

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
 Data File : VU034840.D  
 Acq On : 26 Sep 2019 12:48  
 Operator : JC/SP  
 Sample : K4984-08  
 Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BEDL2

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 0 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
 Title : VOC Analysis

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 1.283       | 56            | 60          | 67           | rBV      | 13688          | 14642         | 0.19%           | 0.011%        |
| 2         | 1.386       | 85            | 92          | 107          | rVB      | 251808         | 305448        | 4.01%           | 0.230%        |
| 3         | 1.630       | 128           | 168         | 179          | rBV      | 441966         | 1418173       | 18.62%          | 1.068%        |
| 4         | 2.222       | 342           | 352         | 365          | rVB      | 542797         | 817877        | 10.74%          | 0.616%        |
| 5         | 2.431       | 413           | 417         | 422          | rBV4     | 1421           | 1419          | 0.02%           | 0.001%        |
| 6         | 2.611       | 470           | 473         | 479          | rVB5     | 2115           | 2080          | 0.03%           | 0.002%        |
| 7         | 2.669       | 486           | 491         | 499          | rVB5     | 5267           | 7477          | 0.10%           | 0.006%        |
| 8         | 2.810       | 523           | 535         | 547          | rBV      | 39322          | 91355         | 1.20%           | 0.069%        |
| 9         | 3.035       | 603           | 605         | 613          | rVB7     | 1329           | 1464          | 0.02%           | 0.001%        |
| 10        | 3.254       | 670           | 673         | 682          | rVB8     | 2040           | 2623          | 0.03%           | 0.002%        |
| 11        | 3.447       | 726           | 733         | 737          | rVV8     | 1303           | 1511          | 0.02%           | 0.001%        |
| 12        | 3.521       | 752           | 756         | 762          | rBV5     | 1341           | 1515          | 0.02%           | 0.001%        |
| 13        | 3.952       | 888           | 890         | 895          | rVB5     | 1739           | 1400          | 0.02%           | 0.001%        |
| 14        | 4.042       | 906           | 918         | 919          | rBV6     | 3494           | 5099          | 0.07%           | 0.004%        |
| 15        | 4.167       | 941           | 957         | 997          | rBV      | 201214         | 695443        | 9.13%           | 0.524%        |
| 16        | 4.617       | 1078          | 1097        | 1124         | rBV      | 373698         | 918479        | 12.06%          | 0.692%        |
| 17        | 4.743       | 1134          | 1136        | 1142         | rBV6     | 1500           | 1696          | 0.02%           | 0.001%        |
| 18        | 4.791       | 1149          | 1151        | 1156         | rBV4     | 2052           | 1979          | 0.03%           | 0.001%        |
| 19        | 4.842       | 1161          | 1167        | 1170         | rBV6     | 2206           | 2487          | 0.03%           | 0.002%        |
| 20        | 4.952       | 1189          | 1201        | 1216         | rVB6     | 22910          | 55879         | 0.73%           | 0.042%        |
| 21        | 5.035       | 1216          | 1227        | 1229         | rBV8     | 7524           | 9591          | 0.13%           | 0.007%        |
| 22        | 5.103       | 1242          | 1248        | 1253         | rVB9     | 2184           | 2163          | 0.03%           | 0.002%        |
| 23        | 5.186       | 1261          | 1274        | 1289         | rVV3     | 22483          | 56722         | 0.74%           | 0.043%        |
| 24        | 5.305       | 1289          | 1311        | 1342         | rVV2     | 844699         | 2419082       | 31.77%          | 1.822%        |
| 25        | 5.424       | 1344          | 1348        | 1349         | rBV5     | 3503           | 2539          | 0.03%           | 0.002%        |
| 26        | 5.444       | 1349          | 1354        | 1362         | rVB6     | 8773           | 13710         | 0.18%           | 0.010%        |
| 27        | 5.588       | 1389          | 1399        | 1414         | rVB7     | 20476          | 48470         | 0.64%           | 0.037%        |
| 28        | 5.752       | 1431          | 1450        | 1468         | rBV3     | 50997          | 111487        | 1.46%           | 0.084%        |
| 29        | 5.858       | 1469          | 1483        | 1505         | rBV2     | 591918         | 1222967       | 16.06%          | 0.921%        |
| 30        | 6.305       | 1609          | 1622        | 1635         | rBV      | 543295         | 1139030       | 14.96%          | 0.858%        |
| 31        | 6.450       | 1661          | 1667        | 1674         | rVB5     | 13895          | 20559         | 0.27%           | 0.015%        |
| 32        | 6.505       | 1677          | 1684        | 1695         | rVB4     | 28457          | 51421         | 0.68%           | 0.039%        |
| 33        | 6.611       | 1707          | 1717        | 1729         | rVB3     | 33283          | 62475         | 0.82%           | 0.047%        |
| 34        | 6.707       | 1734          | 1747        | 1761         | rVB4     | 53150          | 116001        | 1.52%           | 0.087%        |

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Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BEDL2

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 0 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
 Title : VOC Analysis

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 6.878  | 1785 | 1800 | 1820 | rVB5 | 54844   | 147888  | 1.94%  | 0.111% |
| 36 | 7.019  | 1834 | 1844 | 1853 | rBV7 | 13500   | 29811   | 0.39%  | 0.022% |
| 37 | 7.080  | 1854 | 1863 | 1870 | rBV3 | 41841   | 71852   | 0.94%  | 0.054% |
| 38 | 7.154  | 1878 | 1886 | 1893 | rBV  | 163218  | 259267  | 3.40%  | 0.195% |
| 39 | 7.202  | 1893 | 1901 | 1917 | rVB3 | 344682  | 678733  | 8.91%  | 0.511% |
| 40 | 7.315  | 1928 | 1936 | 1957 | rVB2 | 167798  | 360653  | 4.74%  | 0.272% |
| 41 | 7.418  | 1958 | 1968 | 1977 | rVB4 | 35756   | 70697   | 0.93%  | 0.053% |
| 42 | 7.501  | 1981 | 1994 | 1998 | rBV4 | 344874  | 573410  | 7.53%  | 0.432% |
| 43 | 7.543  | 1999 | 2007 | 2025 | rVB  | 1186480 | 2180033 | 28.63% | 1.642% |
| 44 | 7.678  | 2037 | 2049 | 2057 | rBV4 | 111424  | 230694  | 3.03%  | 0.174% |
| 45 | 7.794  | 2077 | 2085 | 2092 | rBV  | 536266  | 831998  | 10.93% | 0.627% |
| 46 | 7.890  | 2107 | 2115 | 2134 | rVB2 | 276547  | 510322  | 6.70%  | 0.384% |
| 47 | 8.022  | 2136 | 2156 | 2167 | rBV  | 209393  | 400083  | 5.25%  | 0.301% |
| 48 | 8.083  | 2168 | 2175 | 2178 | rBV4 | 15968   | 22428   | 0.29%  | 0.017% |
| 49 | 8.202  | 2203 | 2212 | 2223 | rVB3 | 107987  | 185790  | 2.44%  | 0.140% |
| 50 | 8.305  | 2226 | 2244 | 2268 | rBV  | 1139413 | 2225682 | 29.23% | 1.677% |
| 51 | 8.431  | 2275 | 2283 | 2286 | rBV2 | 136234  | 190015  | 2.50%  | 0.143% |
| 52 | 8.521  | 2303 | 2311 | 2320 | rVB  | 434128  | 718363  | 9.43%  | 0.541% |
| 53 | 8.582  | 2320 | 2330 | 2339 | rBV  | 499472  | 886971  | 11.65% | 0.668% |
| 54 | 8.733  | 2368 | 2377 | 2386 | rBV3 | 118872  | 198168  | 2.60%  | 0.149% |
| 55 | 8.791  | 2386 | 2395 | 2400 | rBV2 | 233207  | 376618  | 4.95%  | 0.284% |
| 56 | 8.887  | 2418 | 2425 | 2437 | rVB2 | 654112  | 1092014 | 14.34% | 0.823% |
| 57 | 8.955  | 2438 | 2446 | 2452 | rBV3 | 37616   | 57817   | 0.76%  | 0.044% |
| 58 | 9.009  | 2452 | 2463 | 2472 | rBV  | 574100  | 971426  | 12.76% | 0.732% |
| 59 | 9.070  | 2473 | 2482 | 2496 | rVV  | 859376  | 1483919 | 19.49% | 1.118% |
| 60 | 9.228  | 2520 | 2531 | 2544 | rBV3 | 444791  | 867338  | 11.39% | 0.653% |
| 61 | 9.421  | 2584 | 2591 | 2618 | rVB  | 2007516 | 3274566 | 43.00% | 2.467% |
| 62 | 9.543  | 2619 | 2629 | 2640 | rBV3 | 141952  | 254207  | 3.34%  | 0.191% |
| 63 | 9.694  | 2663 | 2676 | 2686 | rBV3 | 680665  | 1418419 | 18.63% | 1.068% |
| 64 | 9.929  | 2732 | 2749 | 2759 | rBV3 | 962815  | 1844893 | 24.23% | 1.390% |
| 65 | 10.022 | 2770 | 2778 | 2786 | rBV5 | 681494  | 1184771 | 15.56% | 0.892% |
| 66 | 10.144 | 2807 | 2816 | 2825 | rVB  | 460604  | 796636  | 10.46% | 0.600% |
| 67 | 10.418 | 2889 | 2901 | 2911 | rBV2 | 1049539 | 2609646 | 34.27% | 1.966% |
| 68 | 10.566 | 2940 | 2947 | 2955 | rBV  | 544411  | 805039  | 10.57% | 0.606% |
| 69 | 10.662 | 2969 | 2977 | 2991 | rBV  | 1323827 | 3067176 | 40.28% | 2.310% |
| 70 | 10.733 | 2992 | 2999 | 3009 | rVV3 | 714526  | 1788307 | 23.48% | 1.347% |
| 71 | 10.791 | 3010 | 3017 | 3028 | rVB  | 3617718 | 5223541 | 68.60% | 3.935% |

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## Integration Parameters: LSCINT.P

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 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
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 Min Area: 0 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
 Title : VOC Analysis

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 72  | 10.951 | 3058 | 3067 | 3075 | rBV  | 1250028 | 2130998 | 27.98%  | 1.605% |
| 73  | 11.093 | 3104 | 3111 | 3116 | rBV  | 999333  | 1466829 | 19.26%  | 1.105% |
| 74  | 11.125 | 3117 | 3121 | 3133 | rVB  | 1213433 | 1756546 | 23.07%  | 1.323% |
| 75  | 11.376 | 3192 | 3199 | 3204 | rBV3 | 768015  | 1187976 | 15.60%  | 0.895% |
| 76  | 11.459 | 3217 | 3225 | 3233 | rBV3 | 1380540 | 2046161 | 26.87%  | 1.541% |
| 77  | 11.549 | 3247 | 3253 | 3259 | rBV  | 786511  | 1004534 | 13.19%  | 0.757% |
| 78  | 11.588 | 3260 | 3265 | 3275 | rVB  | 786678  | 1058965 | 13.91%  | 0.798% |
| 79  | 11.678 | 3288 | 3293 | 3305 | rVB3 | 865419  | 1238256 | 16.26%  | 0.933% |
| 80  | 11.755 | 3308 | 3317 | 3321 | rBV3 | 723705  | 1231593 | 16.17%  | 0.928% |
| 81  | 11.839 | 3334 | 3343 | 3361 | rVB7 | 2879380 | 6515891 | 85.57%  | 4.908% |
| 82  | 12.003 | 3386 | 3394 | 3401 | rBV  | 4283750 | 6138194 | 80.61%  | 4.624% |
| 83  | 12.125 | 3425 | 3432 | 3436 | rBV  | 782166  | 1139222 | 14.96%  | 0.858% |
| 84  | 12.337 | 3490 | 3498 | 3503 | rBV2 | 955865  | 1531414 | 20.11%  | 1.154% |
| 85  | 12.472 | 3532 | 3540 | 3551 | rBV3 | 1283114 | 2778493 | 36.49%  | 2.093% |
| 86  | 12.591 | 3568 | 3577 | 3582 | rBV4 | 1217209 | 1994062 | 26.19%  | 1.502% |
| 87  | 12.681 | 3598 | 3605 | 3613 | rBV2 | 1315184 | 2413693 | 31.70%  | 1.818% |
| 88  | 12.768 | 3625 | 3632 | 3637 | rBV2 | 918017  | 1401298 | 18.40%  | 1.056% |
| 89  | 13.012 | 3701 | 3708 | 3716 | rBV3 | 933934  | 1451015 | 19.06%  | 1.093% |
| 90  | 13.102 | 3729 | 3736 | 3751 | rBV3 | 4926322 | 7614824 | 100.00% | 5.736% |
| 91  | 13.263 | 3778 | 3786 | 3794 | rBV5 | 2380727 | 3936327 | 51.69%  | 2.965% |
| 92  | 13.434 | 3833 | 3839 | 3847 | rVB3 | 1161696 | 1667144 | 21.89%  | 1.256% |
| 93  | 13.488 | 3848 | 3856 | 3862 | rBV3 | 1505528 | 2388690 | 31.37%  | 1.799% |
| 94  | 13.765 | 3936 | 3942 | 3953 | rBV3 | 1099565 | 1726925 | 22.68%  | 1.301% |
| 95  | 13.835 | 3956 | 3964 | 3971 | rBV2 | 2339162 | 4039382 | 53.05%  | 3.043% |
| 96  | 13.929 | 3986 | 3993 | 4003 | rBV3 | 2351672 | 3998859 | 52.51%  | 3.012% |
| 97  | 14.122 | 4046 | 4053 | 4069 | rBV3 | 3113384 | 5145166 | 67.57%  | 3.876% |
| 98  | 14.327 | 4106 | 4117 | 4128 | rVB4 | 3829266 | 7413717 | 97.36%  | 5.585% |
| 99  | 14.655 | 4213 | 4219 | 4229 | rBV2 | 1915161 | 2901584 | 38.10%  | 2.186% |
| 100 | 14.842 | 4271 | 4277 | 4292 | rVB5 | 3306526 | 5920993 | 77.76%  | 4.460% |

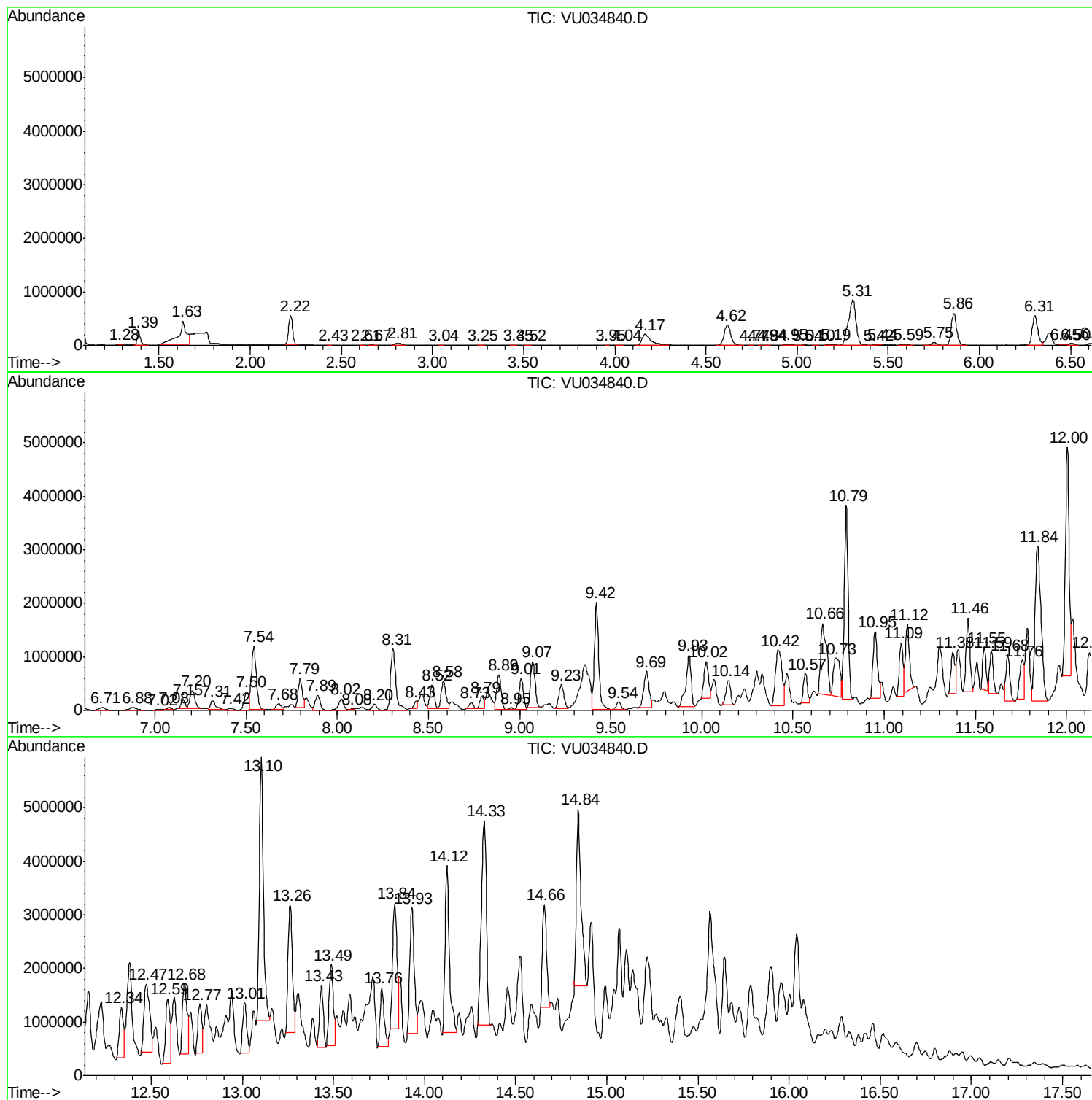
Sum of corrected areas: 132752205

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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



