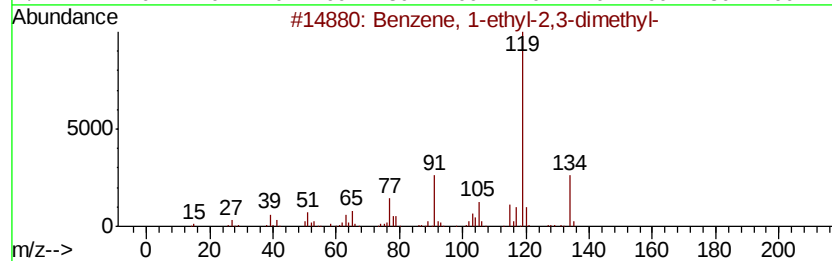
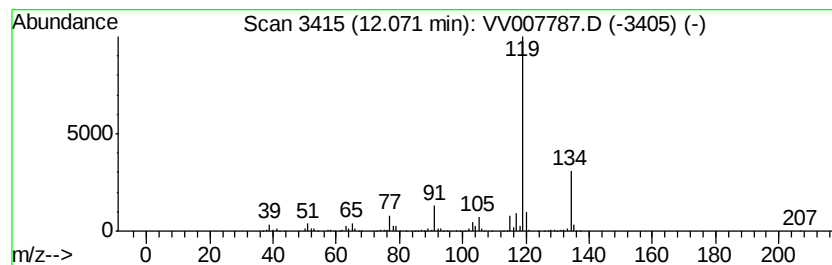
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 12.07 | 59.38 ug/L | 726853 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 96   |
| 2    |    | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 95   |
| 3    |    | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 95   |
| 4    |    | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 95   |
| 5    |    | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

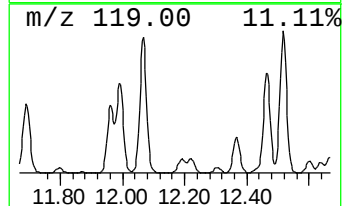
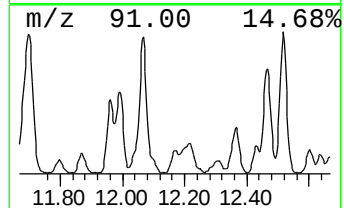
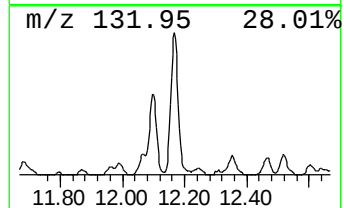
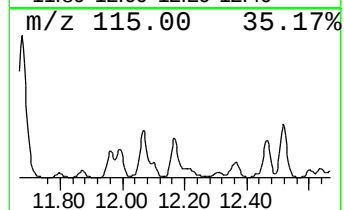
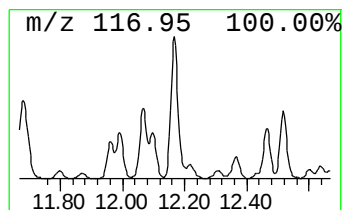
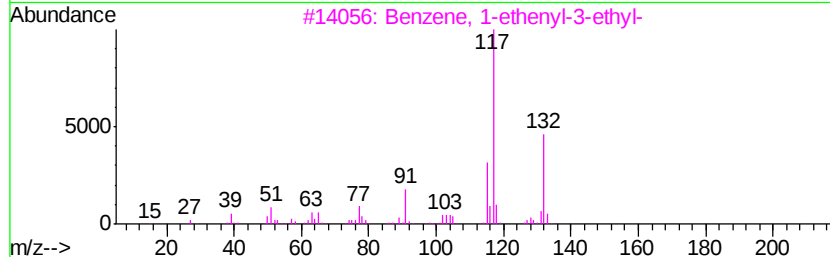
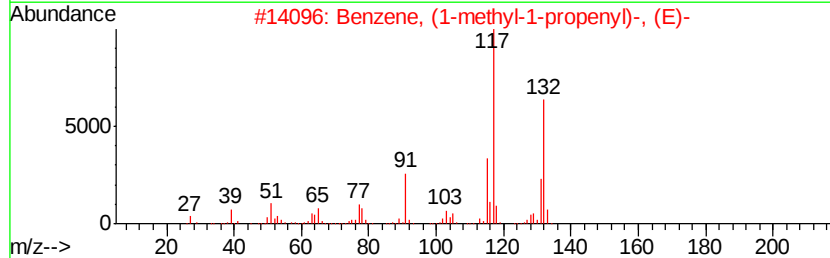
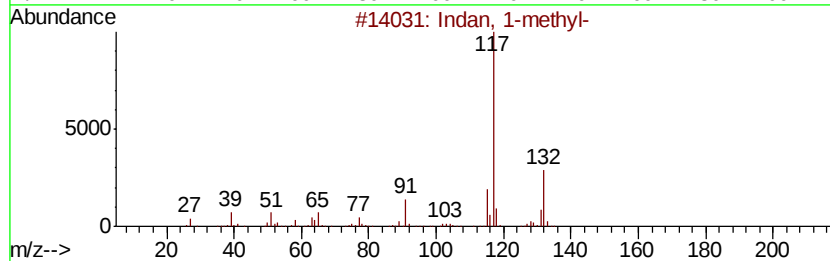
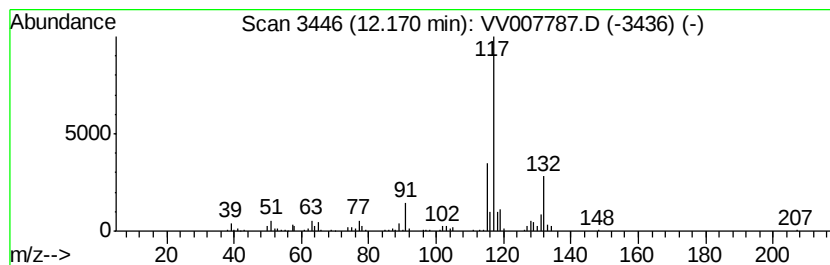
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 23 Indan, 1-methyl- Concentration Rank 19

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 12.17 | 11.83 ug/L | 144755 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Indan, 1-methyl-                     | 132 | C10H12  | 000767-58-8 | 87   |
| 2    |    | Benzene, (1-methyl-1-propenyl)-, ... | 132 | C10H12  | 000768-00-3 | 87   |
| 3    |    | Benzene, 1-ethenyl-3-ethyl-          | 132 | C10H12  | 007525-62-4 | 87   |
| 4    |    | 1-Phenyl-1-butene                    | 132 | C10H12  | 000824-90-8 | 87   |
| 5    |    | Benzene, (2-methyl-2-propenyl)-      | 132 | C10H12  | 003290-53-7 | 86   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

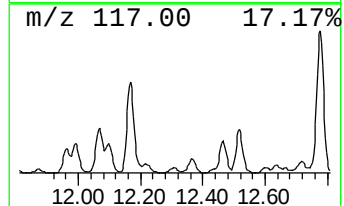
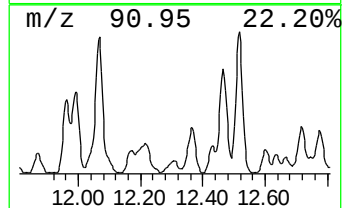
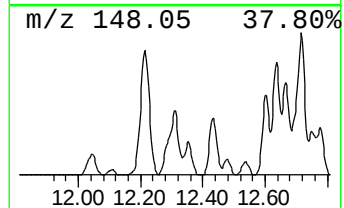
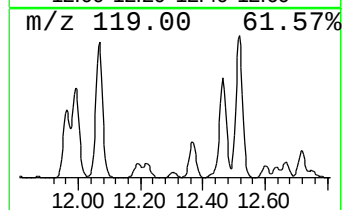
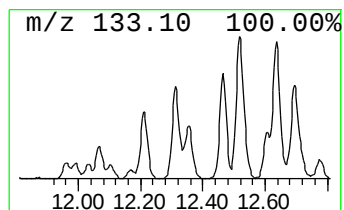
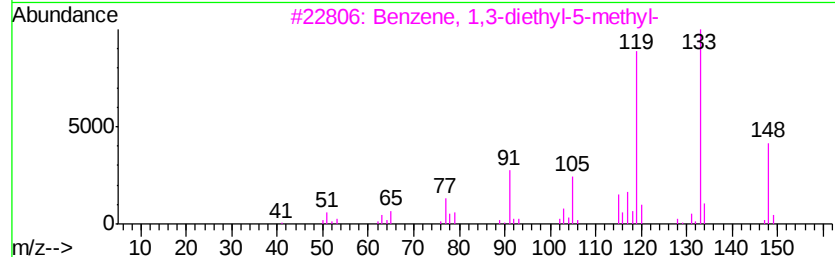
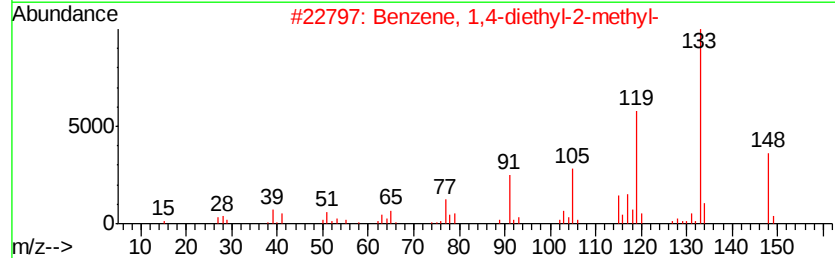
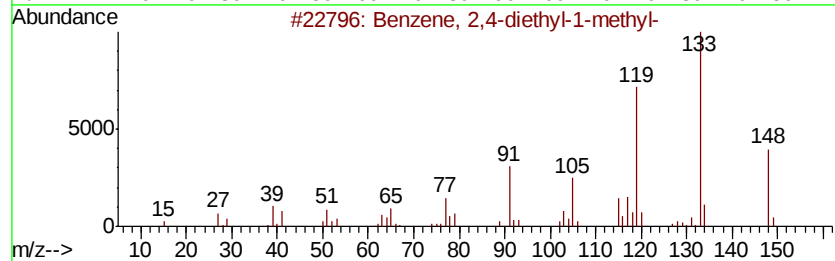
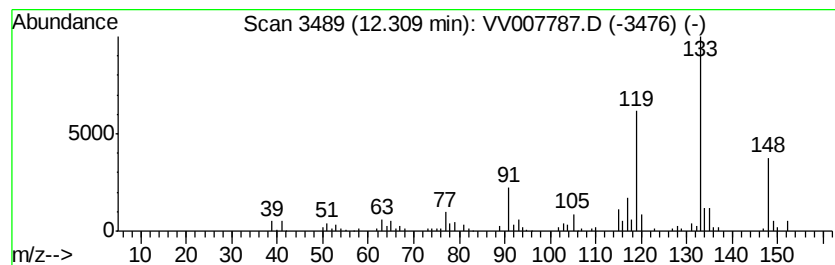
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 Benzene, 2,4-diethyl-1-methyl- Concentration Rank 25

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 12.31 | 7.60 ug/L | 93085 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2,4-diethyl-1-methyl- | 148 | C11H16  | 001758-85-6 | 93   |
| 2    |    |   | Benzene, 1,4-diethyl-2-methyl- | 148 | C11H16  | 013632-94-5 | 90   |
| 3    |    |   | Benzene, 1,3-diethyl-5-methyl- | 148 | C11H16  | 002050-24-0 | 70   |
| 4    |    |   | Benzene, 1,3-diethyl-5-methyl- | 148 | C11H16  | 002050-24-0 | 70   |
| 5    |    |   | Benzene, pentamethyl-          | 148 | C11H16  | 000700-12-9 | 70   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

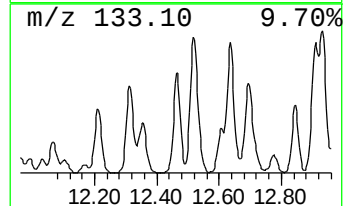
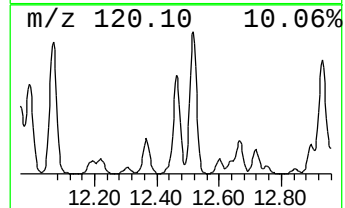
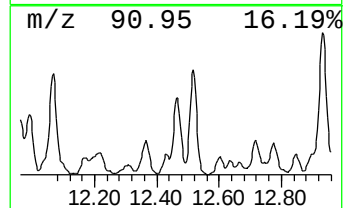
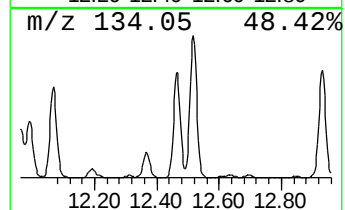
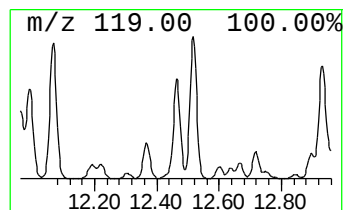
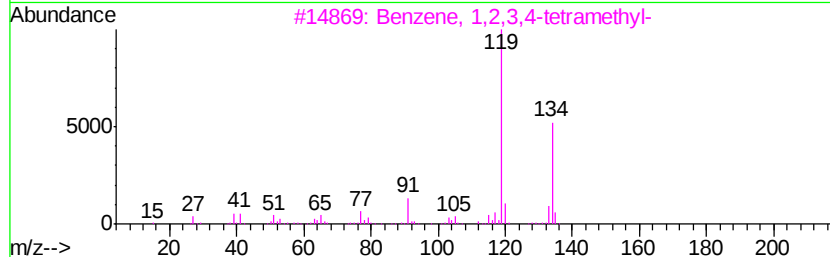
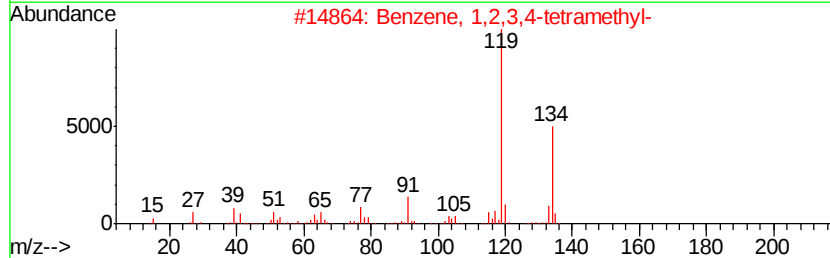
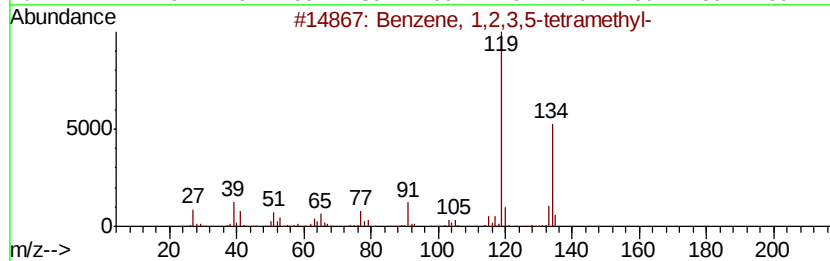
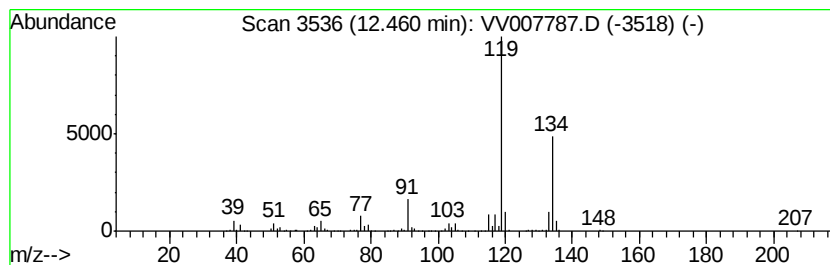
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 12.46 | 50.72 ug/L | 620898 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 97   |
| 2    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 97   |
| 3    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 97   |
| 4    |    | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 96   |
| 5    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

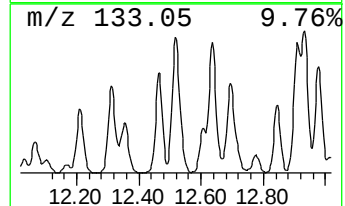
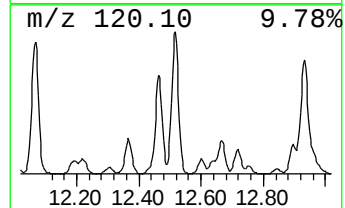
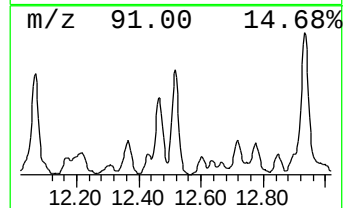
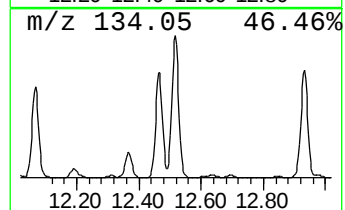
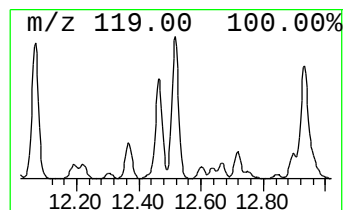
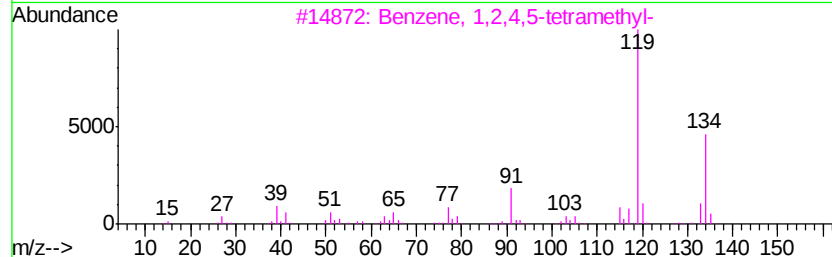
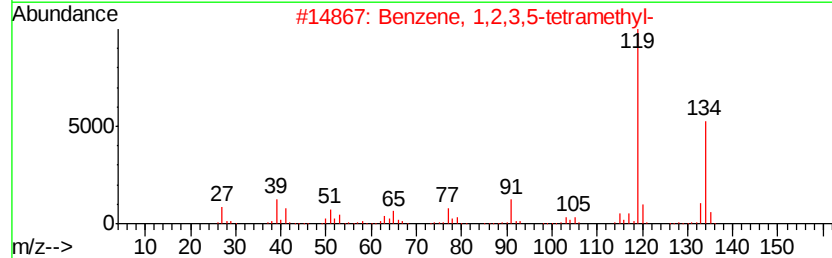
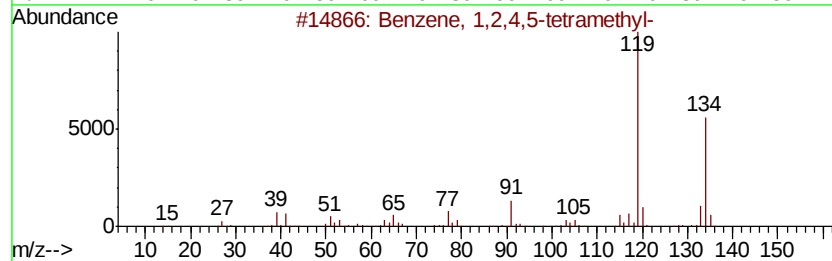
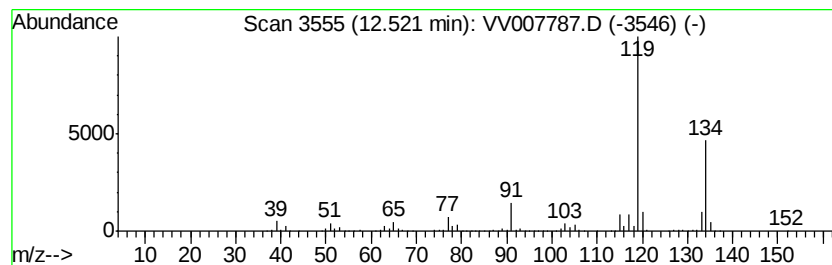
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 12.52 | 61.68 ug/L | 754985 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 95   |
| 2    |    | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 95   |
| 3    |    | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 95   |
| 4    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 95   |
| 5    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