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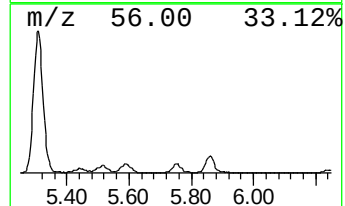
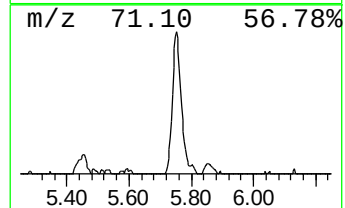
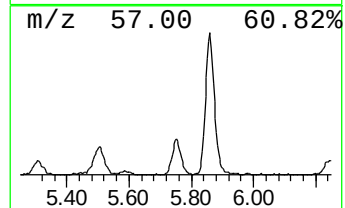
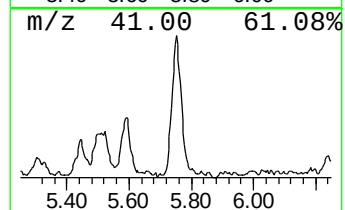
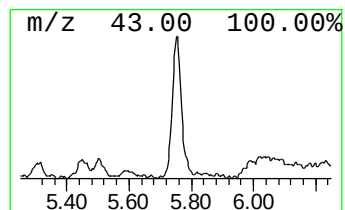
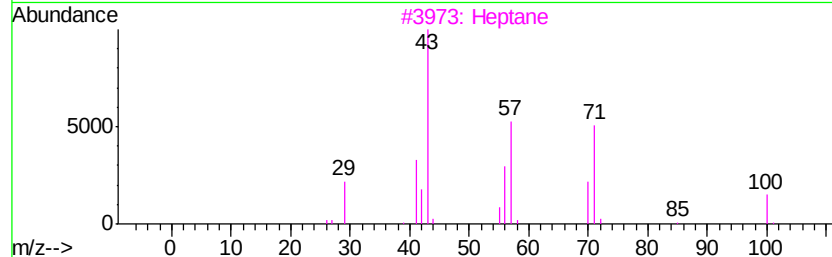
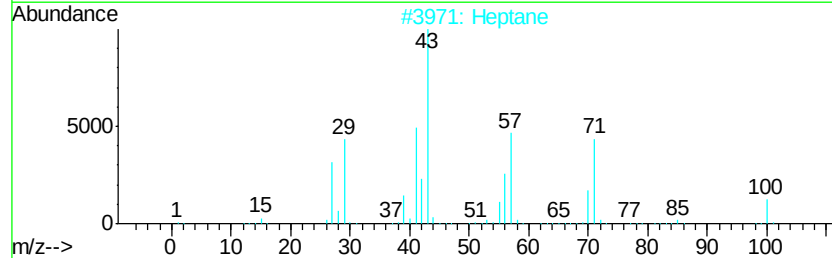
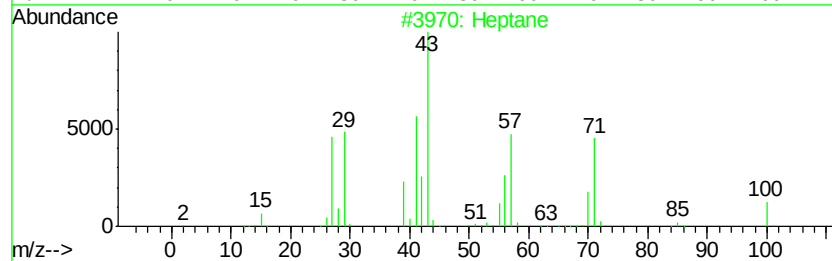
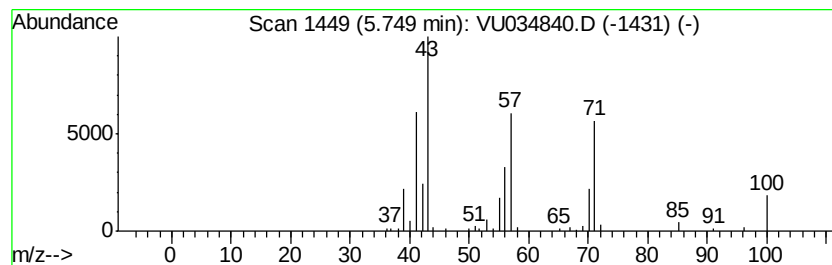
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 53

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 5.75 | 4.56 ug/L | 111487 | 1,4-Difluorobenzene | 5.86 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Heptane               | 100 | C7H16   | 000142-82-5 | 91   |
| 2    |    | Heptane               | 100 | C7H16   | 000142-82-5 | 91   |
| 3    |    | Heptane               | 100 | C7H16   | 000142-82-5 | 90   |
| 4    |    | Heptane               | 100 | C7H16   | 000142-82-5 | 81   |
| 5    |    | Hexane, 3,3-dimethyl- | 114 | C8H18   | 000563-16-6 | 50   |



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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

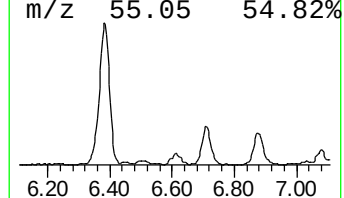
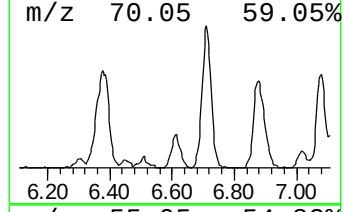
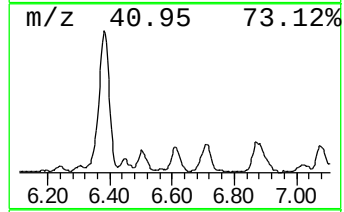
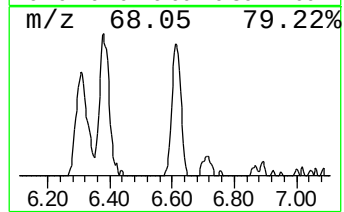
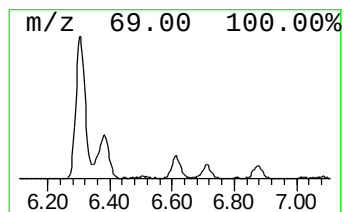
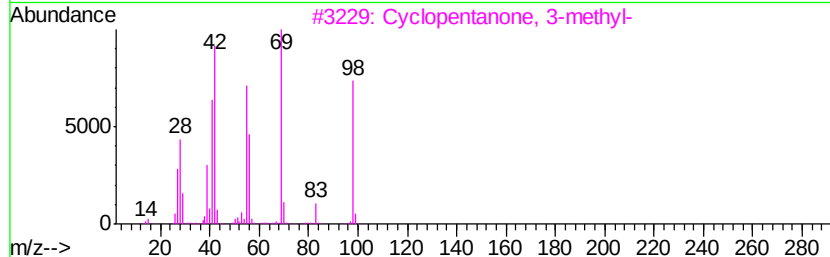
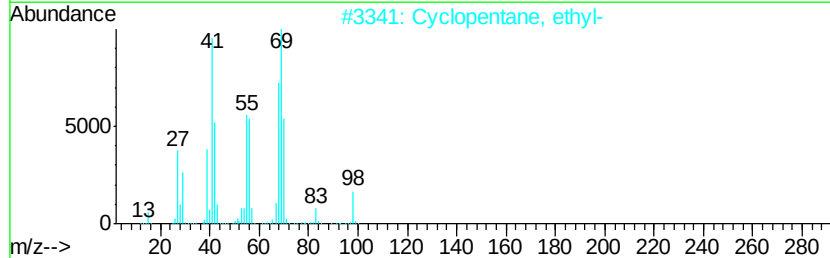
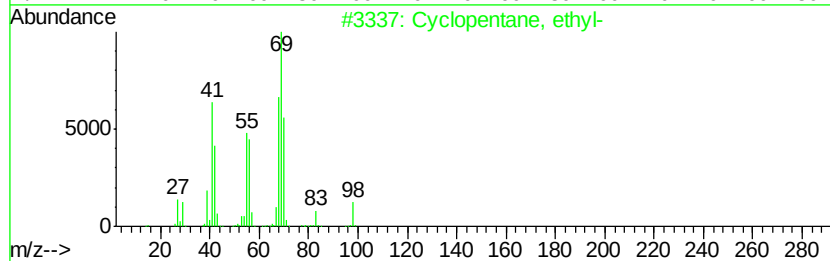
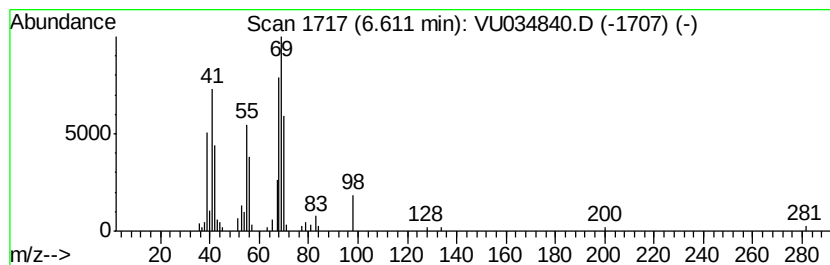
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Cyclic6.61 Concentration Rank 57

| R.T. | EstConc   | Area  | Relative to ISTD    | R.T. |
|------|-----------|-------|---------------------|------|
| 6.61 | 2.55 ug/L | 62475 | 1,4-Difluorobenzene | 5.86 |

| Hit# | of | Tentative ID              | MW | MolForm | CAS#        | Qual |
|------|----|---------------------------|----|---------|-------------|------|
| 1    | 5  | Cyclopentane, ethyl-      | 98 | C7H14   | 001640-89-7 | 80   |
| 2    |    | Cyclopentane, ethyl-      | 98 | C7H14   | 001640-89-7 | 68   |
| 3    |    | Cyclopentanone, 3-methyl- | 98 | C6H10O  | 001757-42-2 | 38   |
| 4    |    | 1-Pentene, 2,3-dimethyl-  | 98 | C7H14   | 003404-72-6 | 38   |
| 5    |    | 1-Pentene, 2,3-dimethyl-  | 98 | C7H14   | 003404-72-6 | 38   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleID :  
BEDL2

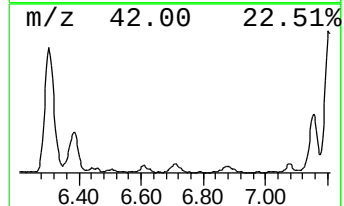
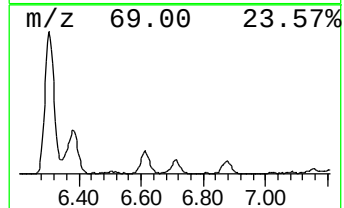
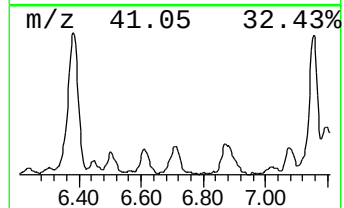
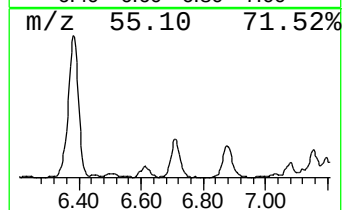
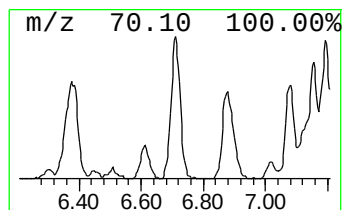
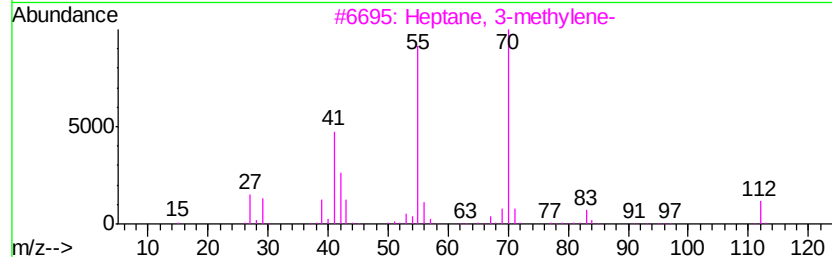
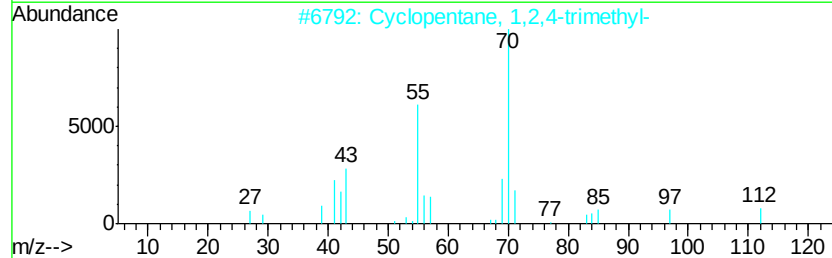
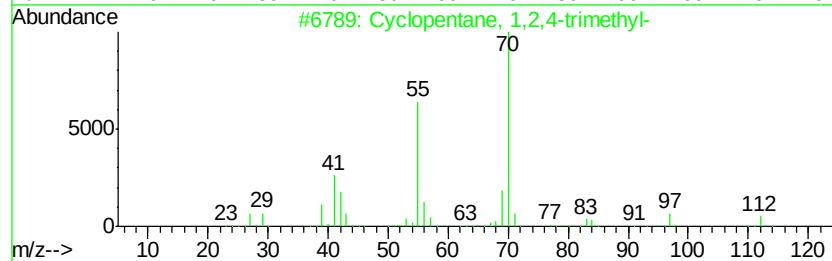
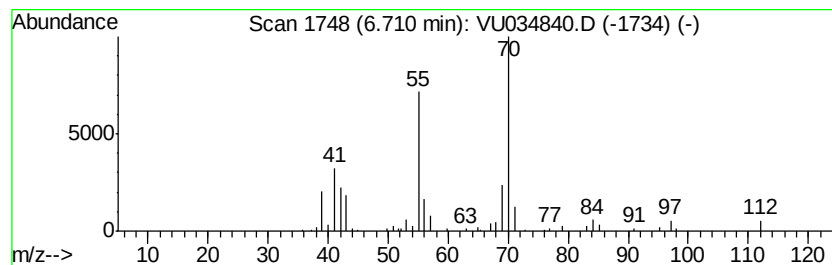
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Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Cyclic6.71 Concentration Rank 52

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 6.71 | 4.74 ug/L | 116001 | 1,4-Difluorobenzene | 5.86 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopentane, 1,2,4-trimethyl-      | 112 | C8H16   | 002815-58-9 | 90   |
| 2    |    | Cyclopentane, 1,2,4-trimethyl-      | 112 | C8H16   | 002815-58-9 | 87   |
| 3    |    | Heptane, 3-methylene-               | 112 | C8H16   | 001632-16-2 | 86   |
| 4    |    | Cyclopentane, 1,2,3-trimethyl-, ... | 112 | C8H16   | 002613-69-6 | 80   |
| 5    |    | Cyclopentane, 1,2,3-trimethyl-, ... | 112 | C8H16   | 002613-69-6 | 80   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

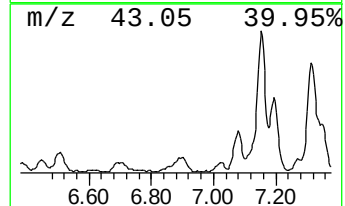
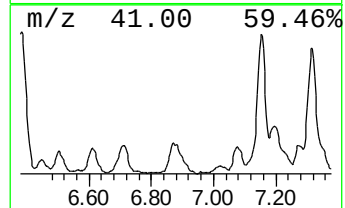
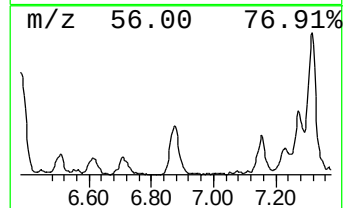
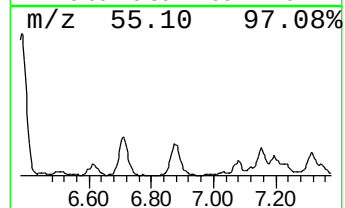
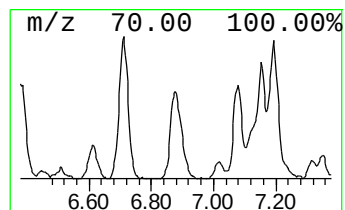
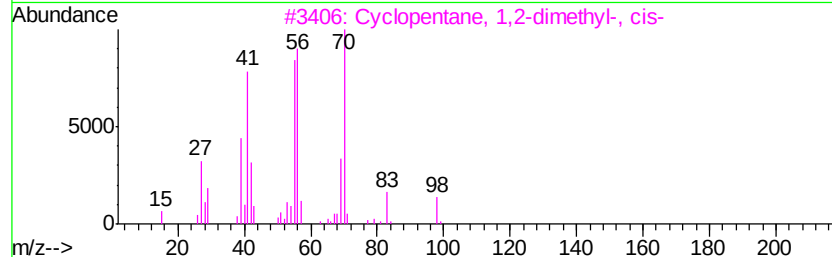
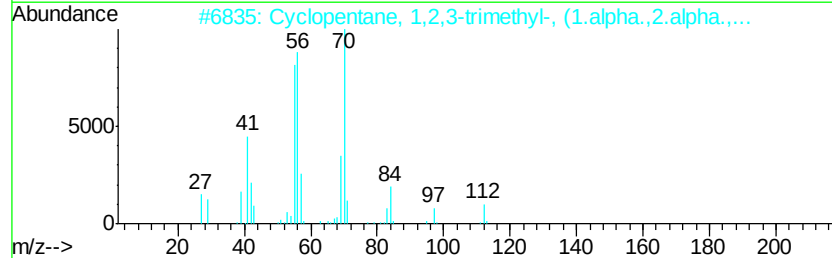
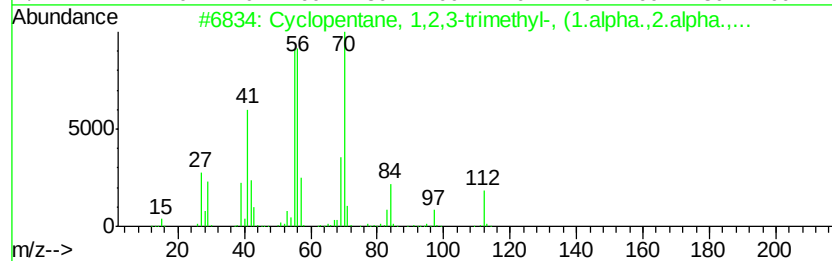
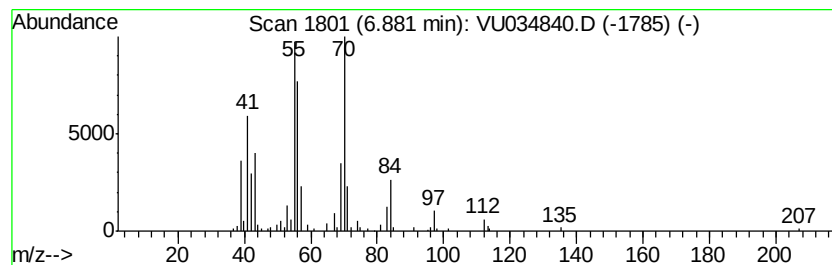
Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Cyclic6.88 Concentration Rank 51

| R.T.      | EstConc                             | Area   | Relative to ISTD    | R.T.        |      |
|-----------|-------------------------------------|--------|---------------------|-------------|------|
| 6.88      | 6.05 ug/L                           | 147888 | 1,4-Difluorobenzene | 5.86        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm             | CAS#        | Qual |
| 1         | Cyclopentane, 1,2,3-trimethyl-, ... | 112    | C8H16               | 015890-40-1 | 90   |
| 2         | Cyclopentane, 1,2,3-trimethyl-, ... | 112    | C8H16               | 015890-40-1 | 78   |
| 3         | Cyclopentane, 1,2-dimethyl-, cis-   | 98     | C7H14               | 001192-18-3 | 72   |
| 4         | 1-Octene, 3-methyl-                 | 126    | C9H18               | 013151-08-1 | 72   |
| 5         | Cyclopentane, 1,2,3-trimethyl-, ... | 112    | C8H16               | 002613-69-6 | 58   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