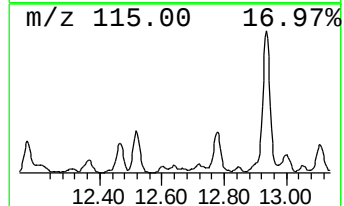
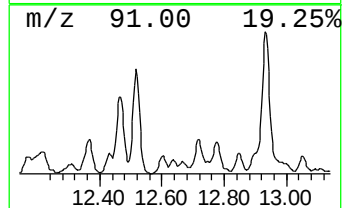
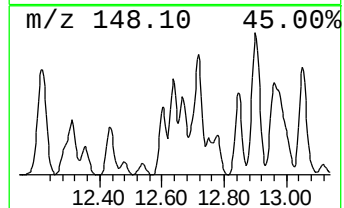
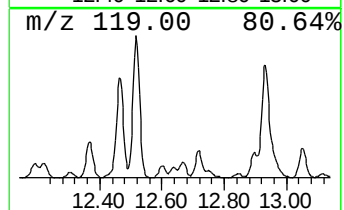
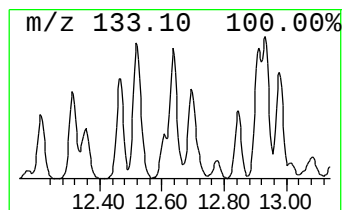
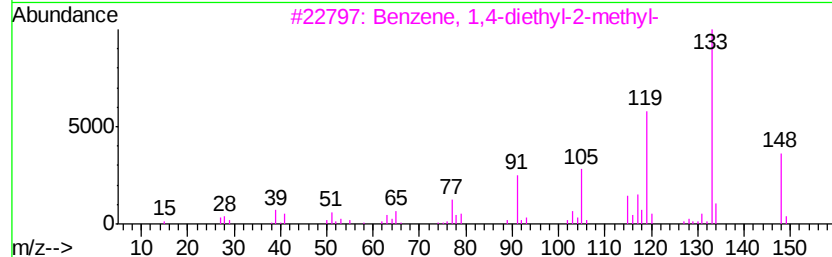
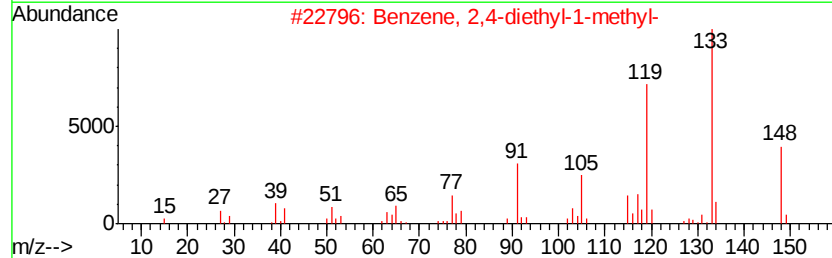
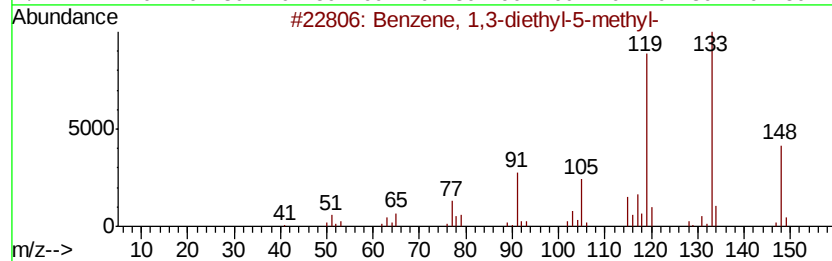
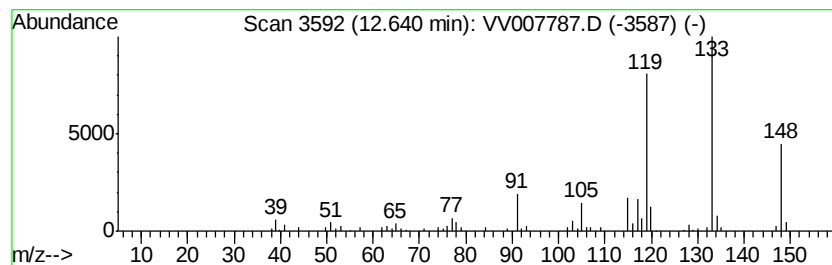
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 42

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 12.64 | 2.54 ug/L | 31080 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,3-diethyl-5-methyl-    | 148 | C11H16  | 002050-24-0 | 94   |
| 2    |    |   | Benzene, 2,4-diethyl-1-methyl-    | 148 | C11H16  | 001758-85-6 | 90   |
| 3    |    |   | Benzene, 1,4-diethyl-2-methyl-    | 148 | C11H16  | 013632-94-5 | 86   |
| 4    |    |   | Benzene, 1,3-diethyl-5-methyl-    | 148 | C11H16  | 002050-24-0 | 81   |
| 5    |    |   | Benzene, 1-ethyl-2,4,5-trimethyl- | 148 | C11H16  | 017851-27-3 | 58   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

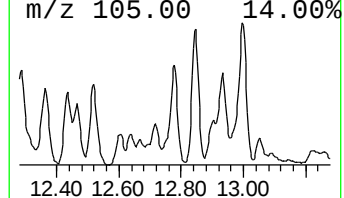
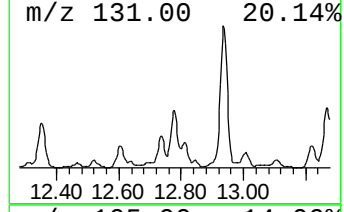
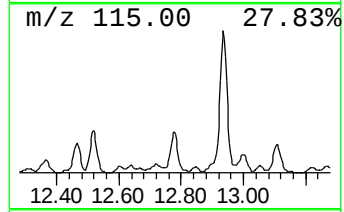
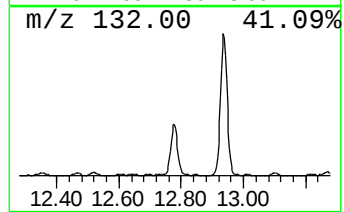
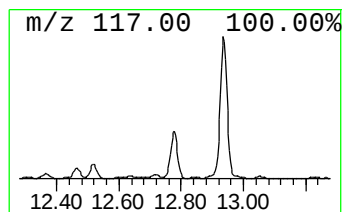
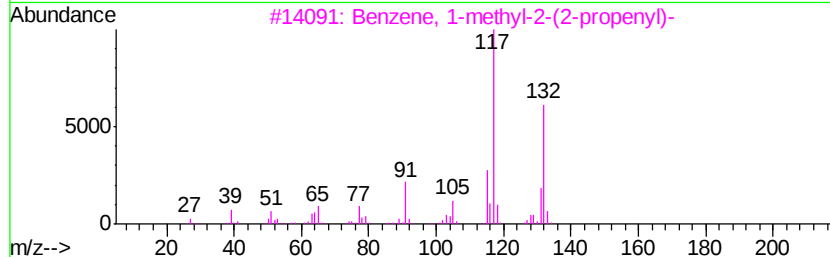
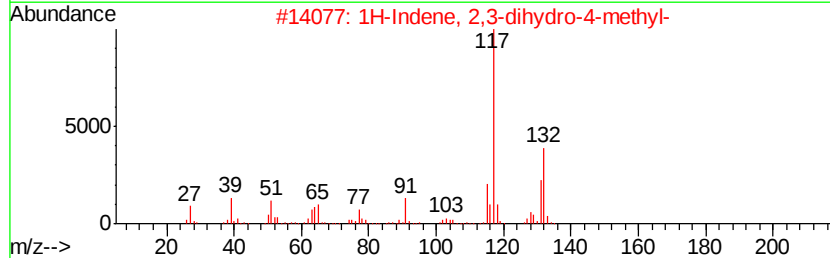
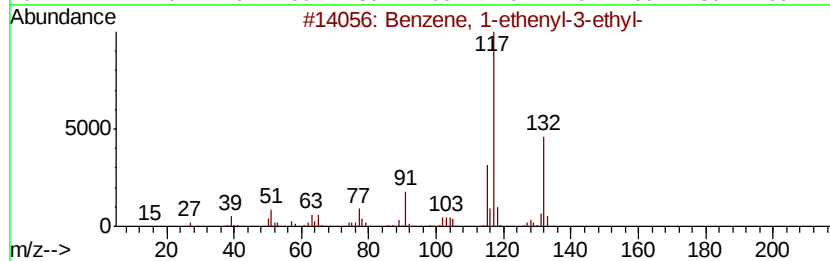
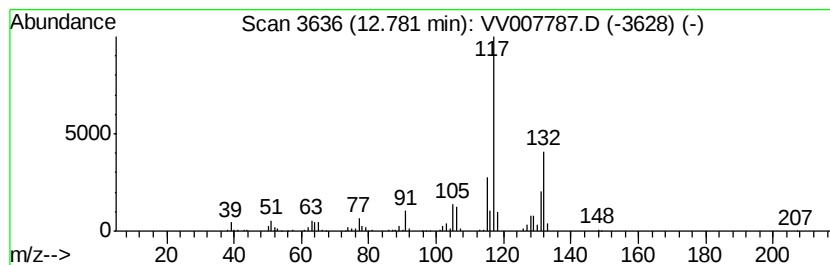
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 12.78 | 21.60 ug/L | 264455 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethenyl-3-ethyl-       | 132 | C10H12  | 007525-62-4 | 87   |
| 2    |    | 1H-Indene, 2,3-dihydro-4-methyl-  | 132 | C10H12  | 000824-22-6 | 87   |
| 3    |    | Benzene, 1-methyl-2-(2-propenyl)- | 132 | C10H12  | 001587-04-8 | 87   |
| 4    |    | 3-Phenylbut-1-ene                 | 132 | C10H12  | 000934-10-1 | 86   |
| 5    |    | Indan, 1-methyl-                  | 132 | C10H12  | 000767-58-8 | 83   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