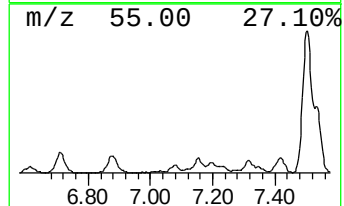
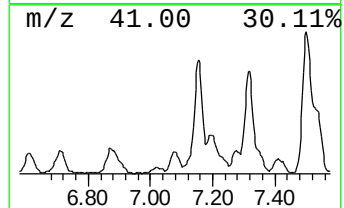
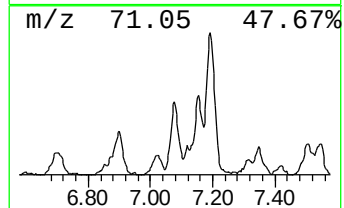
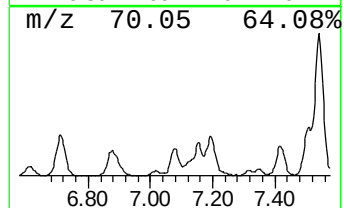
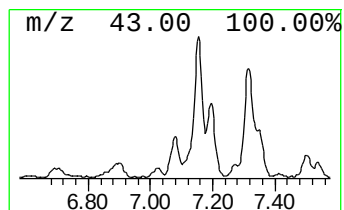
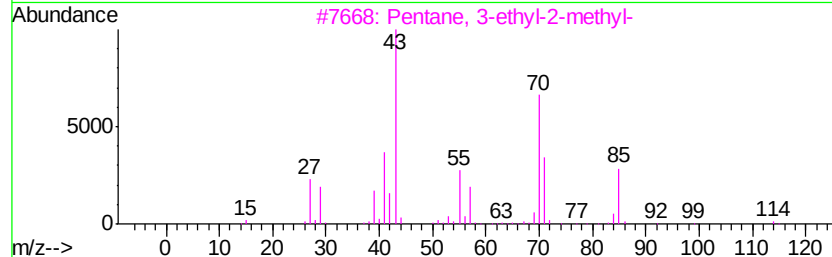
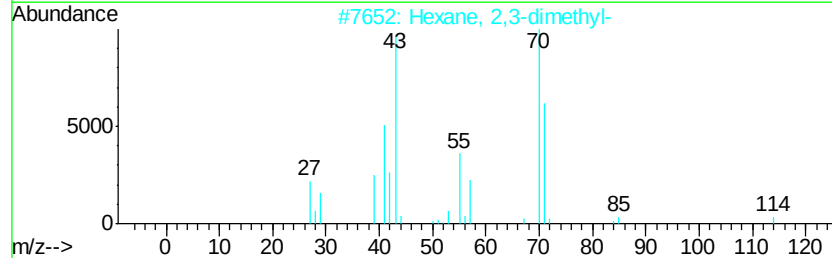
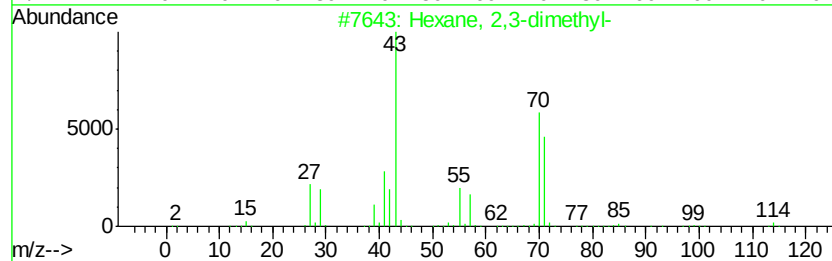
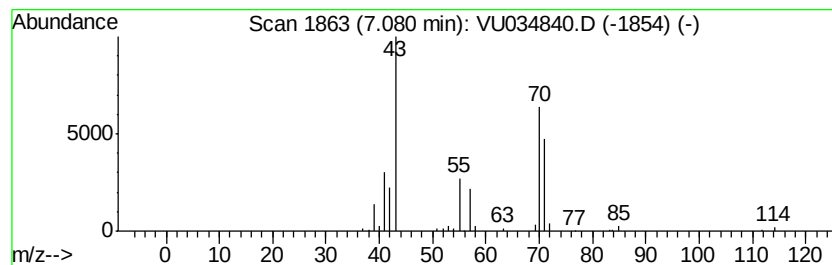
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 55

| R.T. | EstConc   | Area  | Relative to ISTD    | R.T. |
|------|-----------|-------|---------------------|------|
| 7.08 | 2.94 ug/L | 71852 | 1,4-Difluorobenzene | 5.86 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexane, 2,3-dimethyl-      | 114 | C8H18   | 000584-94-1 | 91   |
| 2    |    | Hexane, 2,3-dimethyl-      | 114 | C8H18   | 000584-94-1 | 90   |
| 3    |    | Pentane, 3-ethyl-2-methyl- | 114 | C8H18   | 000609-26-7 | 83   |
| 4    |    | Hexane, 2,3-dimethyl-      | 114 | C8H18   | 000584-94-1 | 78   |
| 5    |    | Pyrrolidine                | 71  | C4H9N   | 000123-75-1 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

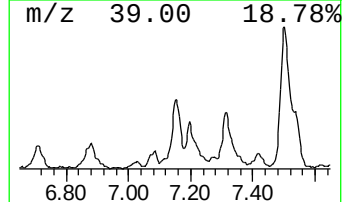
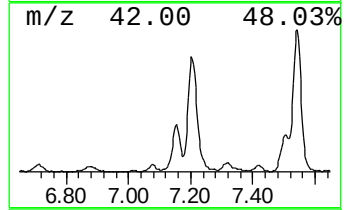
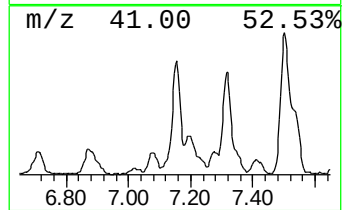
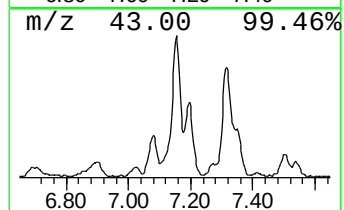
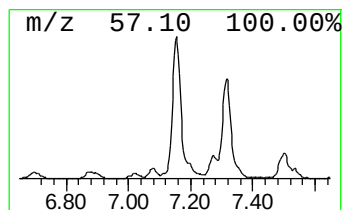
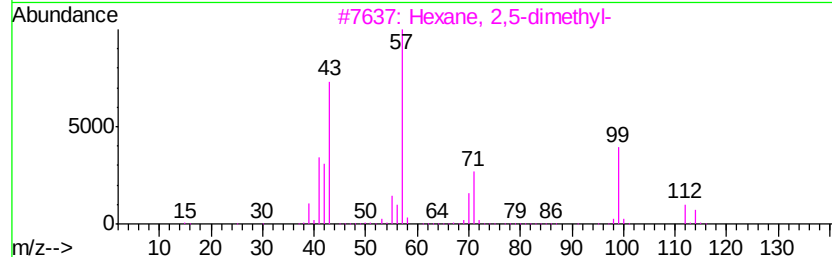
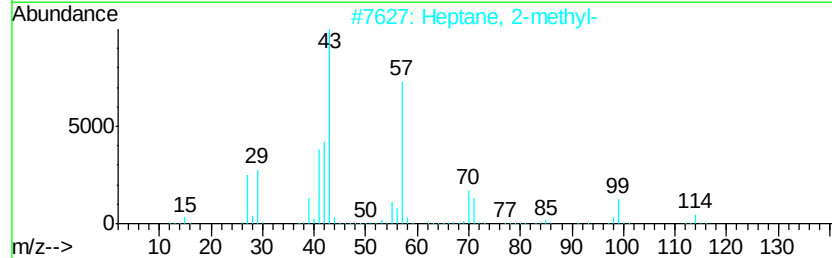
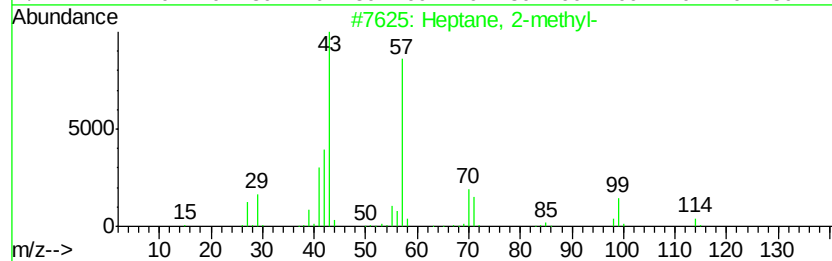
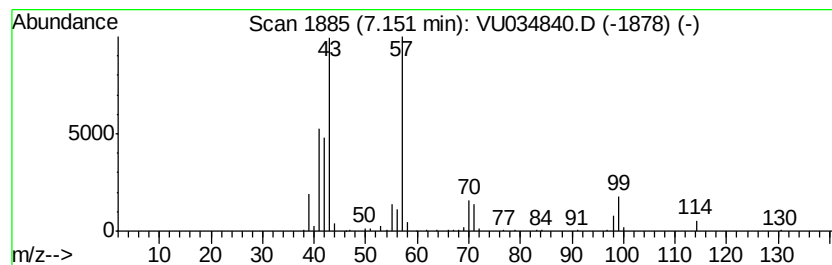
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 45

| R.T. | EstConc    | Area   | Relative to ISTD    | R.T. |
|------|------------|--------|---------------------|------|
| 7.15 | 10.60 ug/L | 259267 | 1,4-Difluorobenzene | 5.86 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Heptane, 2-methyl-    | 114 | C8H18   | 000592-27-8 | 91   |
| 2    |    | Heptane, 2-methyl-    | 114 | C8H18   | 000592-27-8 | 91   |
| 3    |    | Hexane, 2,5-dimethyl- | 114 | C8H18   | 000592-13-2 | 91   |
| 4    |    | Heptane, 2-methyl-    | 114 | C8H18   | 000592-27-8 | 87   |
| 5    |    | Hexane, 2,5-dimethyl- | 114 | C8H18   | 000592-13-2 | 87   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

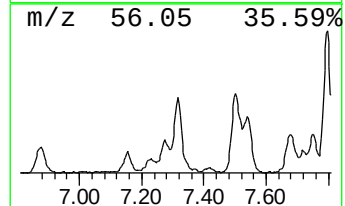
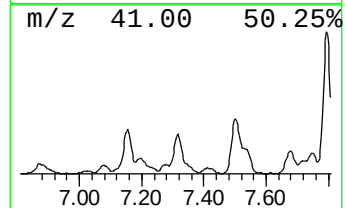
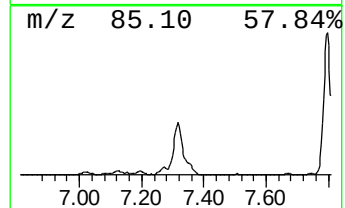
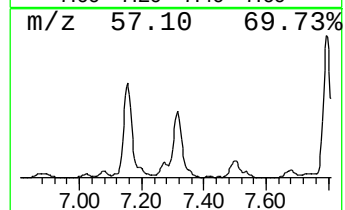
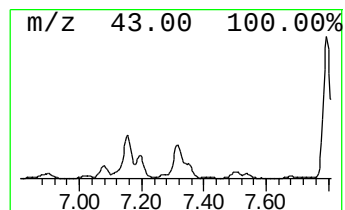
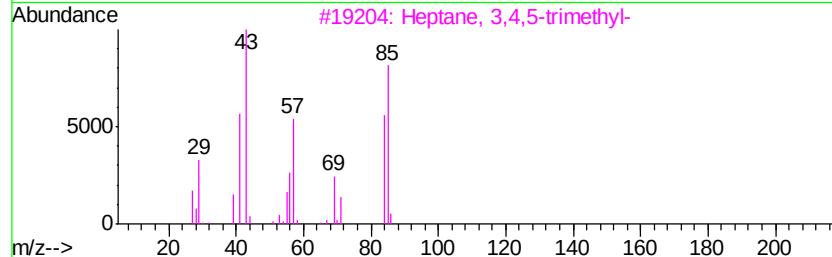
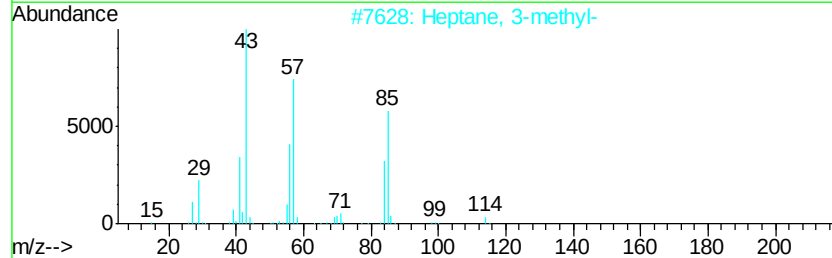
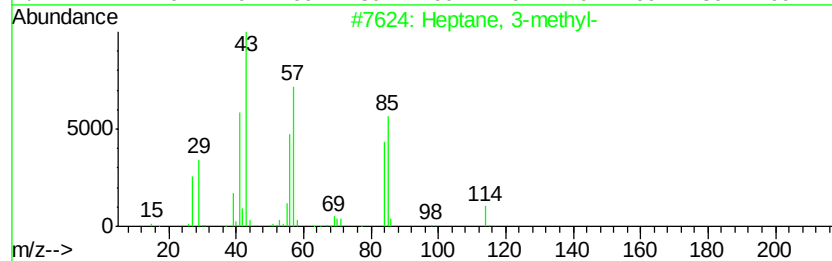
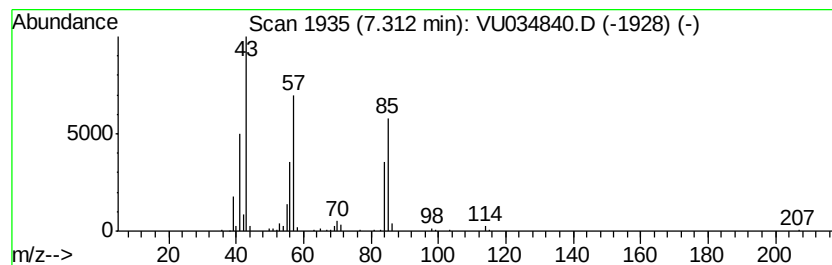
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 42

| R.T. | EstConc    | Area   | Relative to ISTD    | R.T. |
|------|------------|--------|---------------------|------|
| 7.31 | 14.75 ug/L | 360653 | 1,4-Difluorobenzene | 5.86 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptane, 3-methyl-        | 114 | C8H18   | 000589-81-1 | 91   |
| 2    |    | Heptane, 3-methyl-        | 114 | C8H18   | 000589-81-1 | 91   |
| 3    |    | Heptane, 3,4,5-trimethyl- | 142 | C10H22  | 020278-89-1 | 78   |
| 4    |    | Heptane, 3,4,5-trimethyl- | 142 | C10H22  | 020278-89-1 | 64   |
| 5    |    | Hexane, 2-methyl-         | 100 | C7H16   | 000591-76-4 | 59   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

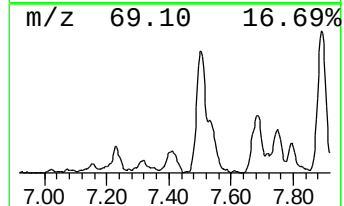
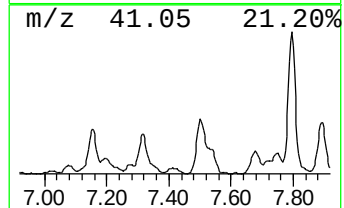
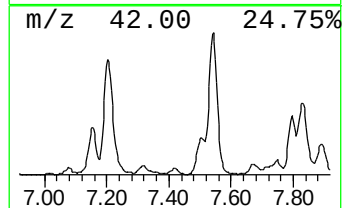
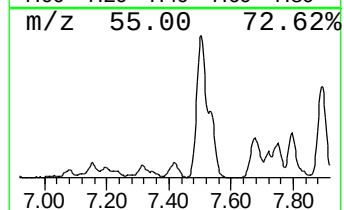
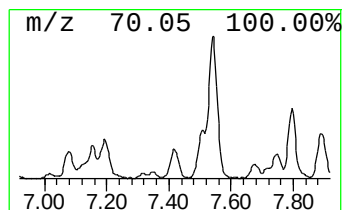
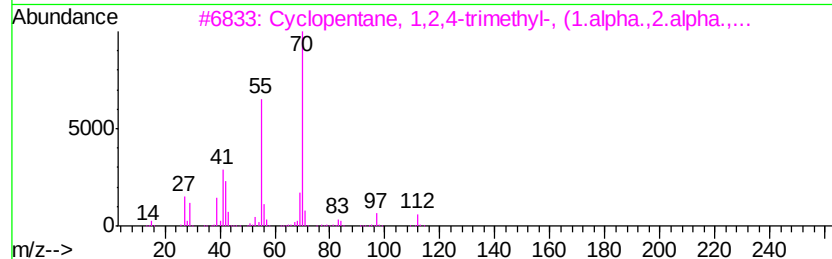
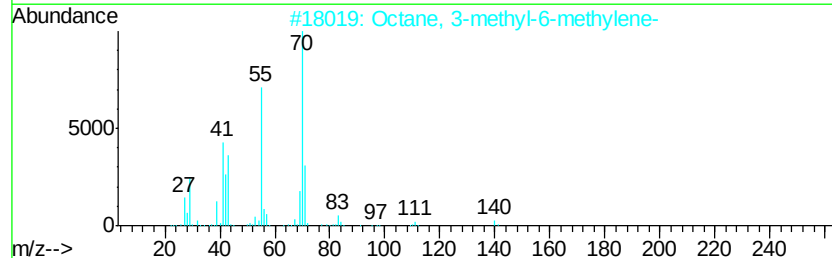
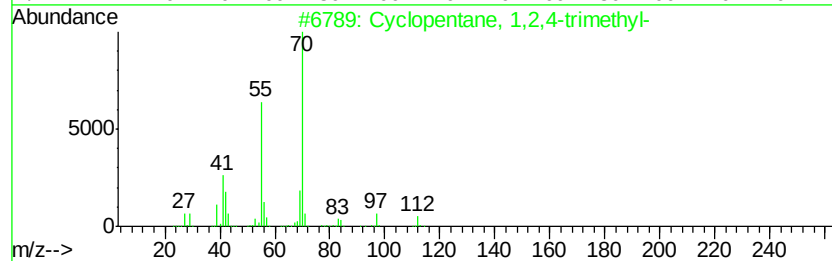
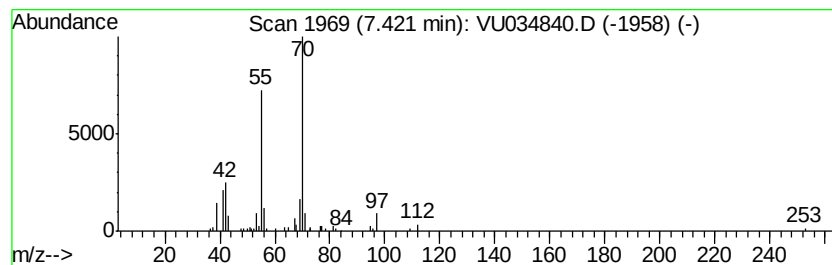
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Cyclic7.42 Concentration Rank 56

| R.T. | EstConc   | Area  | Relative to ISTD    | R.T. |
|------|-----------|-------|---------------------|------|
| 7.42 | 2.89 ug/L | 70697 | 1,4-Difluorobenzene | 5.86 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopentane, 1,2,4-trimethyl-      | 112 | C8H16   | 002815-58-9 | 80   |
| 2    |    | Octane, 3-methyl-6-methylene-       | 140 | C10H20  | 074630-07-2 | 78   |
| 3    |    | Cyclopentane, 1,2,4-trimethyl-, ... | 112 | C8H16   | 004850-28-6 | 78   |
| 4    |    | 2-Butene, 2-methyl-                 | 70  | C5H10   | 000513-35-9 | 72   |
| 5    |    | Hexane, 2-methyl-4-methylene-       | 112 | C8H16   | 003404-80-6 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
 Data File : VU034840.D  
 Acq On : 26 Sep 2019 12:48  
 Operator : JC/SP  
 Sample : K4984-08  
 Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleID :  
 BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
 Quant Title : VOC Analysis