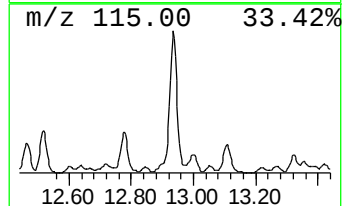
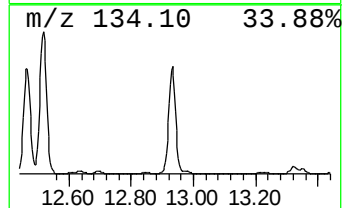
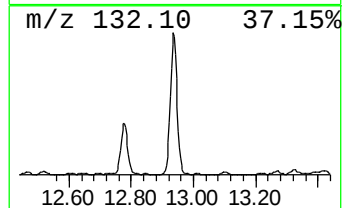
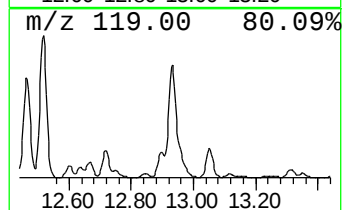
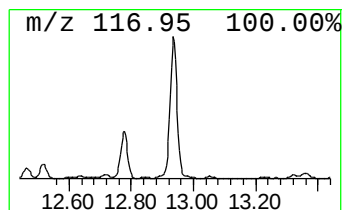
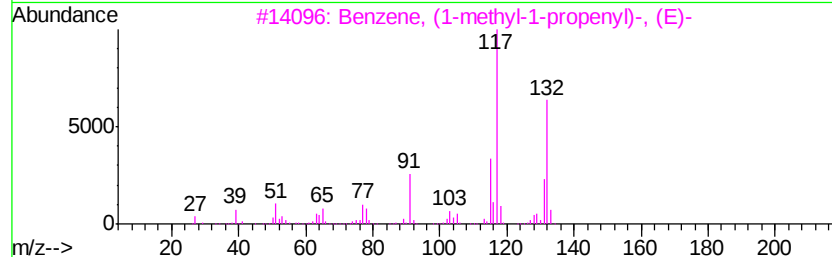
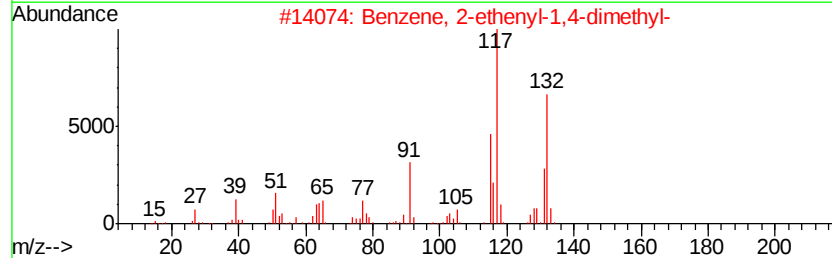
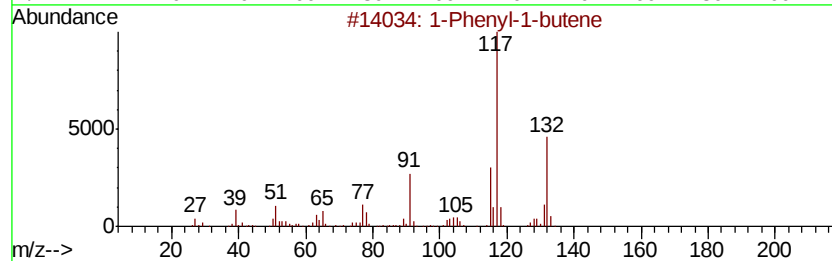
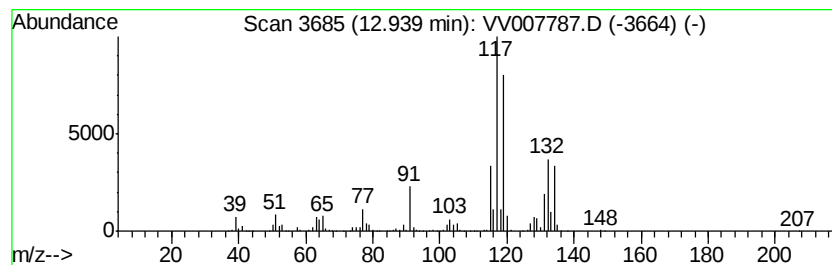
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 1-Phenyl-1-butene Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 12.94 | 122.35 ug/L | 1497620 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1-Phenyl-1-butene                   | 132 | C10H12  | 000824-90-8 | 89   |
| 2    |    | Benzene, 2-ethenyl-1,4-dimethyl-    | 132 | C10H12  | 002039-89-6 | 83   |
| 3    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 64   |
| 4    |    | Benzene, 2-butenyl-                 | 132 | C10H12  | 001560-06-1 | 64   |
| 5    |    | Benzene, 1-ethenyl-3-ethyl-         | 132 | C10H12  | 007525-62-4 | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

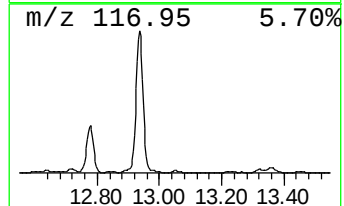
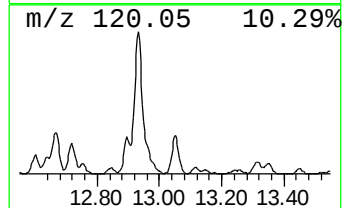
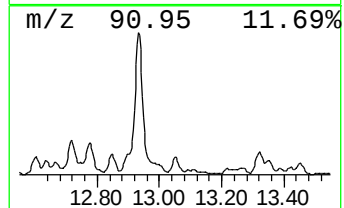
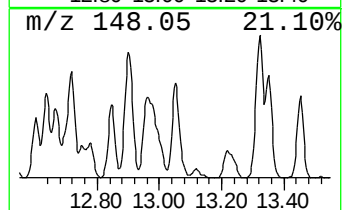
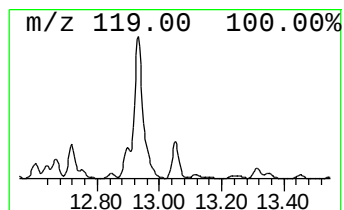
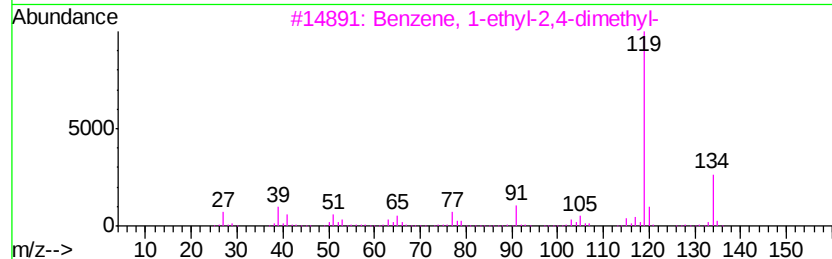
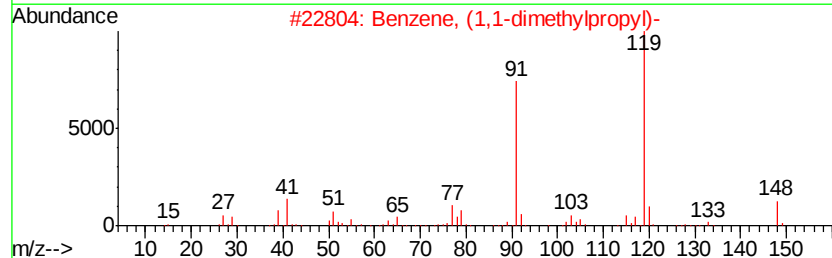
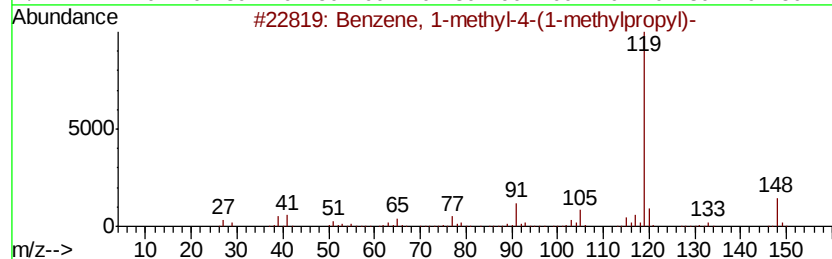
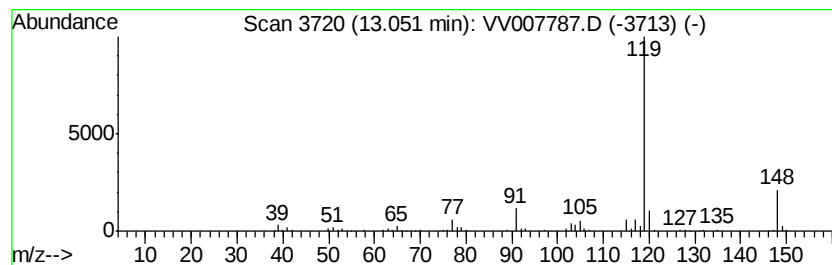
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 Benzene, 1-methyl-4-(1-meth... Concentration Rank 24

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 13.05 | 8.21 ug/L | 100448 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16  | 001595-16-0 | 91   |
| 2    |    | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8 | 83   |
| 3    |    | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9 | 80   |
| 4    |    | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9 | 80   |
| 5    |    | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

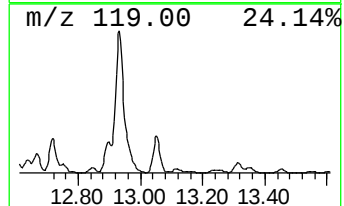
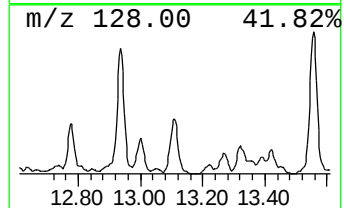
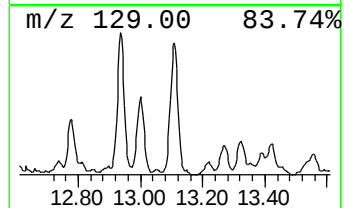
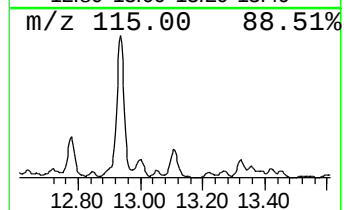
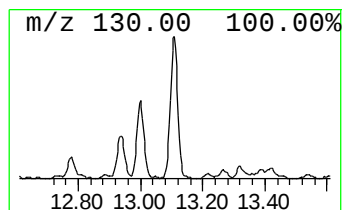
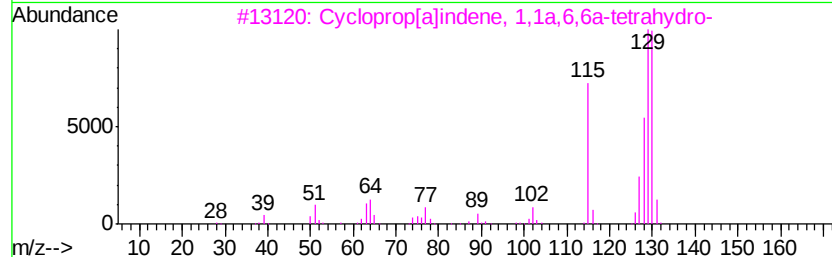
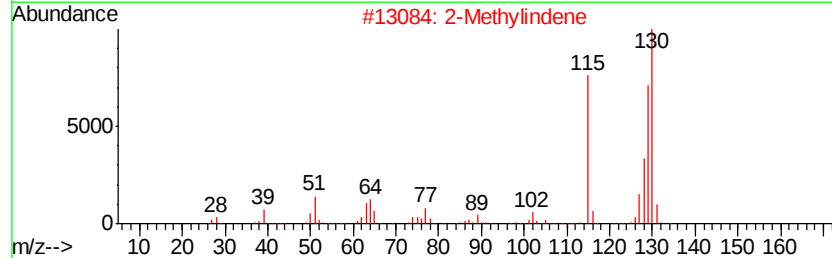
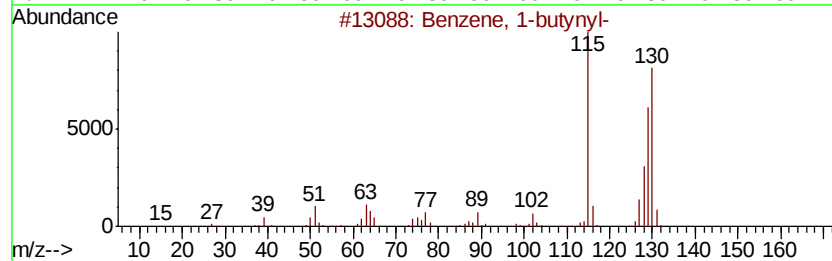
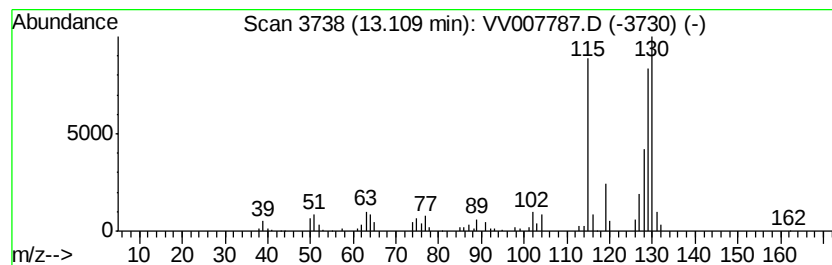
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 Benzene, 1-butynyl- Concentration Rank 21