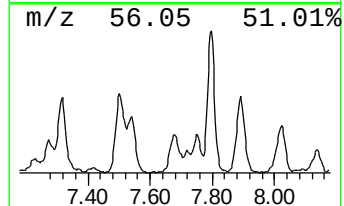
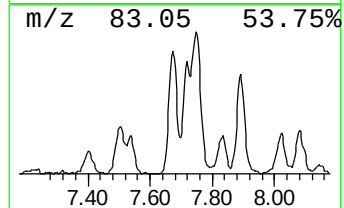
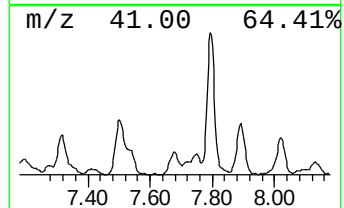
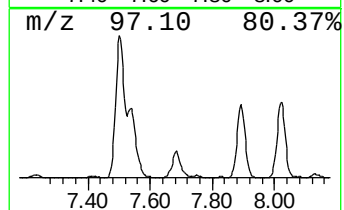
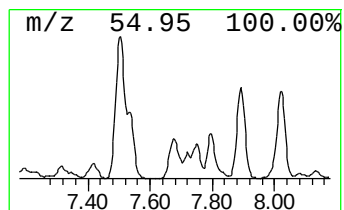
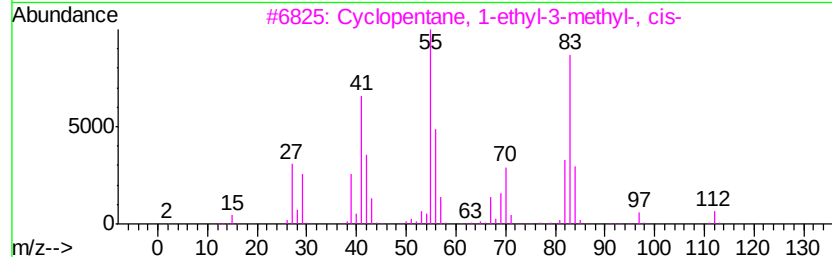
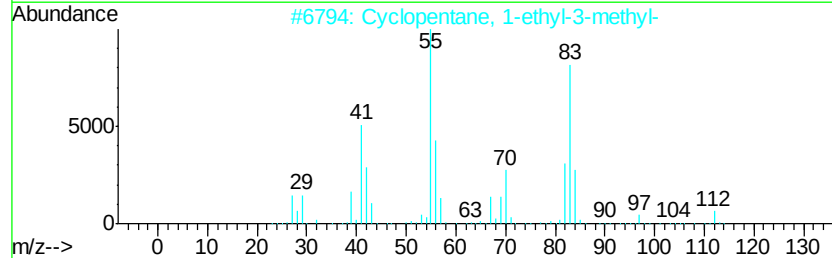
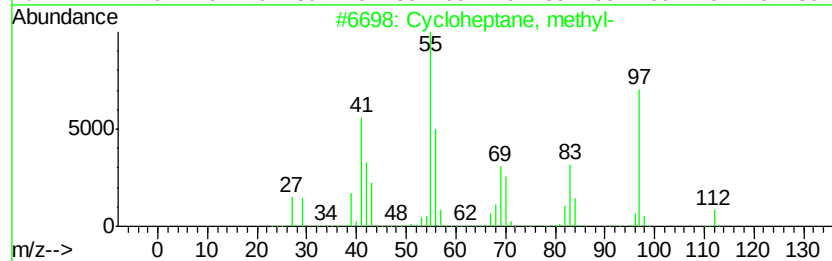
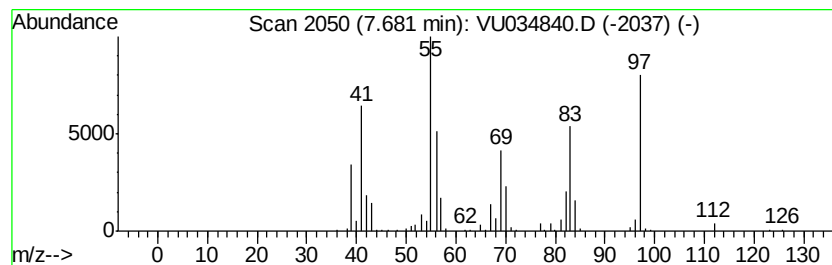
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 11 (DEL) Alkane: Cyclic7.68 Concentration Rank 47

| R.T. | EstConc   | Area   | Relative to ISTD | R.T. |
|------|-----------|--------|------------------|------|
| 7.68 | 7.77 ug/L | 230694 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cycloheptane, methyl-               | 112 | C8H16   | 004126-78-7 | 72   |
| 2    |    | Cyclopentane, 1-ethyl-3-methyl-     | 112 | C8H16   | 003726-47-4 | 70   |
| 3    |    | Cyclopentane, 1-ethyl-3-methyl-,... | 112 | C8H16   | 002613-66-3 | 64   |
| 4    |    | Cyclopentane, 1-ethyl-3-methyl-,... | 112 | C8H16   | 002613-65-2 | 52   |
| 5    |    | Cyclopentane, 1-ethyl-2-methyl-,... | 112 | C8H16   | 000930-89-2 | 43   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

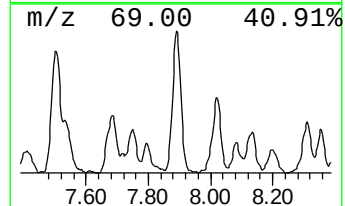
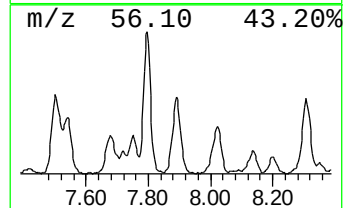
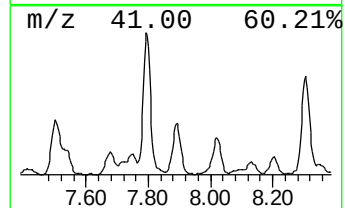
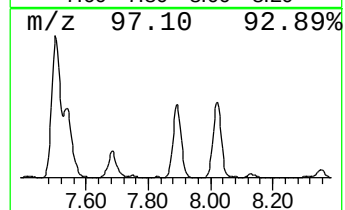
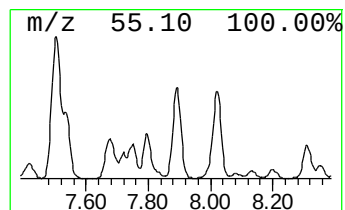
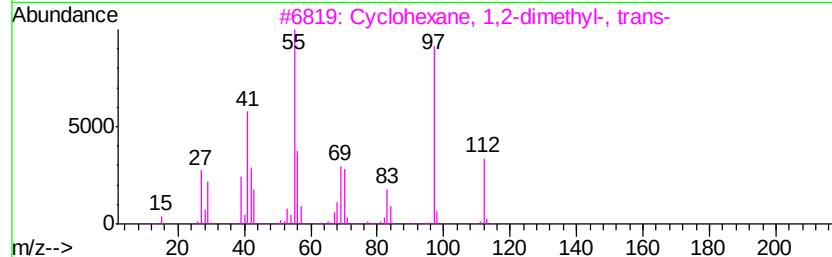
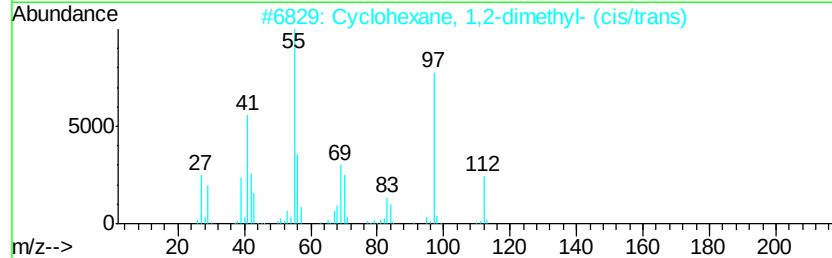
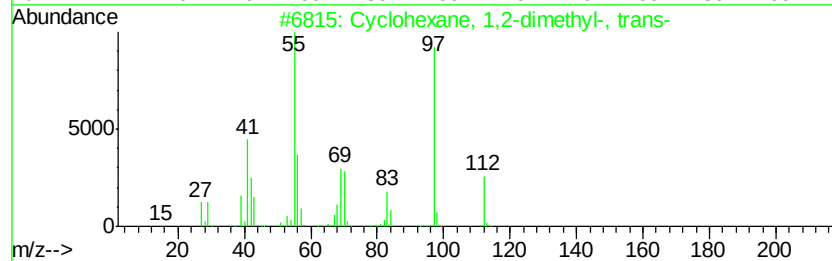
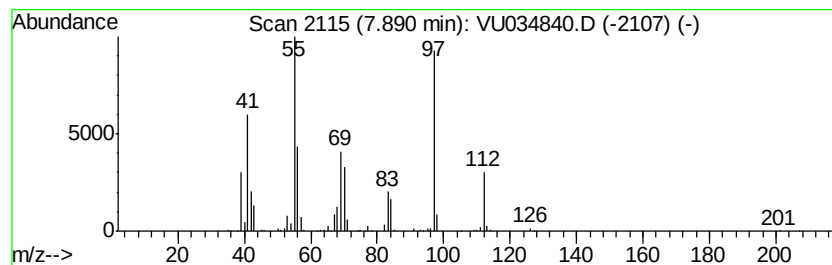
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 (DEL) Alkane: Cyclic7.89 Concentration Rank 41

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 7.89 | 17.20 ug/L | 510322 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1,2-dimethyl-, trans-  | 112 | C8H16   | 006876-23-9 | 95   |
| 2    |    | Cyclohexane, 1,2-dimethyl- (cis/... | 112 | C8H16   | 000583-57-3 | 94   |
| 3    |    | Cyclohexane, 1,2-dimethyl-, trans-  | 112 | C8H16   | 006876-23-9 | 93   |
| 4    |    | Cyclohexane, 1,2-dimethyl- (cis/... | 112 | C8H16   | 000583-57-3 | 90   |
| 5    |    | Cyclohexane, 1,2-dimethyl-, trans-  | 112 | C8H16   | 006876-23-9 | 87   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

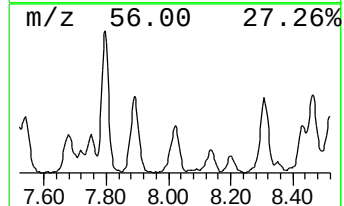
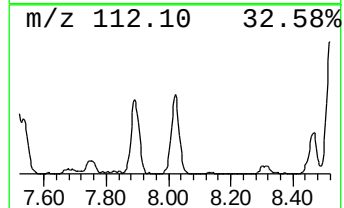
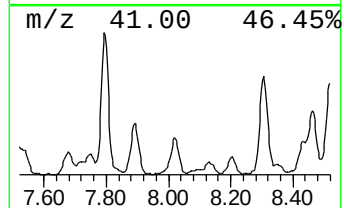
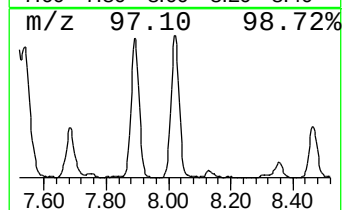
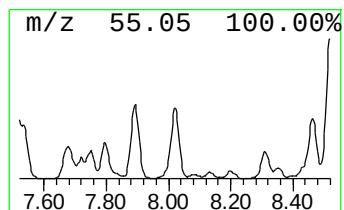
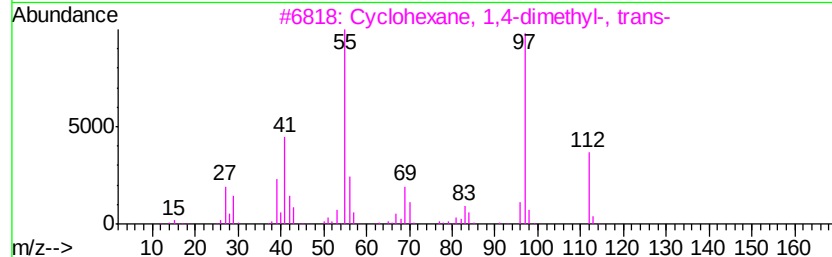
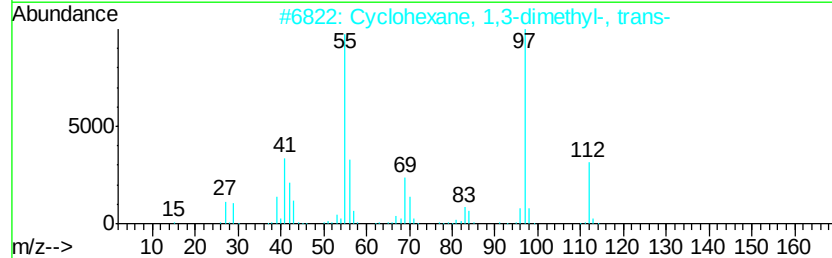
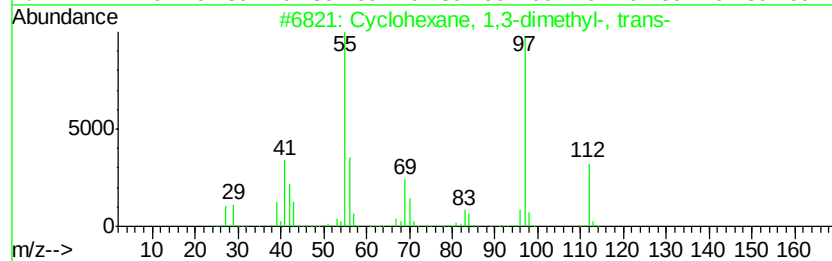
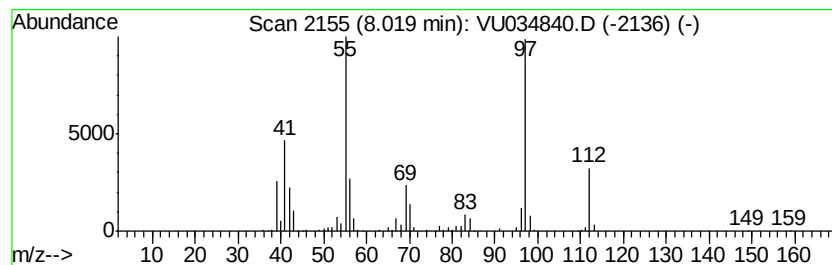
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Cyclic8.02 Concentration Rank 43

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 8.02 | 13.48 ug/L | 400083 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16   | 002207-03-6 | 97   |
| 2    |    | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16   | 002207-03-6 | 97   |
| 3    |    | Cyclohexane, 1,4-dimethyl-, trans- | 112 | C8H16   | 002207-04-7 | 97   |
| 4    |    | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16   | 002207-03-6 | 97   |
| 5    |    | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16   | 002207-03-6 | 96   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

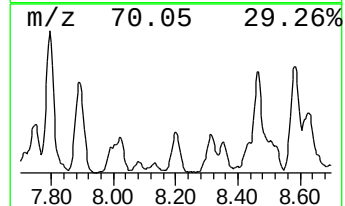
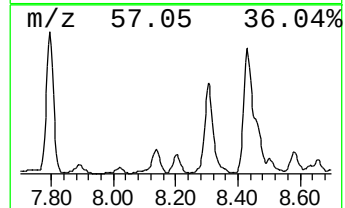
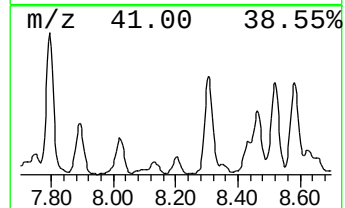
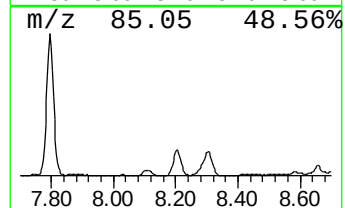
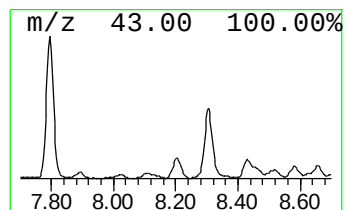
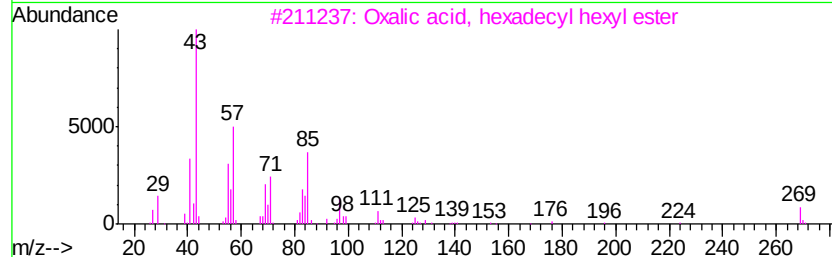
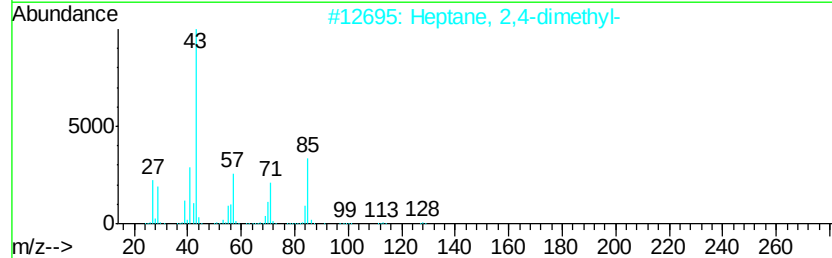
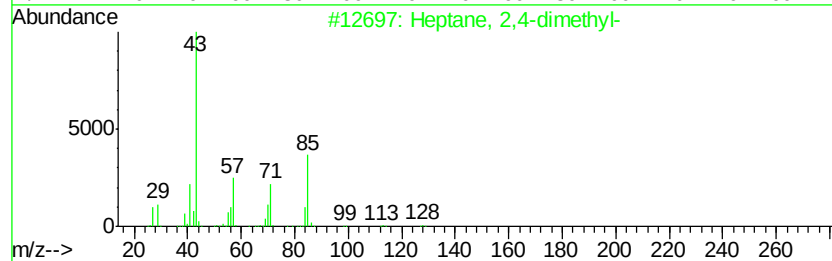
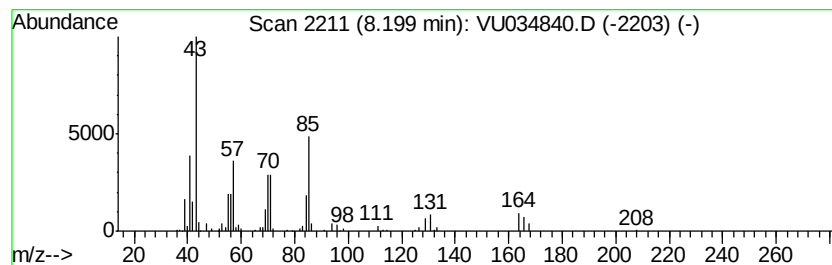
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 50

| R.T. | EstConc   | Area   | Relative to ISTD | R.T. |
|------|-----------|--------|------------------|------|
| 8.20 | 6.26 ug/L | 185790 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|----|------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Heptane, 2,4-dimethyl-             | 128 | C9H20    | 002213-23-2  | 58   |
| 2    |    | Heptane, 2,4-dimethyl-             | 128 | C9H20    | 002213-23-2  | 55   |
| 3    |    | Oxalic acid, hexadecyl hexyl ester | 398 | C24H46O4 | 1000309-25-0 | 50   |
| 4    |    | Octane, 4-methyl-                  | 128 | C9H20    | 002216-34-4  | 46   |
| 5    |    | Heptane, 2,4-dimethyl-             | 128 | C9H20    | 002213-23-2  | 46   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

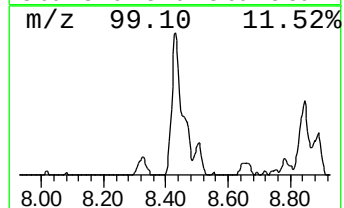
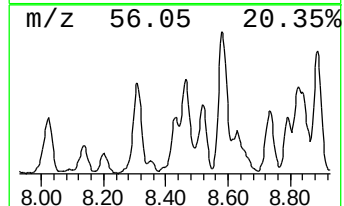
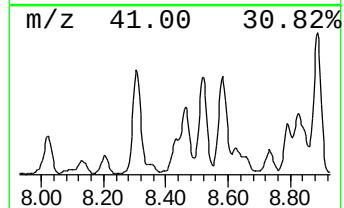
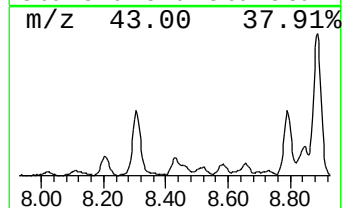
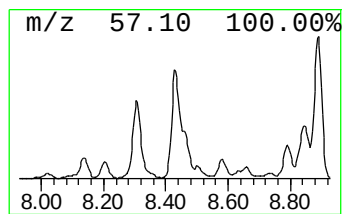
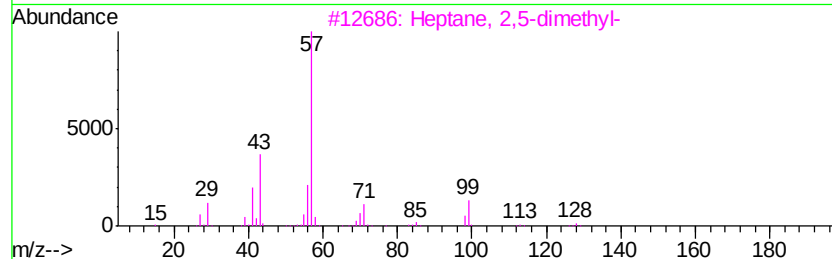
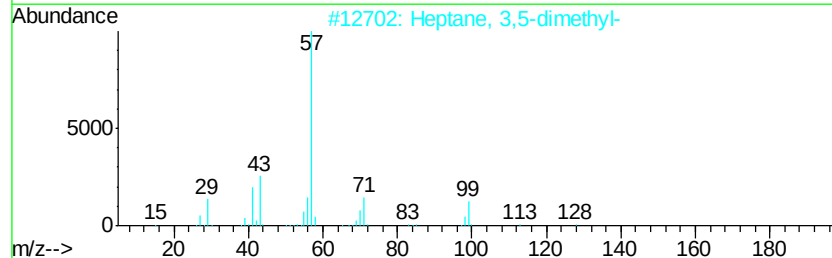
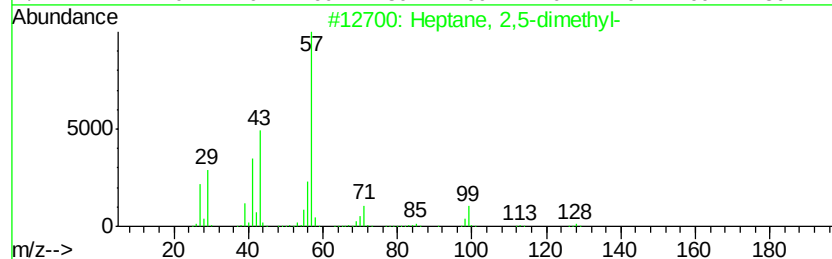
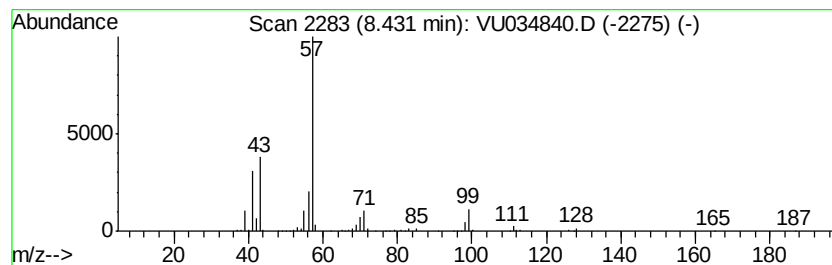
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 49

| R.T. | EstConc   | Area   | Relative to ISTD | R.T. |
|------|-----------|--------|------------------|------|
| 8.43 | 6.40 ug/L | 190015 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 90   |
| 2    |    | Heptane, 3,5-dimethyl- | 128 | C9H20   | 000926-82-9 | 90   |
| 3    |    | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 90   |
| 4    |    | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 87   |
| 5    |    | Octane, 3-methyl-      | 128 | C9H20   | 002216-33-3 | 83   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

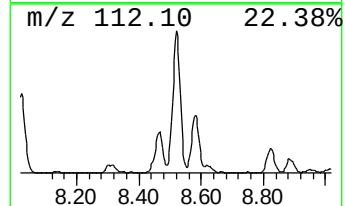
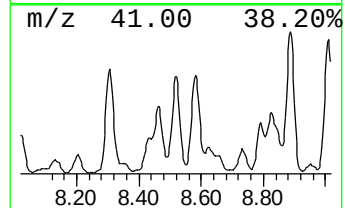
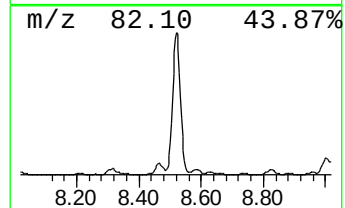
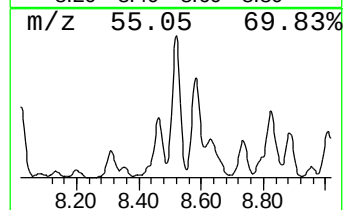
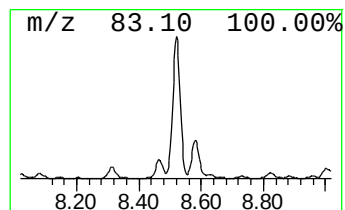
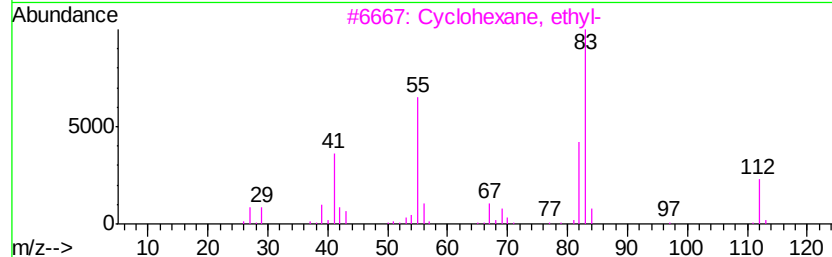
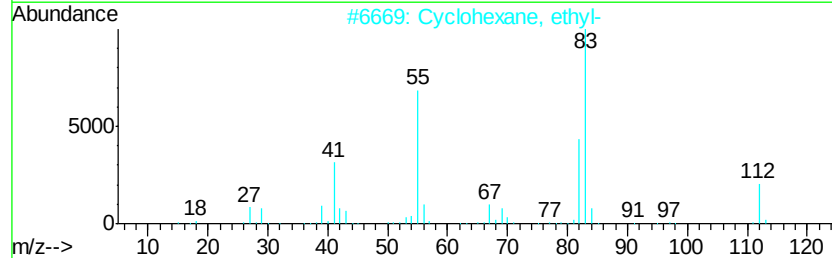
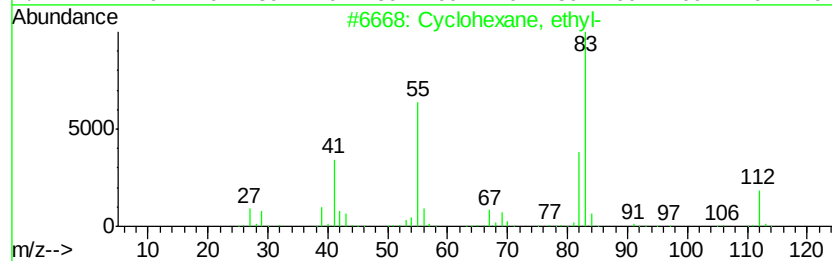
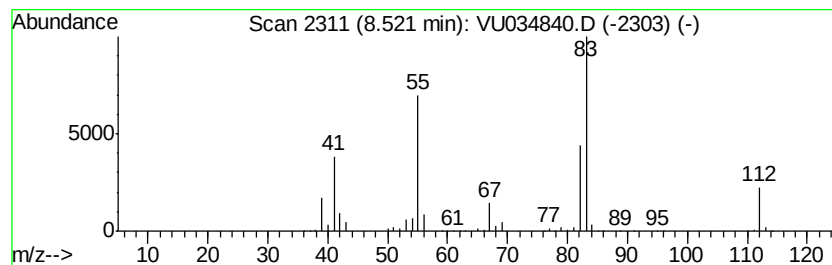
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Cyclic8.52 Concentration Rank 39

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 8.52 | 24.20 ug/L | 718363 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, ethyl- | 112 | C8H16   | 001678-91-7 | 91   |
| 2    |    | Cyclohexane, ethyl- | 112 | C8H16   | 001678-91-7 | 91   |
| 3    |    | Cyclohexane, ethyl- | 112 | C8H16   | 001678-91-7 | 91   |
| 4    |    | Cyclohexane, ethyl- | 112 | C8H16   | 001678-91-7 | 91   |
| 5    |    | Cyclohexane, ethyl- | 112 | C8H16   | 001678-91-7 | 87   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

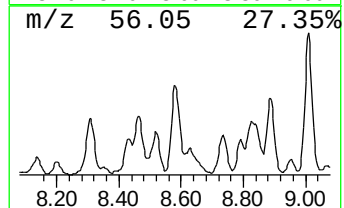
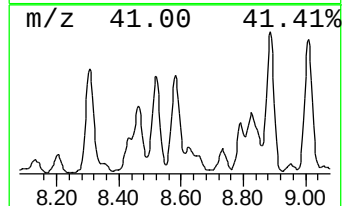
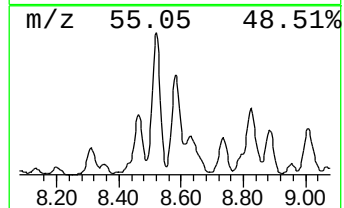
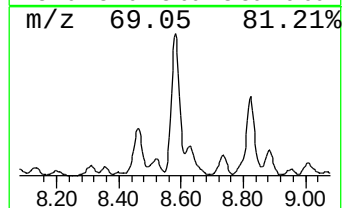
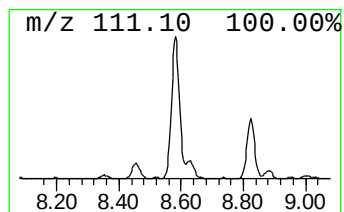
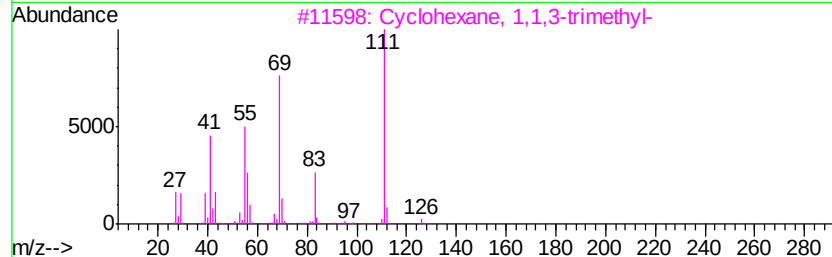
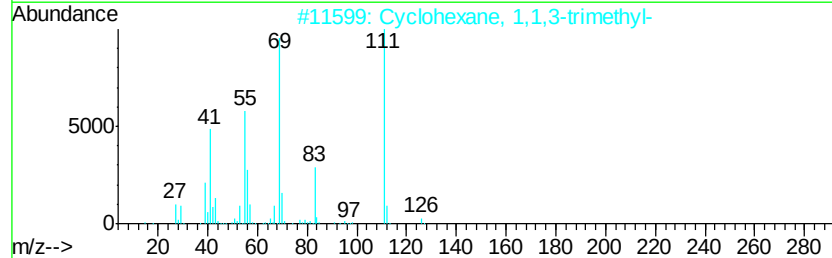
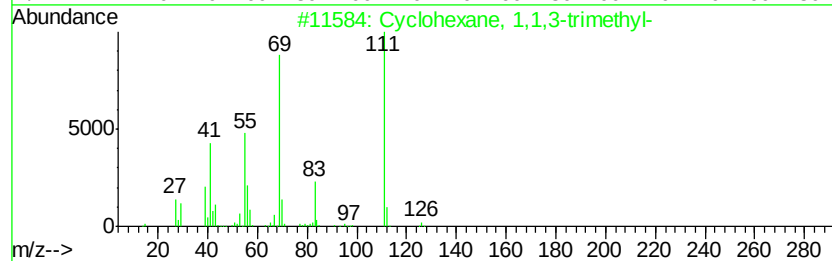
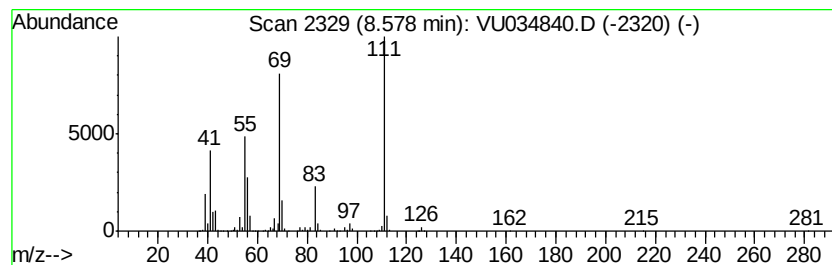
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 (DEL) Alkane: Cyclic8.58 Concentration Rank 33

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 8.58 | 29.89 ug/L | 886971 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1,1,3-trimethyl-        | 126 | C9H18   | 003073-66-3 | 97   |
| 2    |    | Cyclohexane, 1,1,3-trimethyl-        | 126 | C9H18   | 003073-66-3 | 97   |
| 3    |    | Cyclohexane, 1,1,3-trimethyl-        | 126 | C9H18   | 003073-66-3 | 95   |
| 4    |    | Cyclohexane, 1,3,5-trimethyl-, (...) | 126 | C9H18   | 001795-27-3 | 72   |
| 5    |    | Cyclohexane, 1,3,5-trimethyl-        | 126 | C9H18   | 001839-63-0 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
 Data File : VU034840.D  
 Acq On : 26 Sep 2019 12:48  
 Operator : JC/SP  
 Sample : K4984-08  
 Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
 Quant Title : VOC Analysis