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 13.11 | 9.96 ug/L | 121909 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-butynyl-                 | 130 | C10H10  | 000622-76-4 | 94   |
| 2    |    | 2-Methylindene                      | 130 | C10H10  | 002177-47-1 | 93   |
| 3    |    | Cycloprop[alindene, 1,1a,6,6a-te... | 130 | C10H10  | 015677-15-3 | 91   |
| 4    |    | Naphthalene, 1,2-dihydro-           | 130 | C10H10  | 000447-53-0 | 90   |
| 5    |    | 1H-Indene, 1-methyl-                | 130 | C10H10  | 000767-59-9 | 90   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

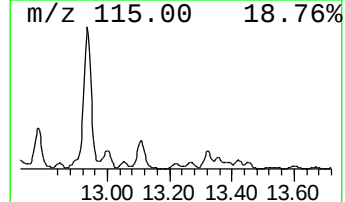
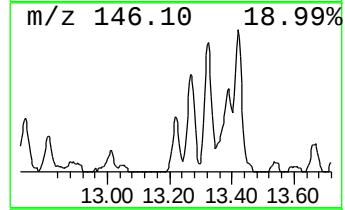
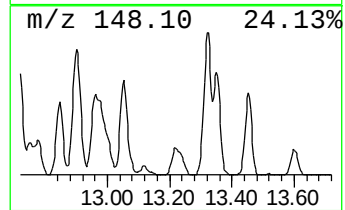
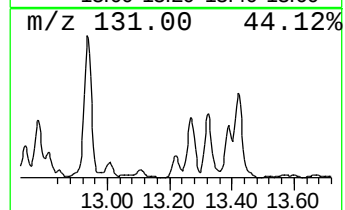
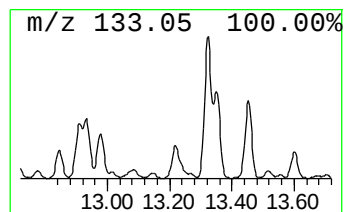
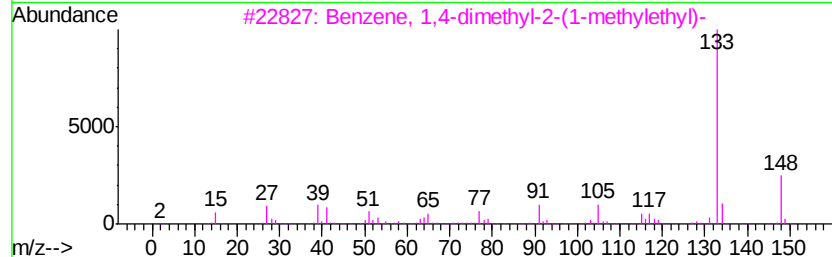
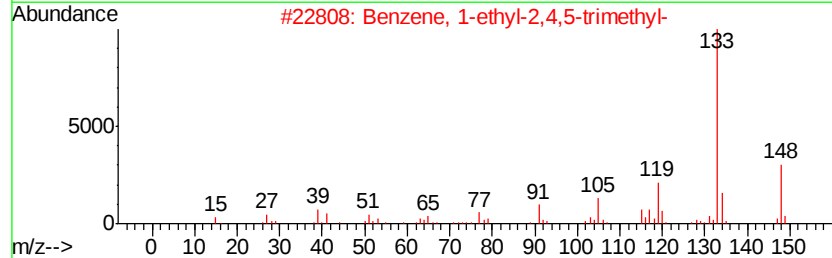
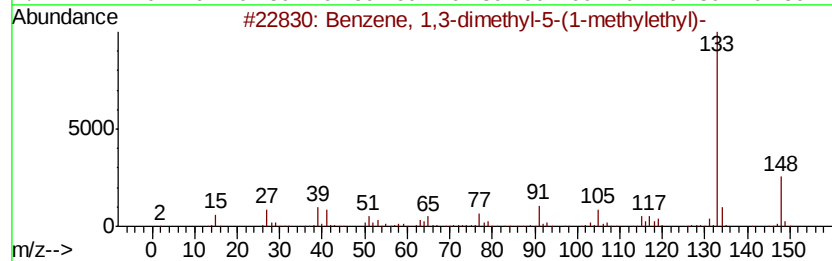
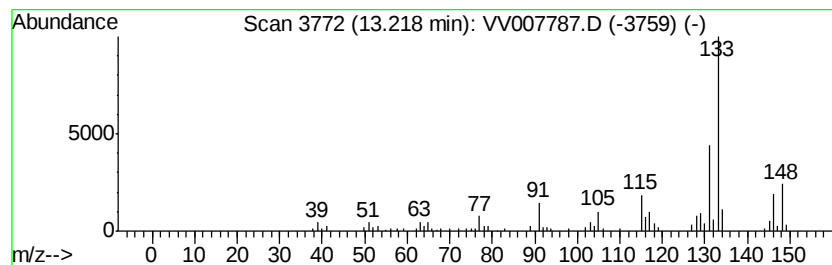
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 Benzene, 1,3-dimethyl-5-(1-... Concentration Rank 29

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 13.22 | 5.69 ug/L | 69661 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Benzene, 1,3-dimethyl-5-(1-methy... | 148 | C11H16  | 004706-90-5  | 64   |
| 2    |    |   | Benzene, 1-ethyl-2,4,5-trimethyl-   | 148 | C11H16  | 017851-27-3  | 58   |
| 3    |    |   | Benzene, 1,4-dimethyl-2-(1-methv... | 148 | C11H16  | 004132-72-3  | 53   |
| 4    |    |   | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9  | 52   |
| 5    |    |   | 3,4-Dimethylcumene                  | 148 | C11H16  | 1000370-34-1 | 52   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

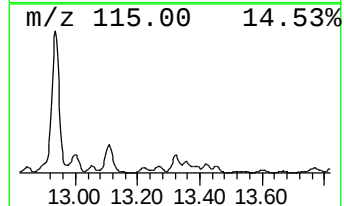
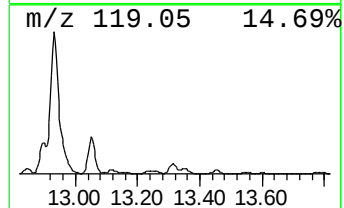
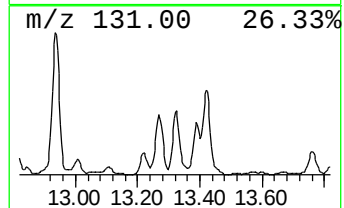
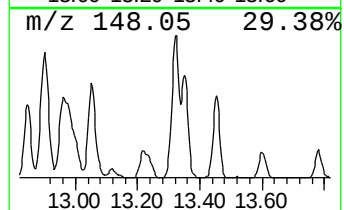
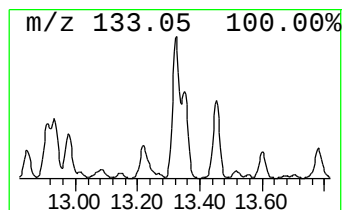
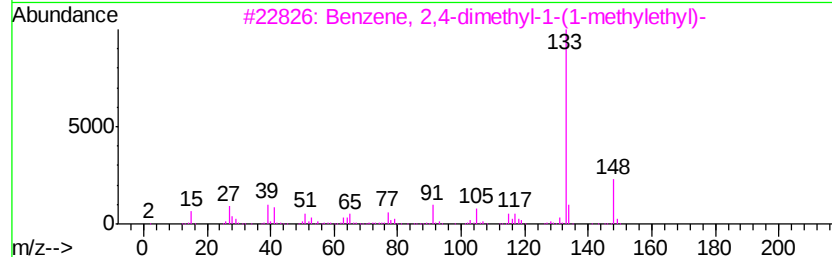
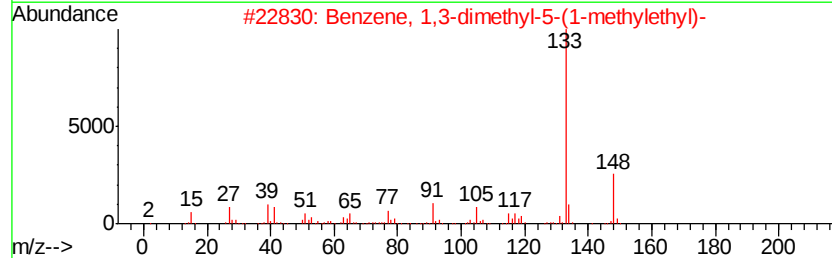
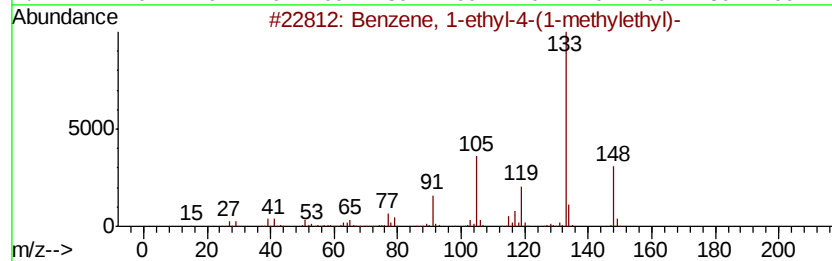
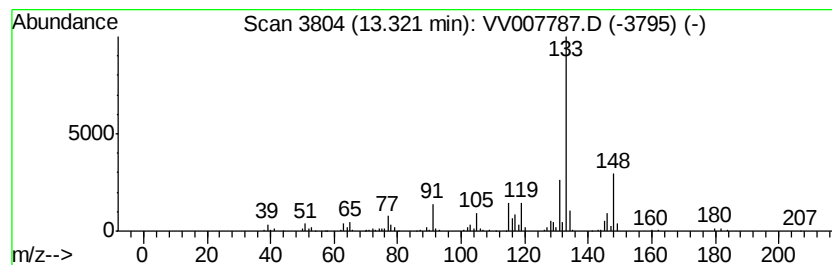
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 36 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 13.32 | 17.51 ug/L | 214386 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 | C11H16  | 004218-48-8 | 76   |
| 2    |    | Benzene, 1,3-dimethyl-5-(1-methy... | 148 | C11H16  | 004706-90-5 | 70   |
| 3    |    | Benzene, 2,4-dimethyl-1-(1-methv... | 148 | C11H16  | 004706-89-2 | 70   |
| 4    |    | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9 | 70   |
| 5    |    | Benzene, 1-ethyl-3-(1-methylethyl)- | 148 | C11H16  | 004920-99-4 | 68   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

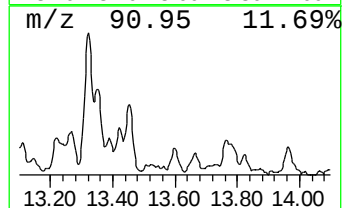
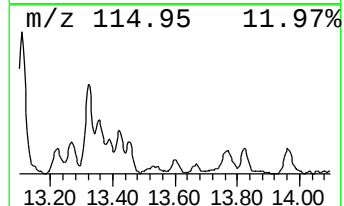
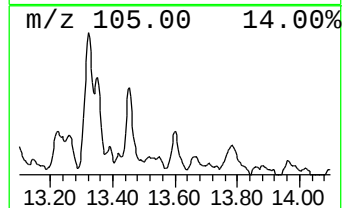
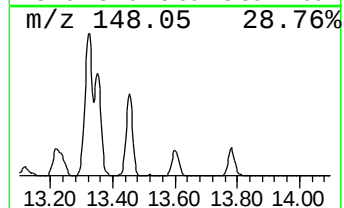
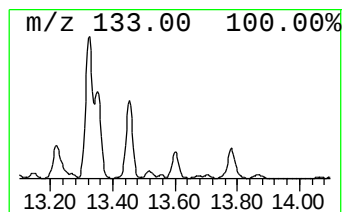
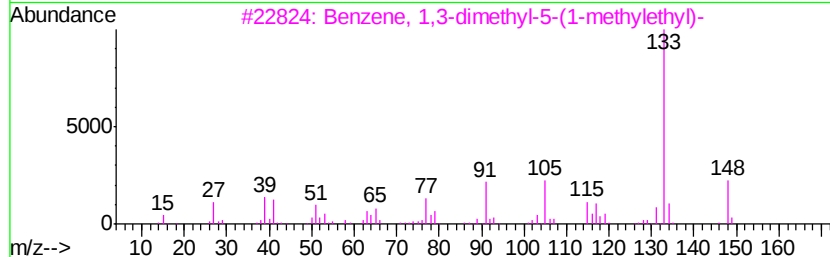
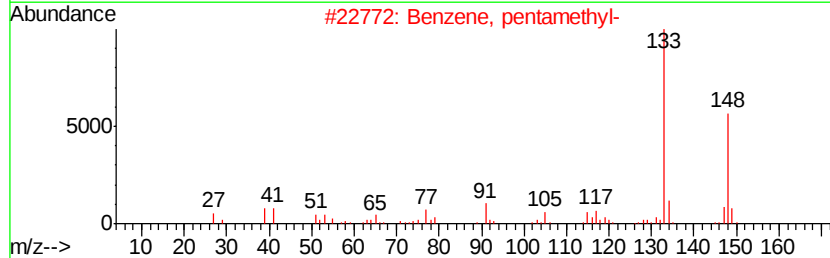
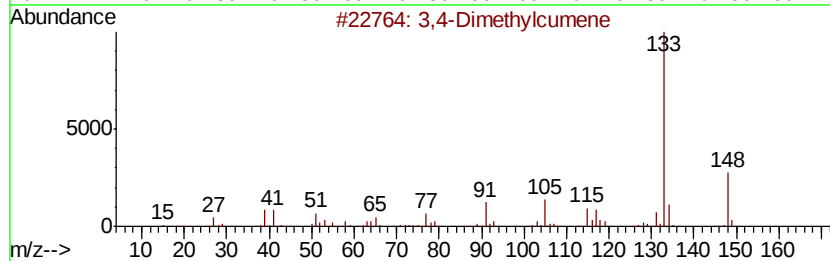
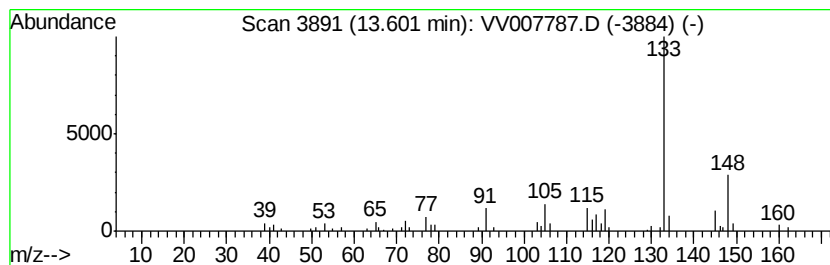
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 37 3,4-Dimethylcumene Concentration Rank 39

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 13.60 | 3.10 ug/L | 37938 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 3,4-Dimethylcumene                  | 148 | C11H16  | 1000370-34-1 | 87   |
| 2    |    | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9  | 83   |
| 3    |    | Benzene, 1,3-dimethyl-5-(1-methv... | 148 | C11H16  | 004706-90-5  | 83   |
| 4    |    | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9  | 83   |
| 5    |    | 2-tert-Butyltoluene                 | 148 | C11H16  | 001074-92-6  | 81   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

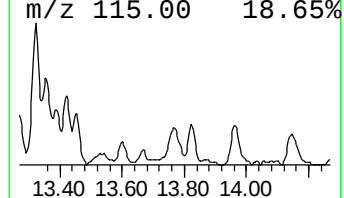
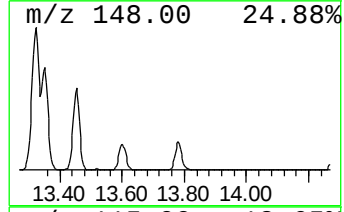
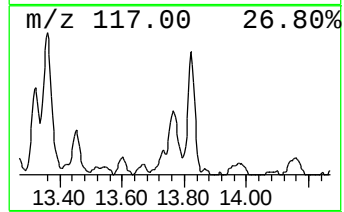
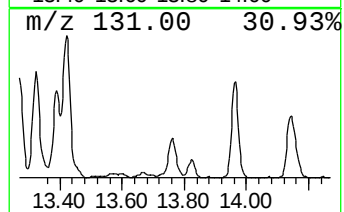
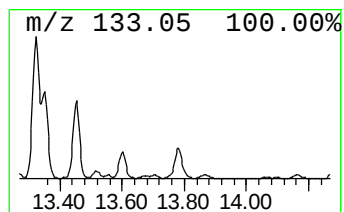
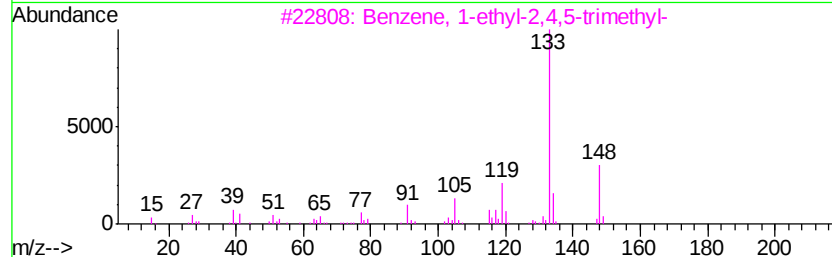
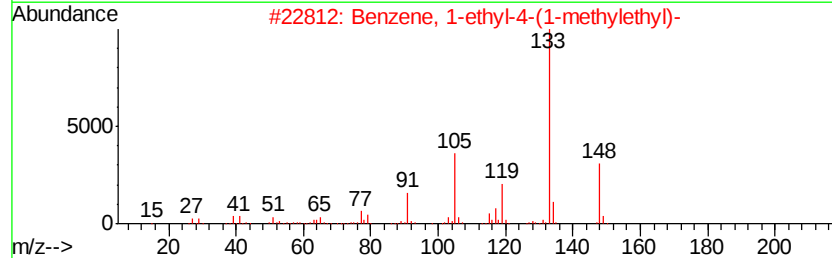
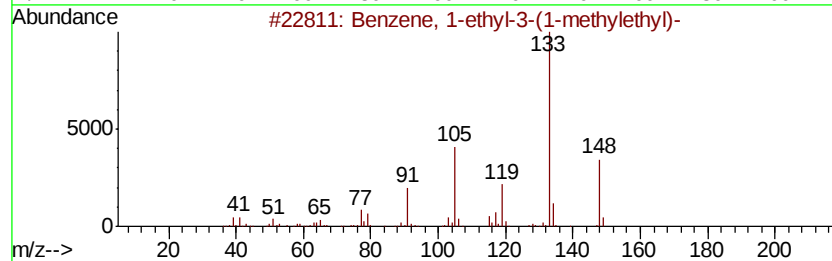
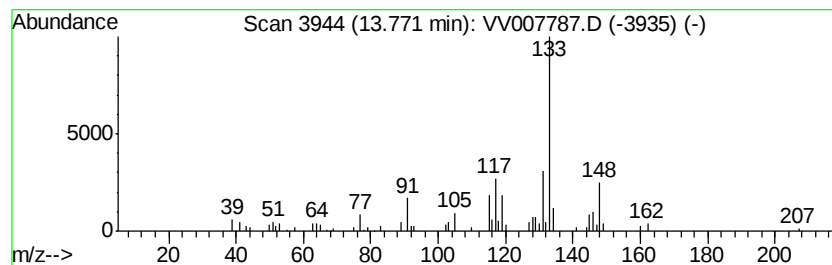
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 38 Benzene, 1-ethyl-3-(1-methy... Concentration Rank 36

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 13.77 | 4.12 ug/L | 50414 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Benzene, 1-ethyl-3-(1-methylethyl)- | 148 | C11H16   | 004920-99-4  | 58   |
| 2    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 | C11H16   | 004218-48-8  | 58   |
| 3    |    |   | Benzene, 1-ethyl-2,4,5-trimethyl-   | 148 | C11H16   | 017851-27-3  | 58   |
| 4    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 | C11H16   | 004218-48-8  | 58   |
| 5    |    |   | Benzene, 1,3-bis(1-formylethyl)-    | 190 | C12H14O2 | 1000160-34-1 | 53   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