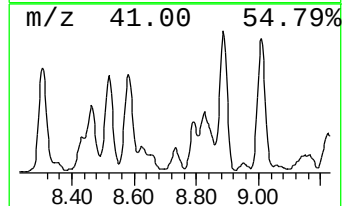
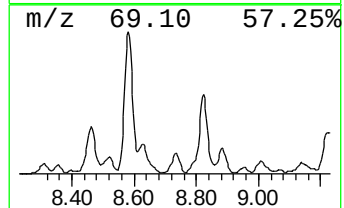
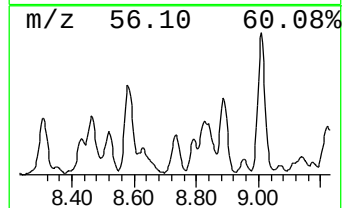
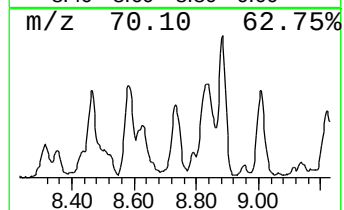
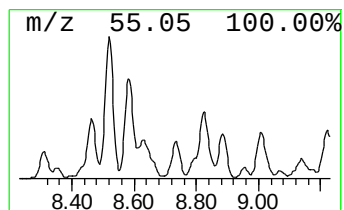
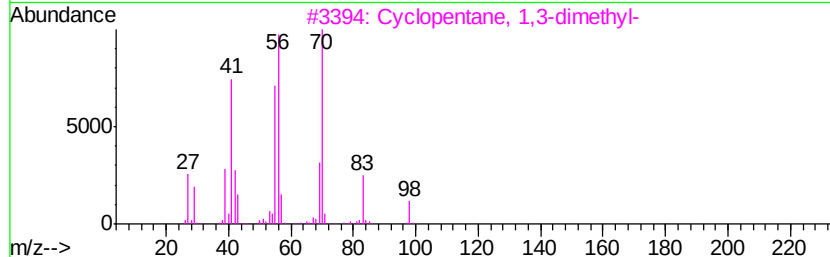
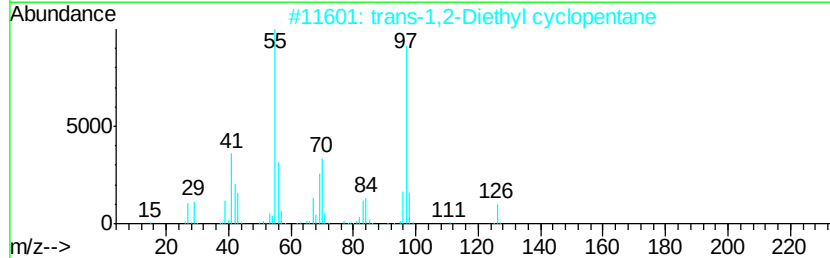
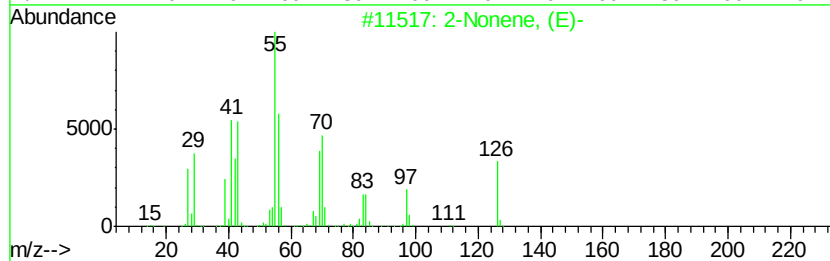
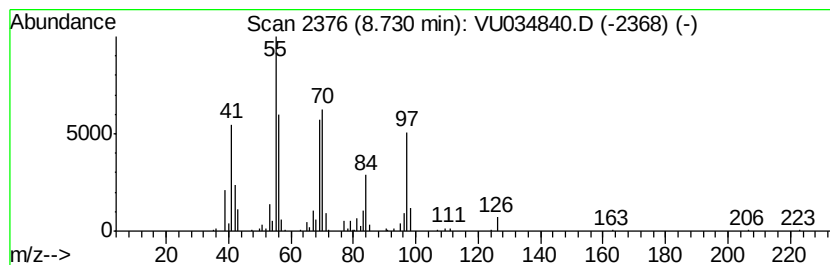
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 18 (DEL) Alkane: Cyclic8.73 Concentration Rank 48

| R.T. | EstConc   | Area   | Relative to ISTD | R.T. |
|------|-----------|--------|------------------|------|
| 8.73 | 6.68 ug/L | 198168 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#         | Qual |
|------|----|--------------------------------|-----|---------|--------------|------|
| 1    | 5  | 2-Nonene, (E)-                 | 126 | C9H18   | 006434-78-2  | 74   |
| 2    |    | trans-1,2-Diethyl cyclopentane | 126 | C9H18   | 000932-40-1  | 46   |
| 3    |    | Cyclopentane, 1,3-dimethyl-    | 98  | C7H14   | 002453-00-1  | 46   |
| 4    |    | Cyclobutanone, 2,2-dimethyl-   | 98  | C6H10O  | 001192-14-9  | 46   |
| 5    |    | trans-1,3-Diethylcyclopentane  | 126 | C9H18   | 1000113-87-1 | 43   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

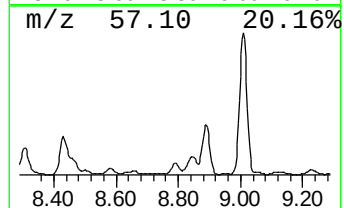
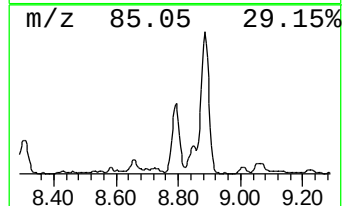
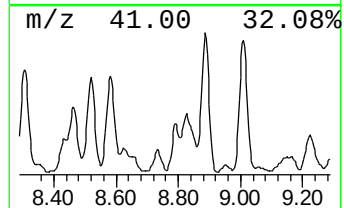
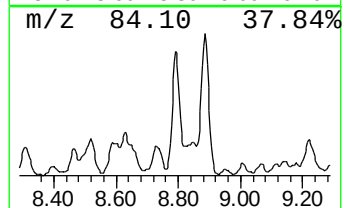
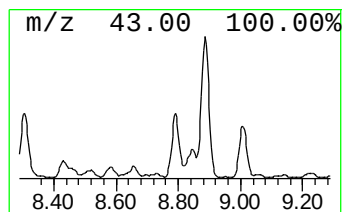
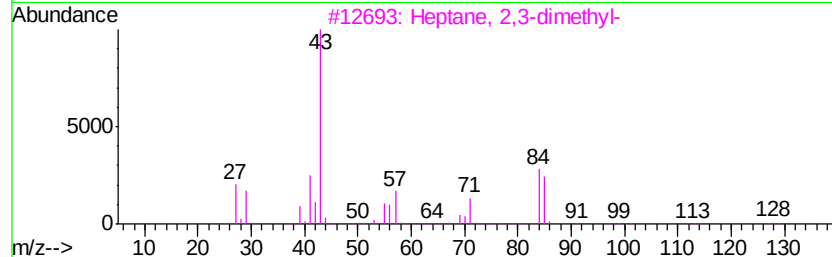
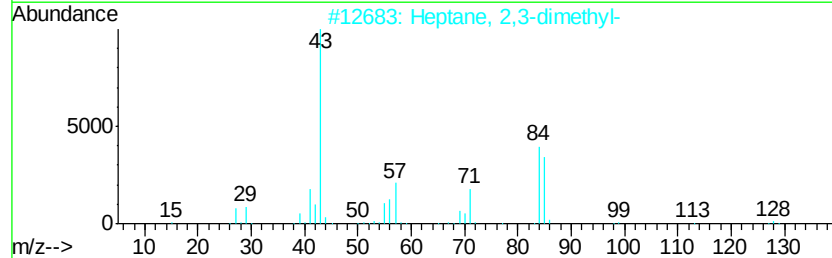
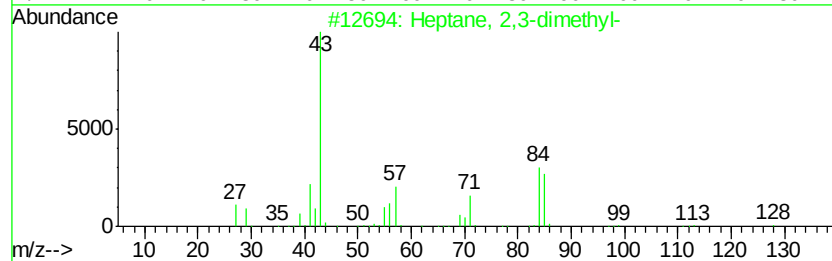
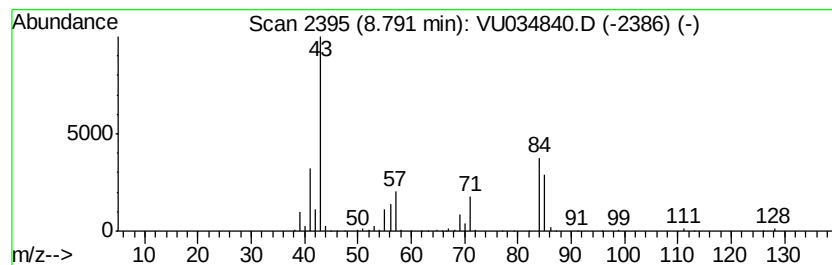
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 44

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 8.79 | 12.69 ug/L | 376618 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptane, 2,3-dimethyl-    | 128 | C9H20   | 003074-71-3 | 94   |
| 2    |    |   | Heptane, 2,3-dimethyl-    | 128 | C9H20   | 003074-71-3 | 91   |
| 3    |    |   | Heptane, 2,3-dimethyl-    | 128 | C9H20   | 003074-71-3 | 91   |
| 4    |    |   | Hexane, 2,3,5-trimethyl-  | 128 | C9H20   | 001069-53-0 | 87   |
| 5    |    |   | Hexane, 3-ethyl-2-methyl- | 128 | C9H20   | 016789-46-1 | 86   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

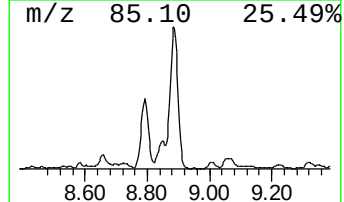
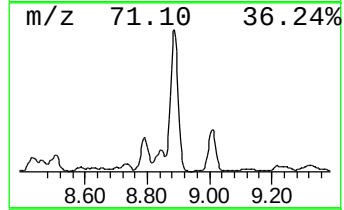
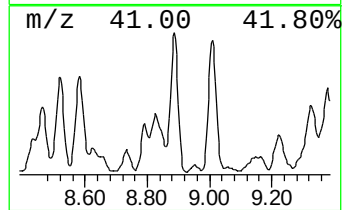
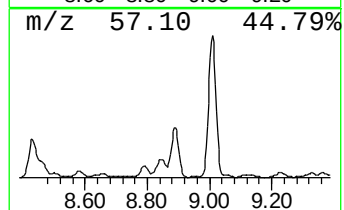
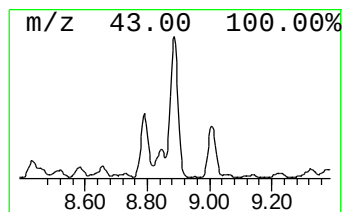
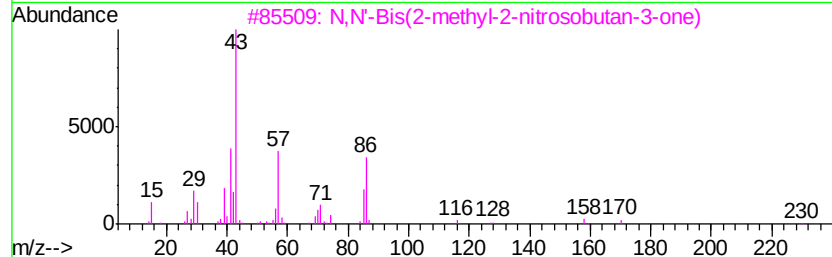
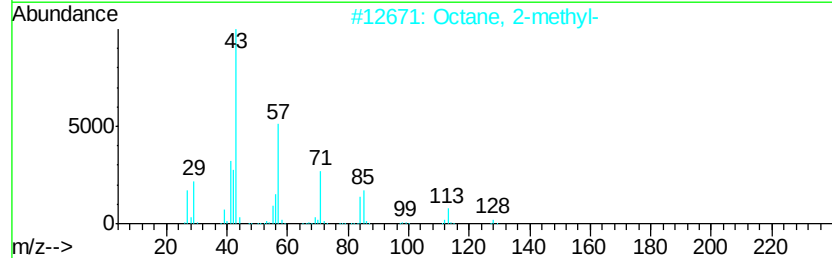
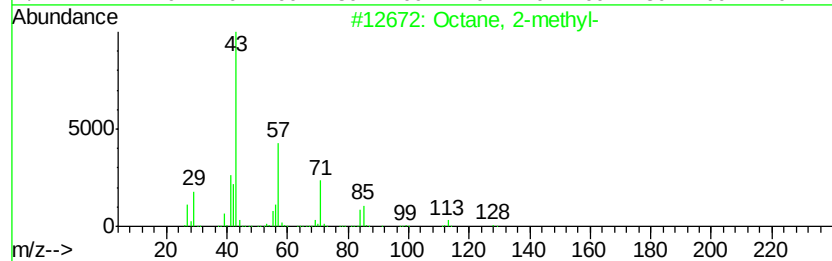
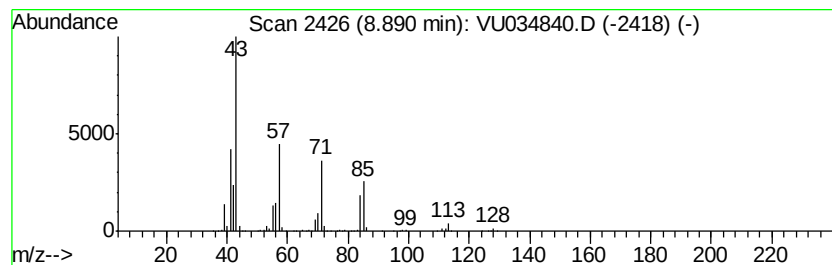
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 26

| R.T. | EstConc    | Area    | Relative to ISTD | R.T. |
|------|------------|---------|------------------|------|
| 8.89 | 36.79 ug/L | 1092010 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Octane, 2-methyl-                   | 128 | C9H20      | 003221-61-2 | 64   |
| 2    |    | Octane, 2-methyl-                   | 128 | C9H20      | 003221-61-2 | 64   |
| 3    |    | N,N'-Bis(2-methyl-2-nitrosobutan... | 230 | C10H18N2O4 | 034946-73-1 | 64   |
| 4    |    | Ether, hexyl pentyl                 | 172 | C11H24O    | 032357-83-8 | 59   |
| 5    |    | Undecane, 4,7-dimethyl-             | 184 | C13H28     | 017301-32-5 | 53   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

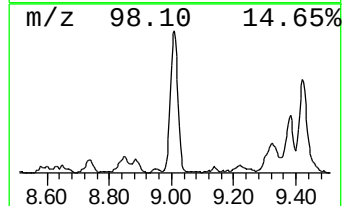
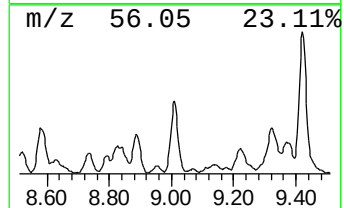
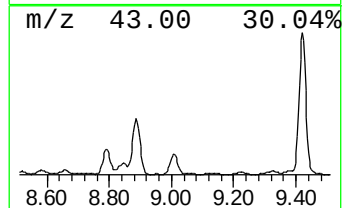
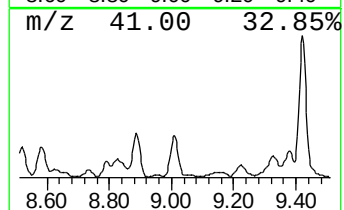
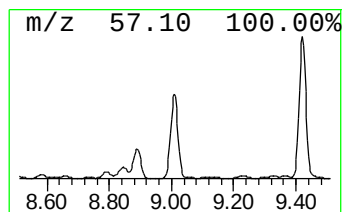
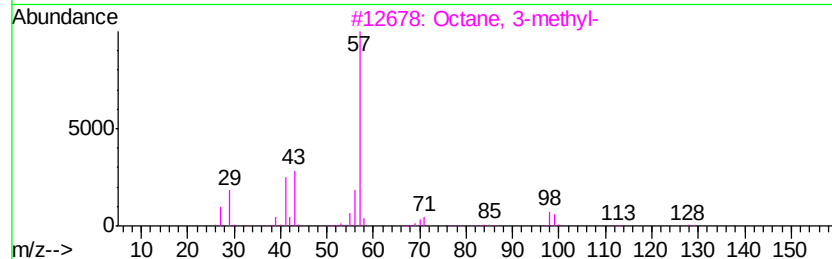
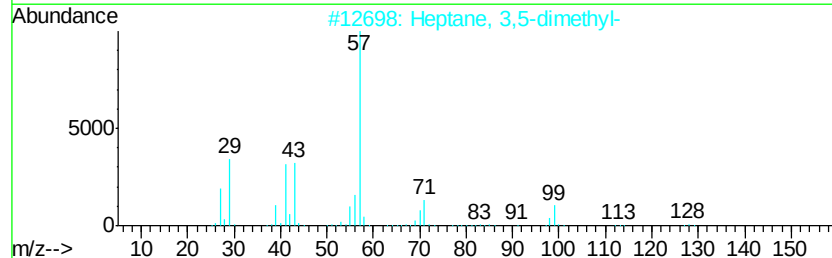
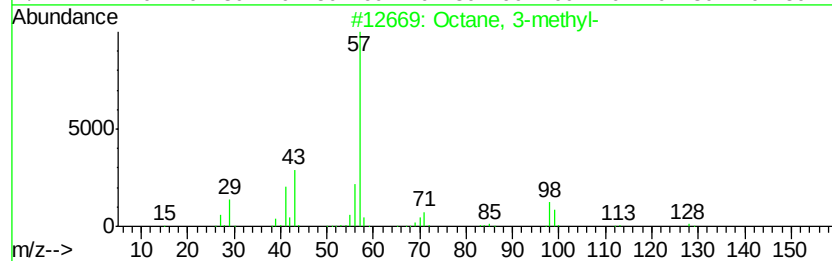
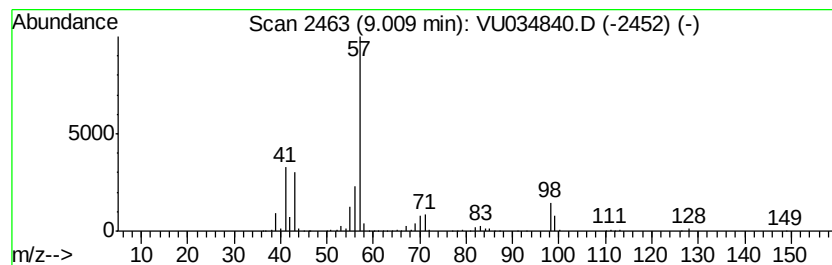
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 30

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 9.01 | 32.73 ug/L | 971426 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 3-methyl-      | 128 | C9H20   | 002216-33-3 | 91   |
| 2    |    | Heptane, 3,5-dimethyl- | 128 | C9H20   | 000926-82-9 | 64   |
| 3    |    | Octane, 3-methyl-      | 128 | C9H20   | 002216-33-3 | 64   |
| 4    |    | Decane, 2,5-dimethyl-  | 170 | C12H26  | 017312-50-4 | 59   |
| 5    |    | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 59   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