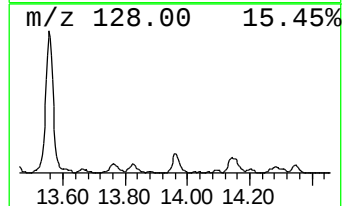
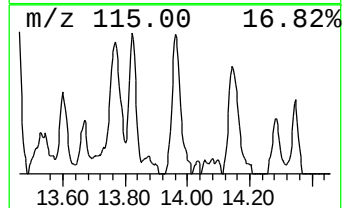
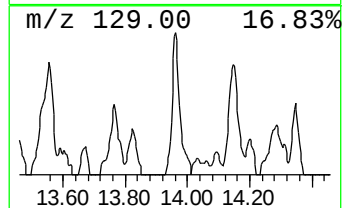
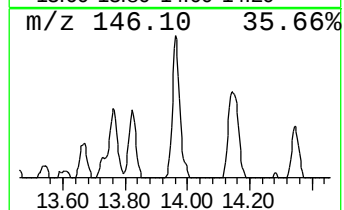
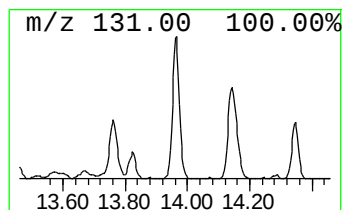
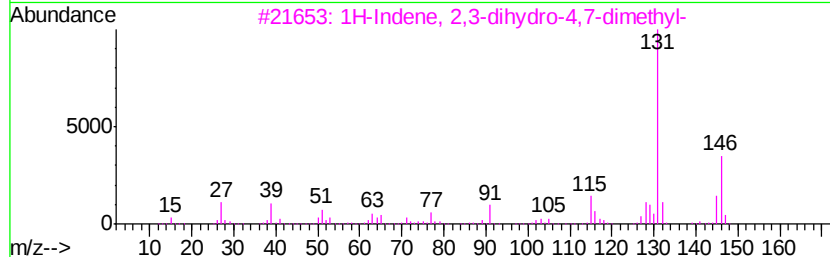
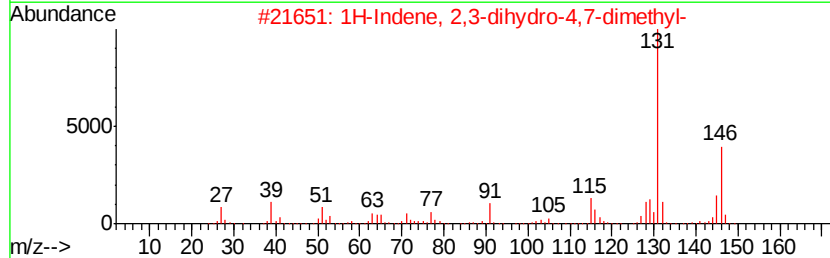
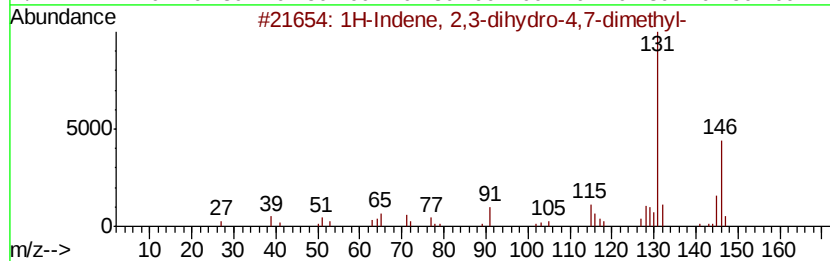
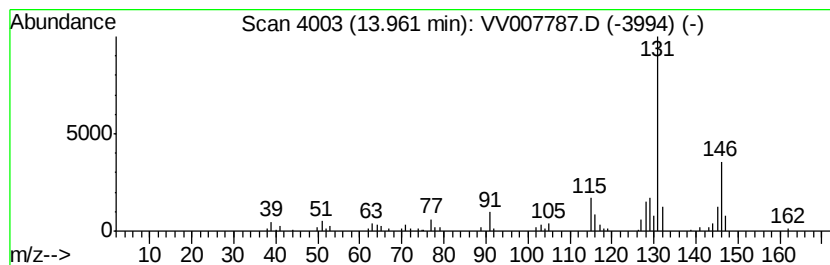
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 32

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 13.96 | 4.78 ug/L | 58488 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 96   |
| 2    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 94   |
| 3    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 94   |
| 4    |    | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 94   |
| 5    |    | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146 | C11H14  | 004175-53-5 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

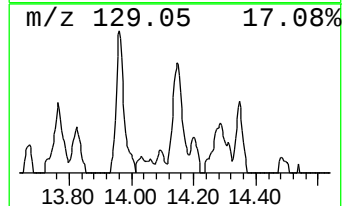
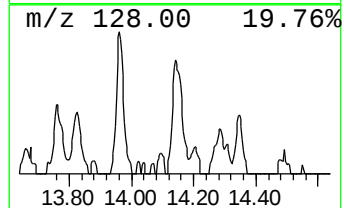
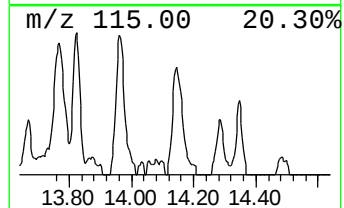
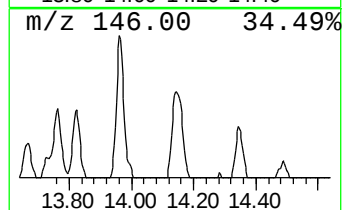
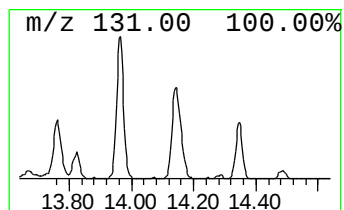
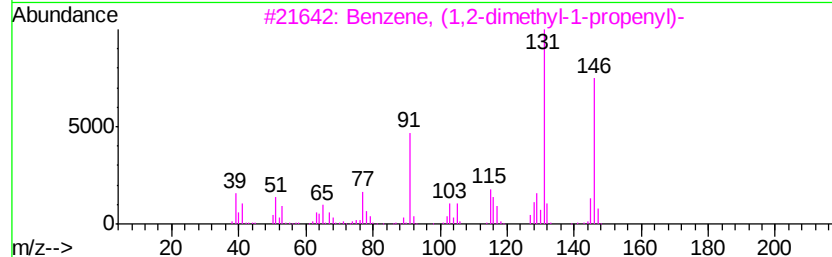
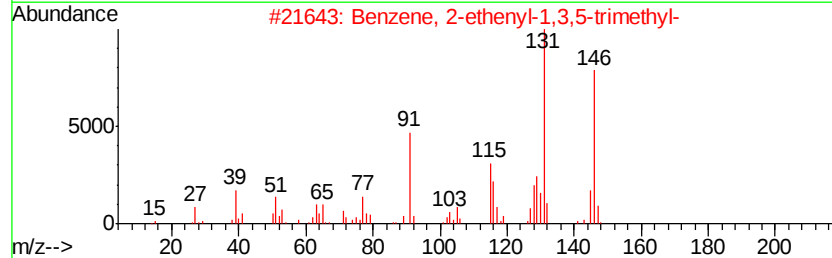
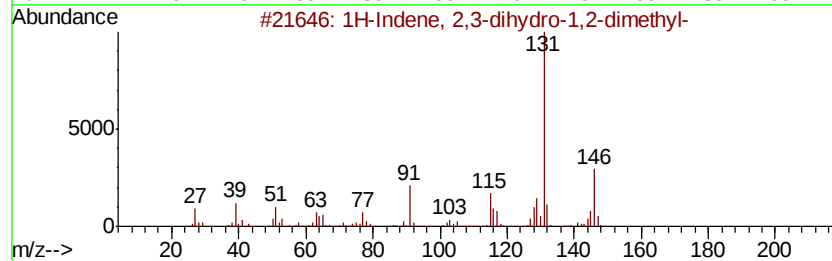
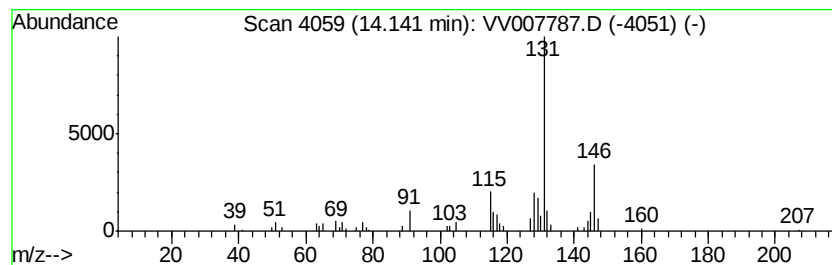
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 40 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 35

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 14.14 | 4.25 ug/L | 51991 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 95   |
| 2    |    | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 | C11H14  | 000769-25-5 | 90   |
| 3    |    | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 87   |
| 4    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 87   |
| 5    |    | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14  | 017059-48-2 | 87   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

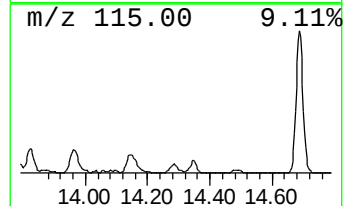
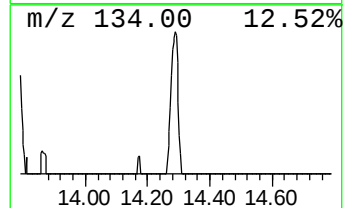
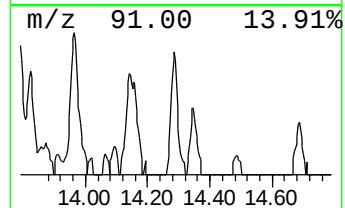
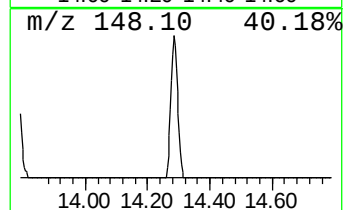
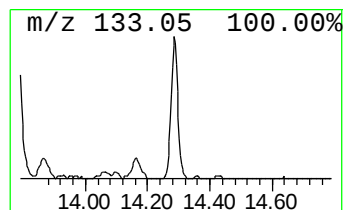
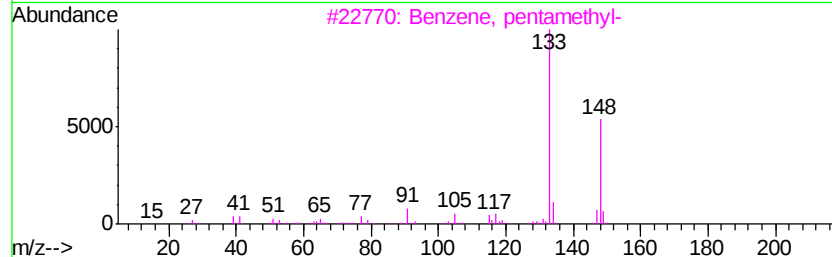
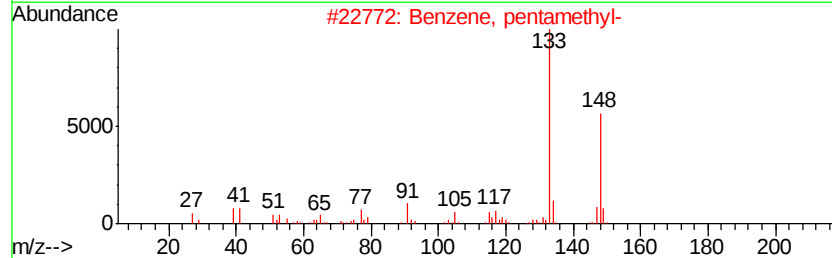
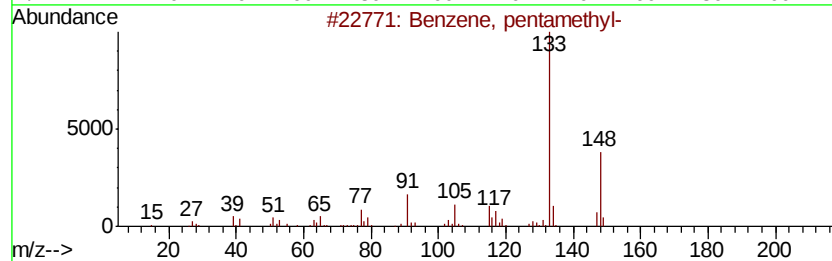
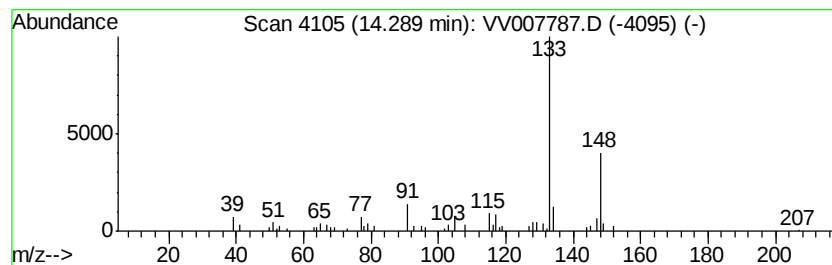
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 41 Benzene, pentamethyl- Concentration Rank 40