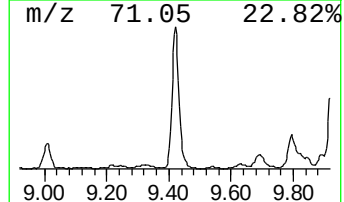
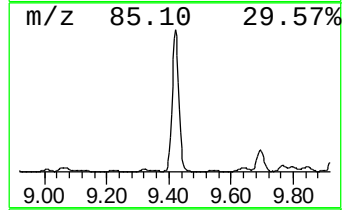
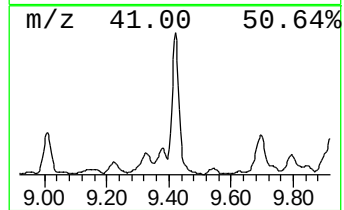
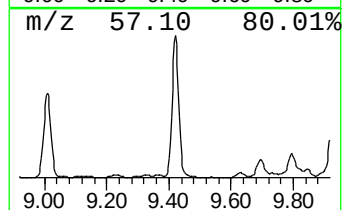
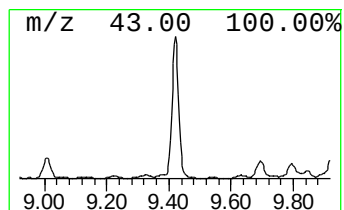
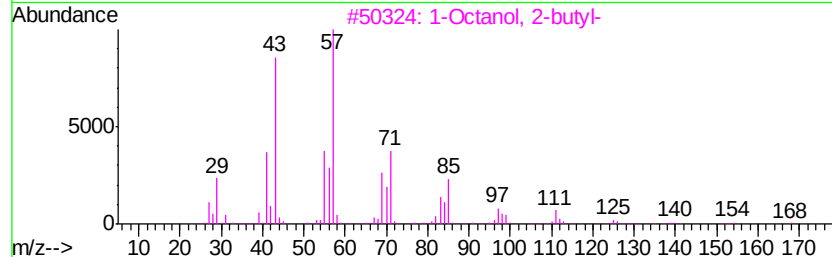
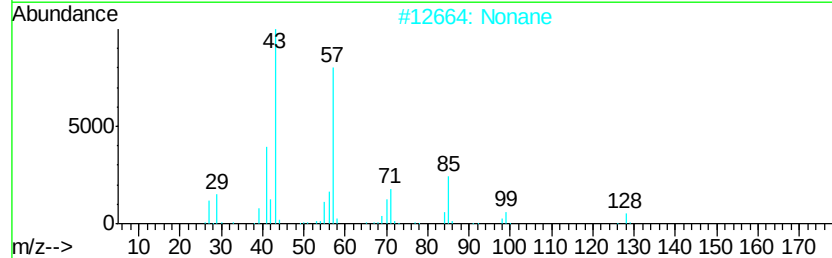
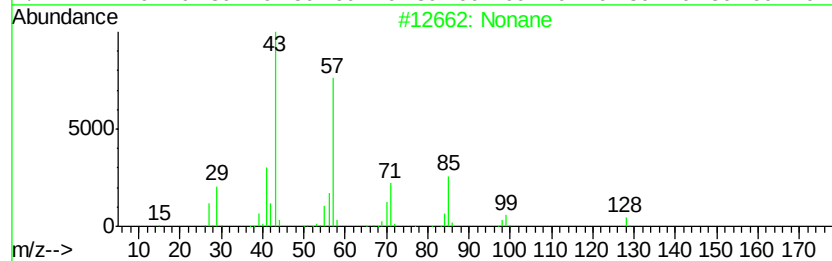
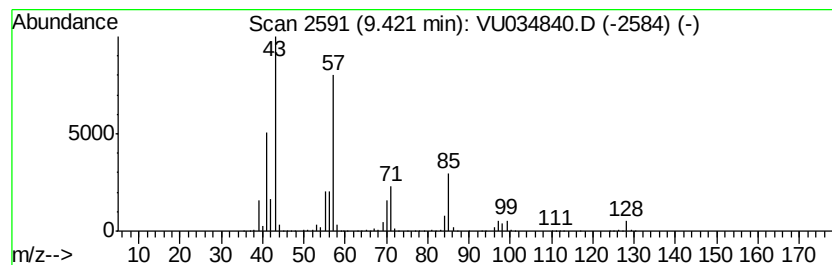
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T. | EstConc     | Area    | Relative to ISTD | R.T. |
|------|-------------|---------|------------------|------|
| 9.42 | 110.34 ug/L | 3274570 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane                | 128 | C9H20   | 000111-84-2 | 97   |
| 2    |    | Nonane                | 128 | C9H20   | 000111-84-2 | 94   |
| 3    |    | 1-Octanol, 2-butyl-   | 186 | C12H26O | 003913-02-8 | 72   |
| 4    |    | Nonane                | 128 | C9H20   | 000111-84-2 | 58   |
| 5    |    | Hexane, 2,4-dimethyl- | 114 | C8H18   | 000589-43-5 | 53   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

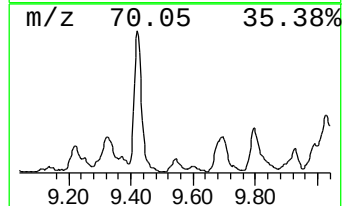
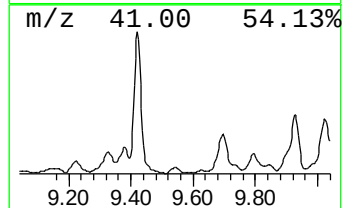
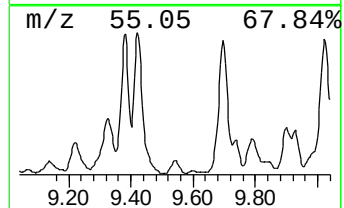
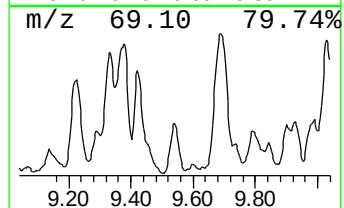
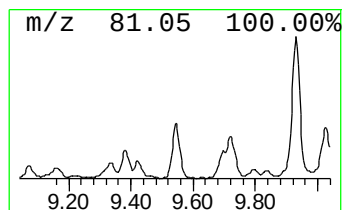
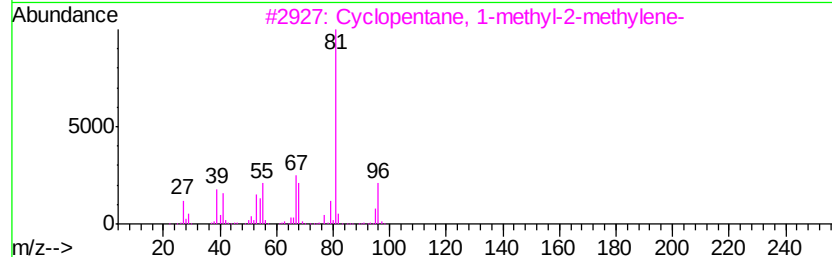
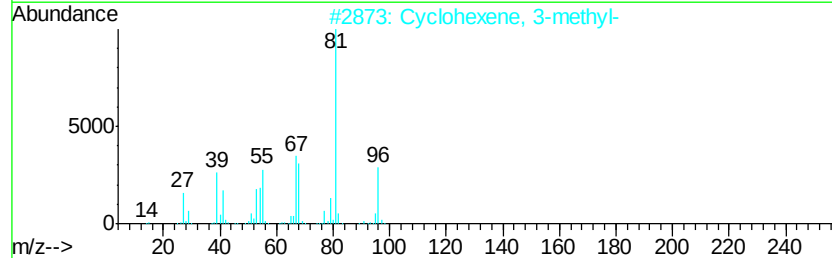
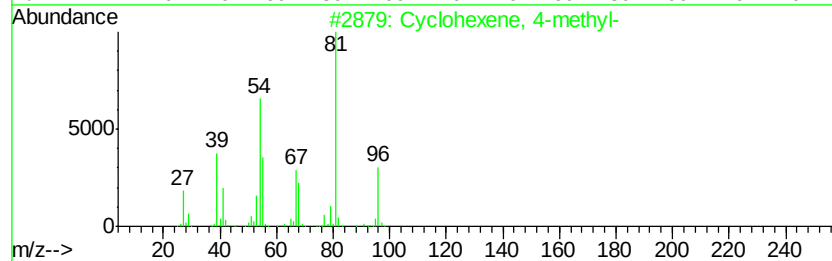
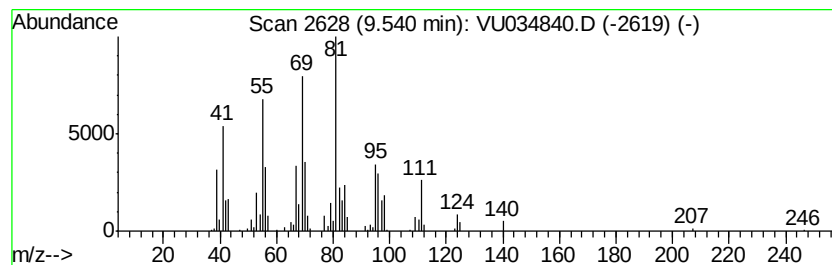
Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 23 unknown-01 Concentration Rank 46

| R.T.    | EstConc   | Area                                | Relative to ISTD | R.T.    |             |      |
|---------|-----------|-------------------------------------|------------------|---------|-------------|------|
| 9.54    | 8.57 ug/L | 254207                              | Chlorobenzene-d5 | 9.07    |             |      |
| Hit# of | 5         | Tentative ID                        | MW               | MolForm | CAS#        | Qual |
| 1       |           | Cyclohexene, 4-methyl-              | 96               | C7H12   | 000591-47-9 | 46   |
| 2       |           | Cyclohexene, 3-methyl-              | 96               | C7H12   | 000591-48-0 | 46   |
| 3       |           | Cyclopentane, 1-methyl-2-methylene- | 96               | C7H12   | 041158-41-2 | 46   |
| 4       |           | Cyclohexene, 3-methyl-              | 96               | C7H12   | 000591-48-0 | 46   |
| 5       |           | Cyclopropane, trimethylmethylene-   | 96               | C7H12   | 034462-28-7 | 43   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

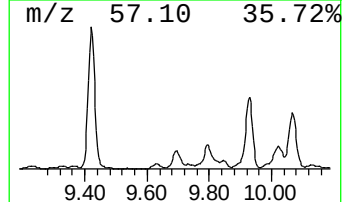
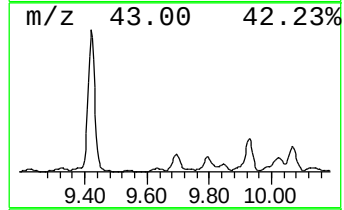
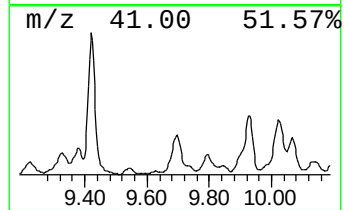
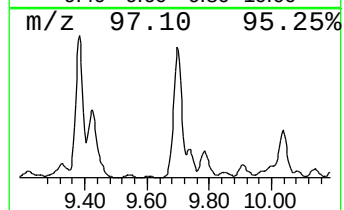
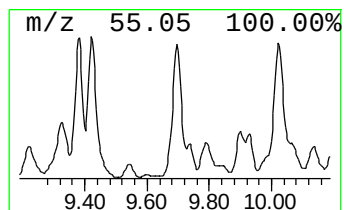
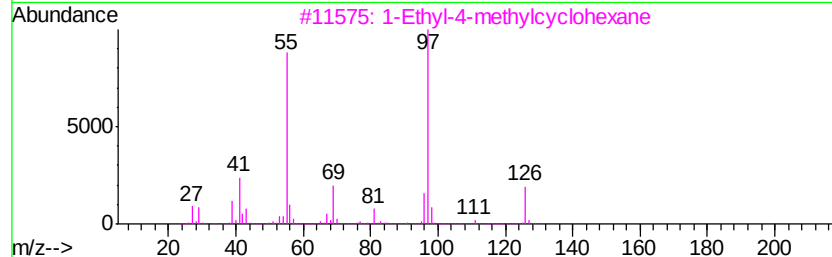
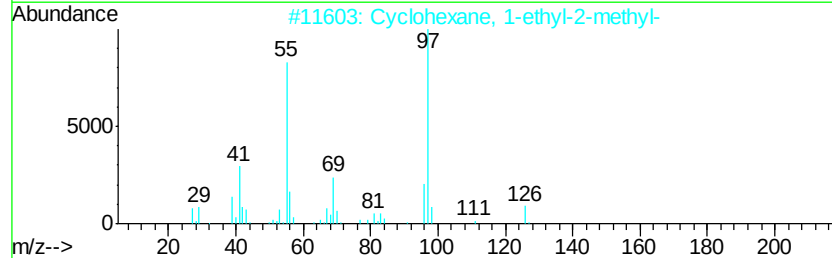
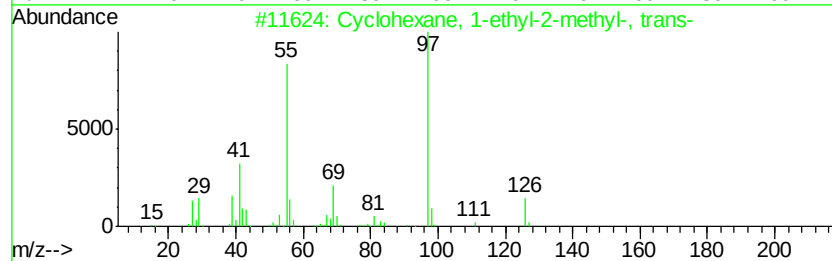
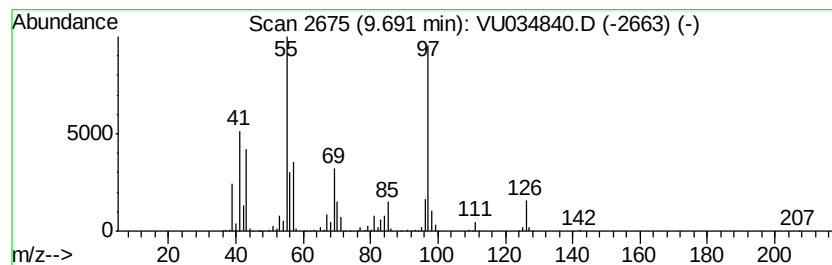
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 (DEL) Alkane: Cyclic9.69 Concentration Rank 19

| R.T. | EstConc    | Area    | Relative to ISTD | R.T. |
|------|------------|---------|------------------|------|
| 9.69 | 47.79 ug/L | 1418420 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Cyclohexane, 1-ethyl-2-methyl-, ... | 126 | C9H18     | 004923-78-8  | 81   |
| 2    |    | Cyclohexane, 1-ethyl-2-methyl-      | 126 | C9H18     | 003728-54-9  | 81   |
| 3    |    | 1-Ethyl-4-methylcyclohexane         | 126 | C9H18     | 003728-56-1  | 76   |
| 4    |    | 3,5-Dimethyl-3-heptene              | 126 | C9H18     | 059643-68-4  | 64   |
| 5    |    | Sulfurous acid, cyclohexylmethyl... | 402 | C23H46O3S | 1000309-22-4 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

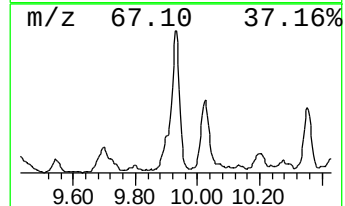
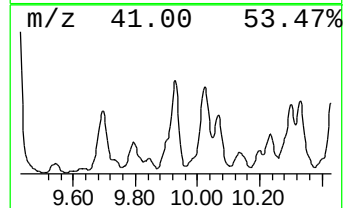
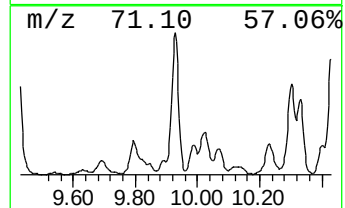
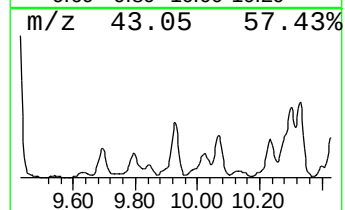
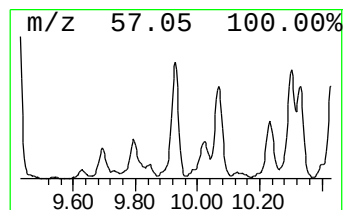
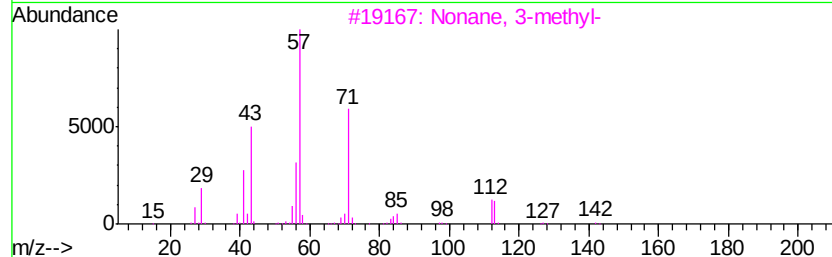
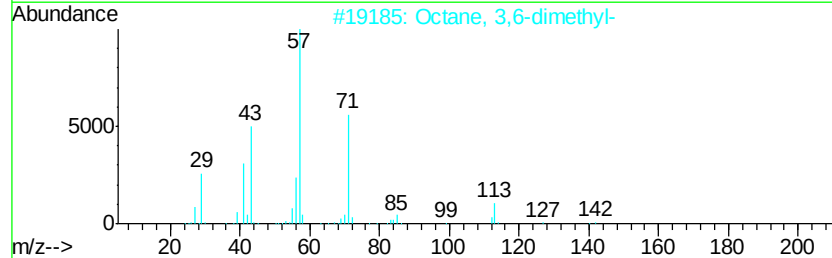
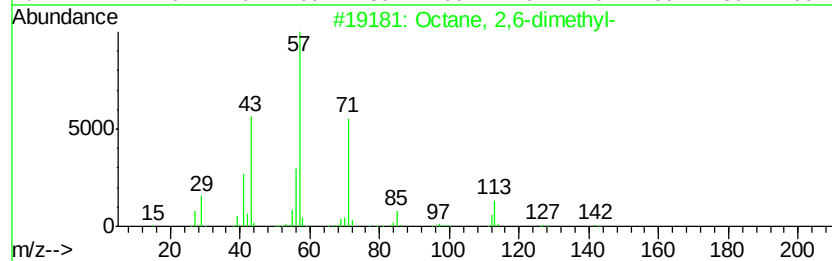
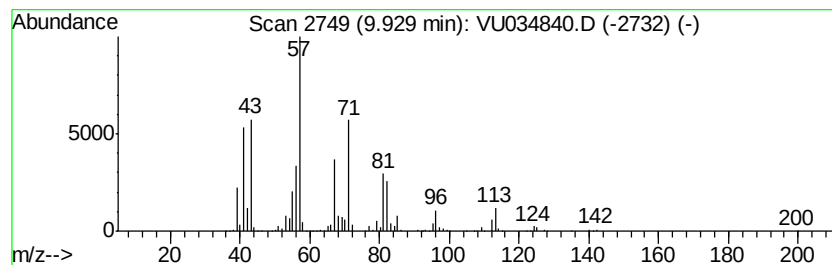
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T. | EstConc    | Area    | Relative to ISTD | R.T. |
|------|------------|---------|------------------|------|
| 9.93 | 62.16 ug/L | 1844890 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 90   |
| 2    |    | Octane, 3,6-dimethyl- | 142 | C10H22  | 015869-94-0 | 60   |
| 3    |    | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 60   |
| 4    |    | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 60   |
| 5    |    | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 53   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