| R.T.  | EstConc   | Area  | Relative to ISTD       | R.T.  |
|-------|-----------|-------|------------------------|-------|
| 14.29 | 2.94 ug/L | 35969 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9 | 94   |
| 2    |    | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9 | 91   |
| 3    |    | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9 | 91   |
| 4    |    | Benzene, 1,3-dimethyl-5-(1-methy... | 148 | C11H16  | 004706-90-5 | 90   |
| 5    |    | Benzene, 2,4-dimethyl-1-(1-methy... | 148 | C11H16  | 004706-89-2 | 87   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV092318\  
Data File : VV007787.D  
Acq On : 23 Sep 2018 14:43  
Operator : SY/MD  
Sample : J5014-02  
Misc : 5.0 mL/MSVOA V/WATER  
ALS Vial : 65 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
E8JB2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.M  
Quant Title : VOC Analysis

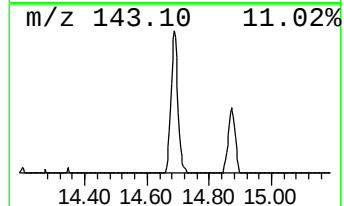
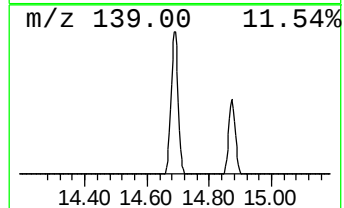
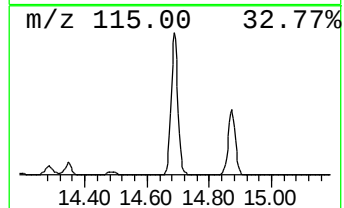
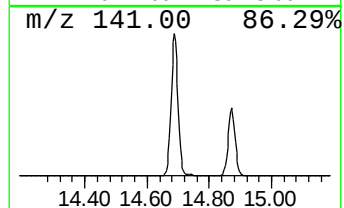
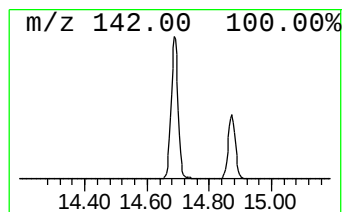
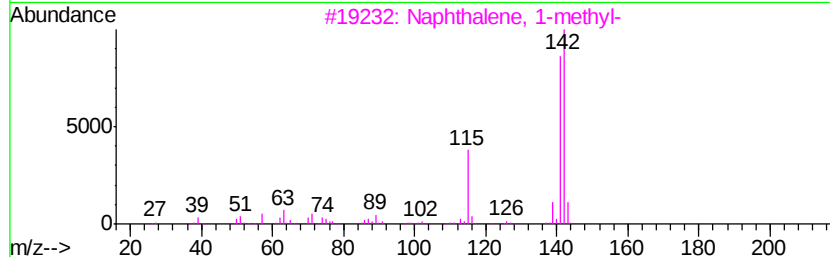
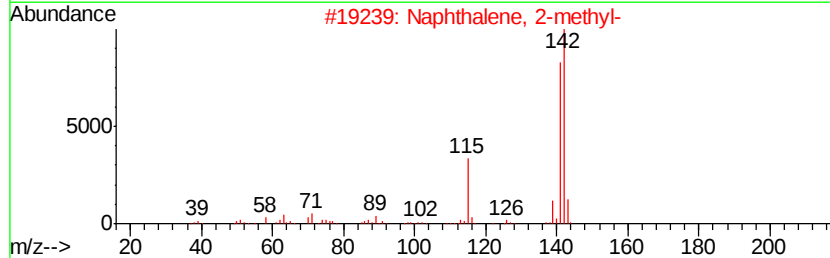
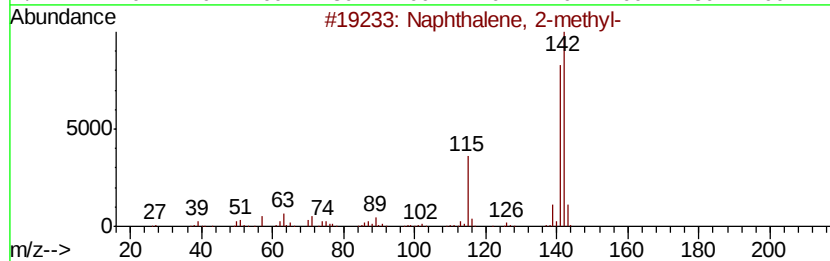
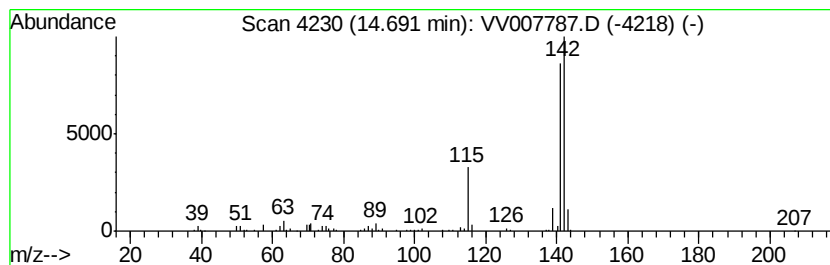
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 42 Naphthalene, 1-methyl- Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 14.69 | 12.77 ug/L | 156353 | 1,4-Dichlorobenzene-d4 | 11.30 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 96   |
| 2    |    |   | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 96   |
| 3    |    |   | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 96   |
| 4    |    |   | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 94   |
| 5    |    |   | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 94   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_V\DATA\VV092318\  
 Data File : VV007787.D  
 Acq On : 23 Sep 2018 14:43  
 Operator : SY/MD  
 Sample : J5014-02  
 Misc : 5.0 mL/MSVOA\_V/WATER  
 ALS Vial : 65 Sample Multiplier: 1

Instrument :  
 MSVOA\_V  
 ClientSampleId :  
 E8JB2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVLM092318WMA.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: Cyc... | 8.40  | 2.9     | ug/L  | 29062    | 2                     | 8.90  | 496230 | 50.0 |
| (DEL) Alkane: Cyc... | 8.64  | 8.7     | ug/L  | 85959    | 2                     | 8.90  | 496230 | 50.0 |
| (DEL) Alkane: Str... | 8.70  | 3.2     | ug/L  | 31314    | 2                     | 8.90  | 496230 | 50.0 |
| (DEL) Alkane: Str... | 8.83  | 3.9     | ug/L  | 38309    | 2                     | 8.90  | 496230 | 50.0 |
| Cyclohexene, 4-pr... | 9.37  | 2.5     | ug/L  | 25176    | 2                     | 8.90  | 496230 | 50.0 |
| (DEL) Alkane: Cyc... | 9.52  | 28.5    | ug/L  | 283289   | 2                     | 8.90  | 496230 | 50.0 |
| (DEL) Alkane: Cyc... | 9.72  | 4.8     | ug/L  | 47383    | 2                     | 8.90  | 496230 | 50.0 |
| Pentalene, octahy... | 9.76  | 10.9    | ug/L  | 107960   | 2                     | 8.90  | 496230 | 50.0 |
| (DEL) Alkane: Cyc... | 9.85  | 18.2    | ug/L  | 180808   | 2                     | 8.90  | 496230 | 50.0 |
| unknown-01           | 10.18 | 4.8     | ug/L  | 58588    | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1-ethyl-... | 10.50 | 16.5    | ug/L  | 202288   | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1-ethyl-... | 10.78 | 19.2    | ug/L  | 235270   | 3                     | 11.30 | 612028 | 50.0 |
| (DEL) Alkane: Str... | 10.93 | 4.9     | ug/L  | 60414    | 3                     | 11.30 | 612028 | 50.0 |
| (DEL) Alkane: Cyc... | 11.20 | 6.6     | ug/L  | 81156    | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1,4-diet... | 11.80 | 4.5     | ug/L  | 55705    | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1-methyl... | 11.87 | 15.5    | ug/L  | 189682   | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1-ethyl-... | 12.07 | 59.4    | ug/L  | 726853   | 3                     | 11.30 | 612028 | 50.0 |
| Indan, 1-methyl-     | 12.17 | 11.8    | ug/L  | 144755   | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 2,4-diet... | 12.31 | 7.6     | ug/L  | 93085    | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1,2,3,5-... | 12.46 | 50.7    | ug/L  | 620898   | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1,2,4,5-... | 12.52 | 61.7    | ug/L  | 754985   | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1,3-diet... | 12.64 | 2.5     | ug/L  | 31080    | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1-etheny... | 12.78 | 21.6    | ug/L  | 264455   | 3                     | 11.30 | 612028 | 50.0 |
| 1-Phenyl-1-butene    | 12.94 | 122.3   | ug/L  | 1497620  | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1-methyl... | 13.05 | 8.2     | ug/L  | 100448   | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1-butyryl-  | 13.11 | 10.0    | ug/L  | 121909   | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1,3-dime... | 13.22 | 5.7     | ug/L  | 69661    | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1-ethyl-... | 13.32 | 17.5    | ug/L  | 214386   | 3                     | 11.30 | 612028 | 50.0 |
| 3,4-Dimethylcumene   | 13.60 | 3.1     | ug/L  | 37938    | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, 1-ethyl-... | 13.77 | 4.1     | ug/L  | 50414    | 3                     | 11.30 | 612028 | 50.0 |
| 1H-Indene, 2,3-di... | 13.96 | 4.8     | ug/L  | 58488    | 3                     | 11.30 | 612028 | 50.0 |
| 1H-Indene, 2,3-di... | 14.14 | 4.3     | ug/L  | 51991    | 3                     | 11.30 | 612028 | 50.0 |
| Benzene, pentamet... | 14.29 | 2.9     | ug/L  | 35969    | 3                     | 11.30 | 612028 | 50.0 |
| Naphthalene, 1-me... | 14.69 | 12.8    | ug/L  | 156353   | 3                     | 11.30 | 612028 | 50.0 |