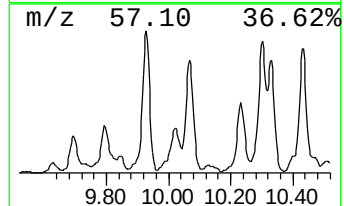
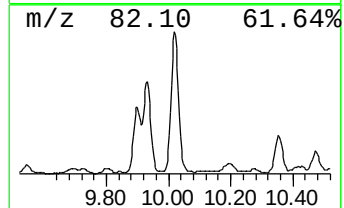
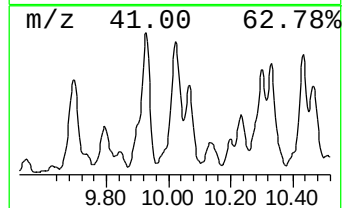
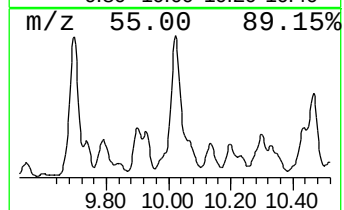
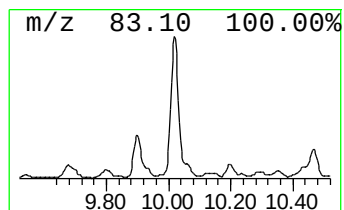
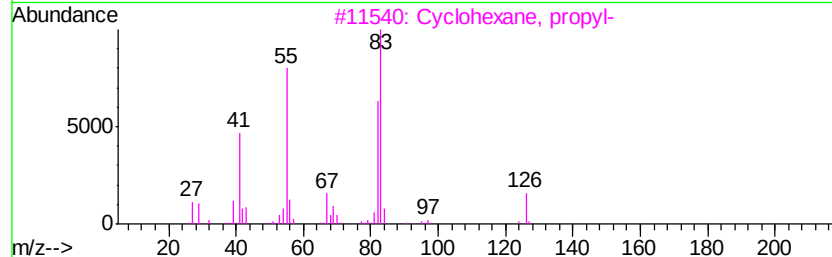
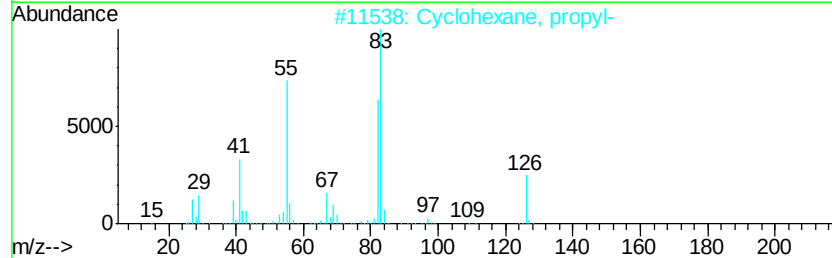
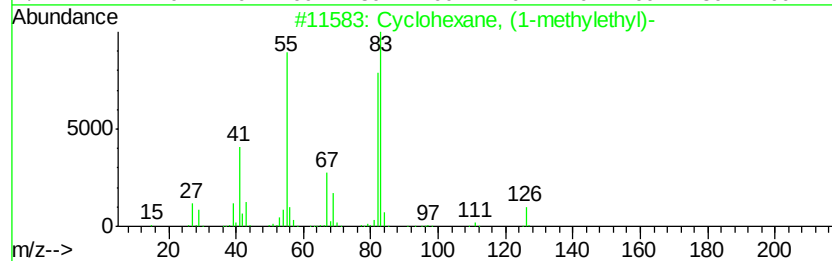
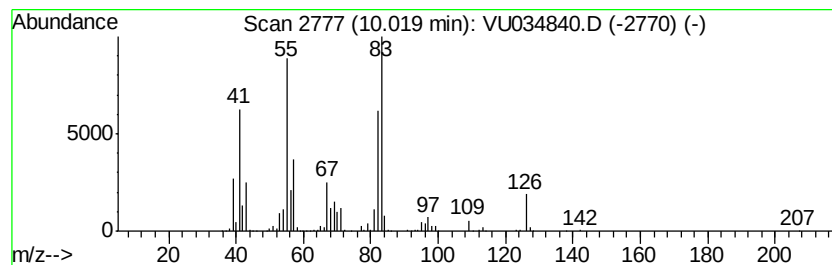
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 (DEL) Alkane: Cyclic10.02 Concentration Rank 24

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T. |
|-------|------------|---------|------------------|------|
| 10.02 | 39.92 ug/L | 1184770 | Chlorobenzene-d5 | 9.07 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 83   |
| 2    |    | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 72   |
| 3    |    | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 72   |
| 4    |    | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 72   |
| 5    |    | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

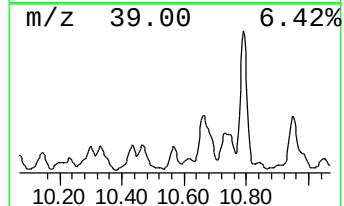
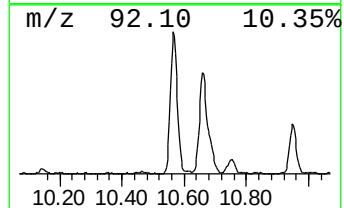
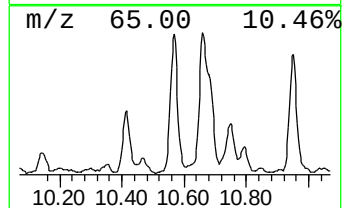
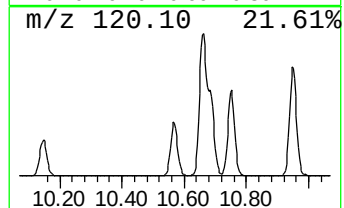
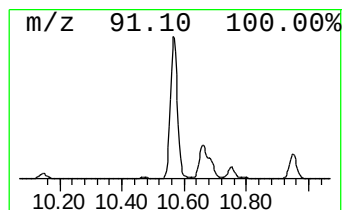
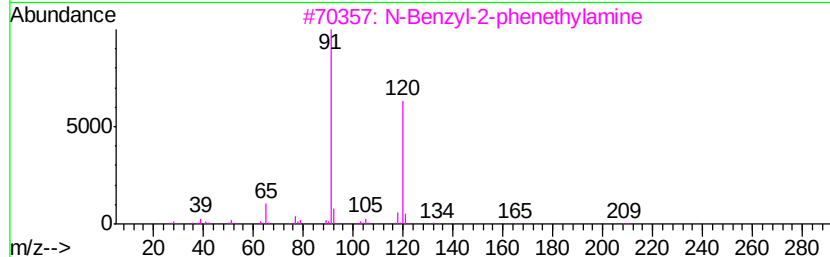
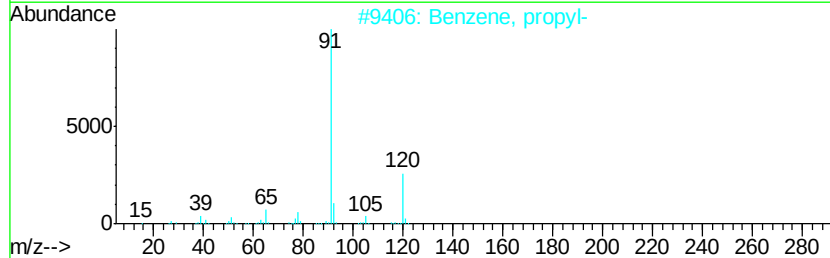
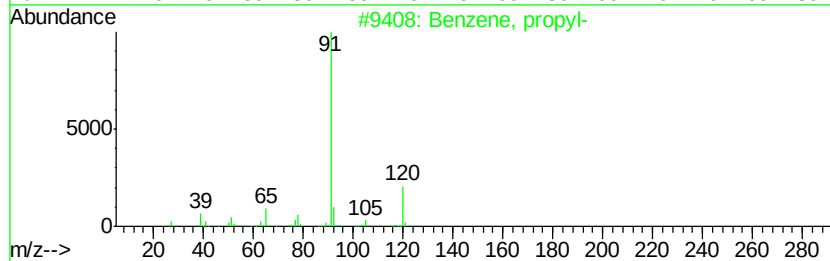
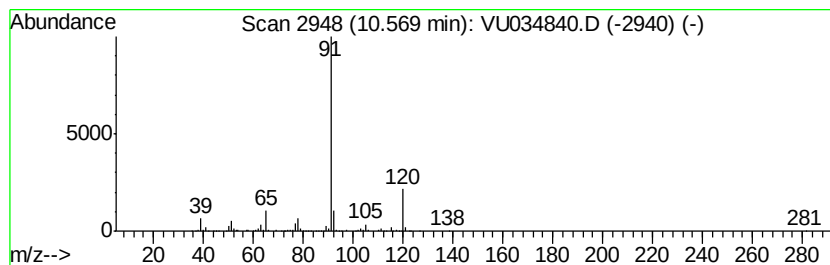
Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

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Peak Number 27 Benzene, propyl- Concentration Rank 40

| R.T.      | EstConc                            | Area   | Relative to ISTD       | R.T.        |      |
|-----------|------------------------------------|--------|------------------------|-------------|------|
| 10.57     | 19.67 ug/L                         | 805039 | 1,4-Dichlorobenzene-d4 | 11.46       |      |
| Hit# of 5 | Tentative ID                       | MW     | MolForm                | CAS#        | Qual |
| 1         | Benzene, propyl-                   | 120    | C9H12                  | 000103-65-1 | 90   |
| 2         | Benzene, propyl-                   | 120    | C9H12                  | 000103-65-1 | 86   |
| 3         | N-Benzyl-2-phenethylamine          | 211    | C15H17N                | 003647-71-0 | 78   |
| 4         | Phenethylamine, N-benzyl-p-chloro- | 245    | C15H16ClN              | 013622-43-0 | 78   |
| 5         | N-Benzyl-2-phenethylamine          | 211    | C15H17N                | 003647-71-0 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

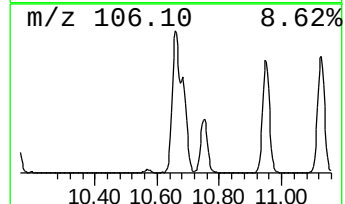
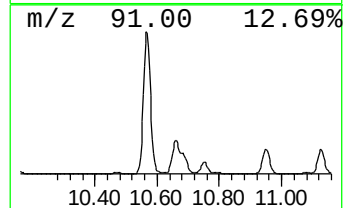
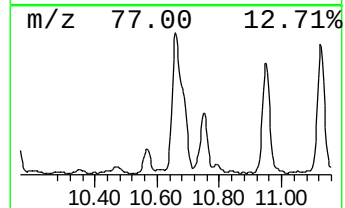
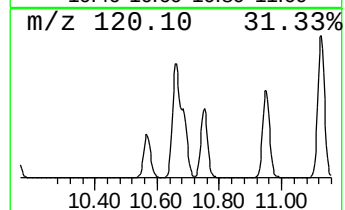
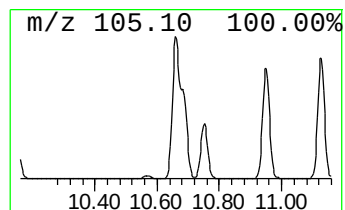
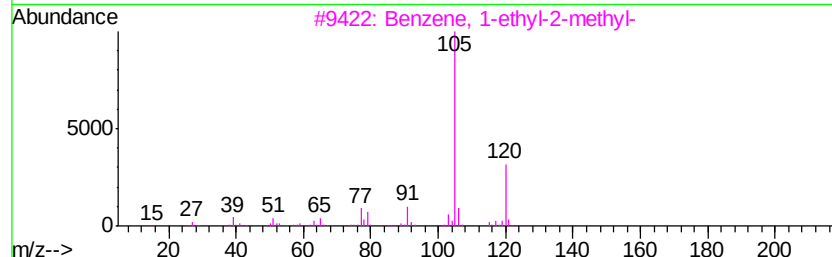
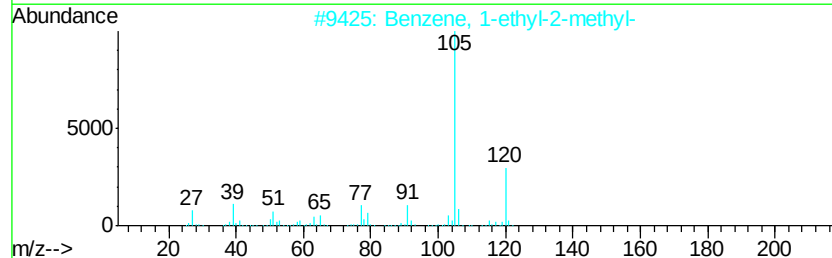
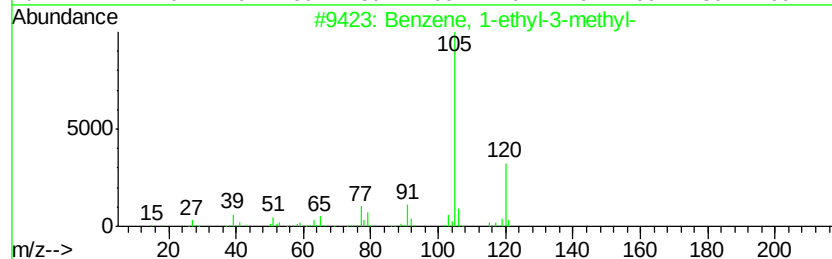
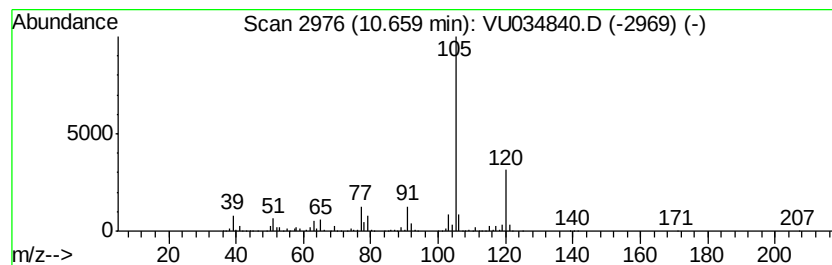
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 Benzene, 1-ethyl-3-methyl- Concentration Rank 11

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 10.66 | 74.95 ug/L | 3067180 | 1,4-Dichlorobenzene-d4 | 11.46 |

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

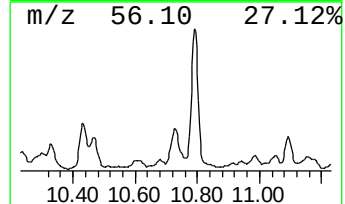
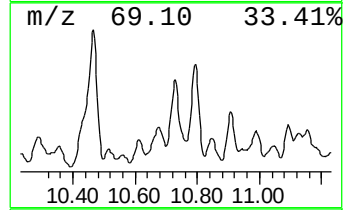
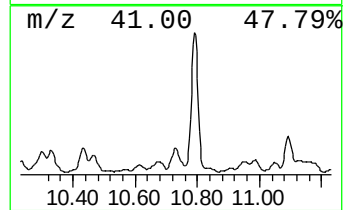
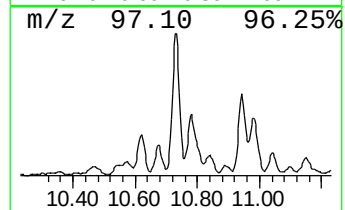
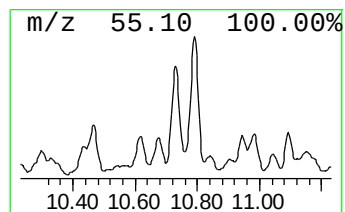
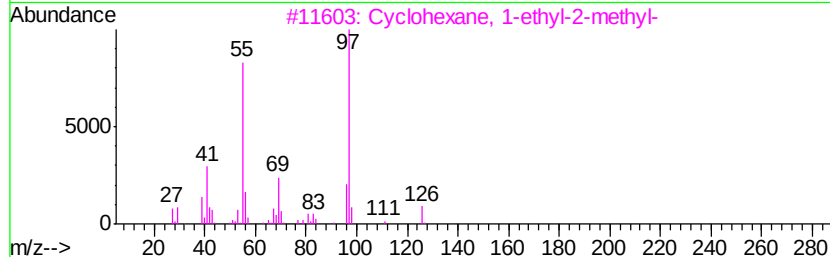
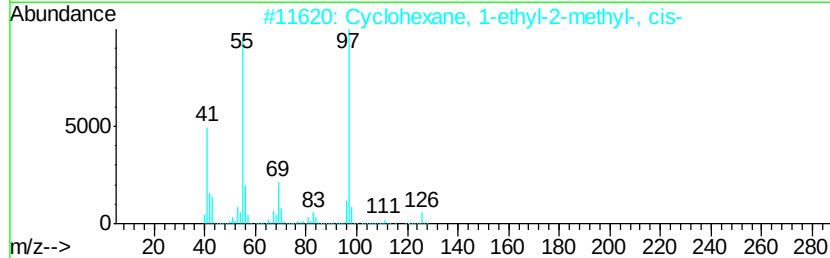
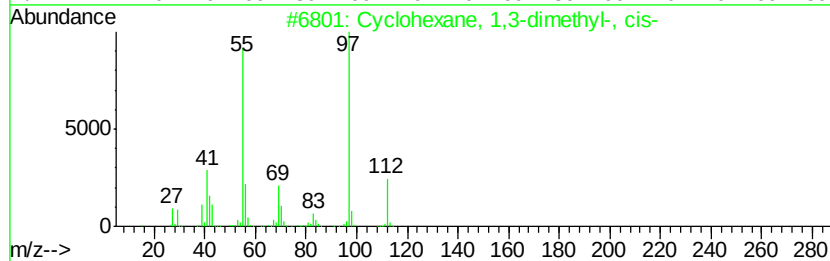
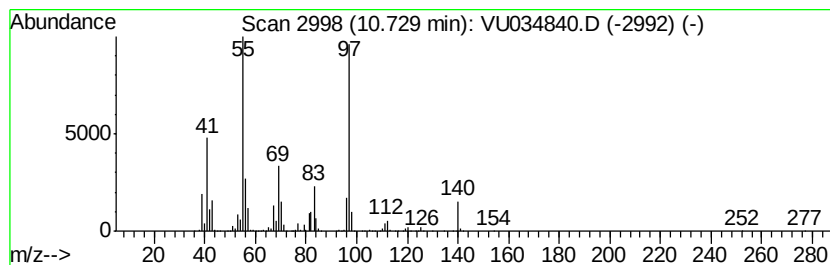
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 (DEL) Alkane: Cyclic10.73 Concentration Rank 20

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 10.73 | 43.70 ug/L | 1788310 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1,3-dimethyl-, cis-    | 112 | C8H16   | 000638-04-0 | 64   |
| 2    |    | Cyclohexane, 1-ethyl-2-methyl-, ... | 126 | C9H18   | 004923-77-7 | 64   |
| 3    |    | Cyclohexane, 1-ethyl-2-methyl-      | 126 | C9H18   | 003728-54-9 | 53   |
| 4    |    | cis-1-Ethyl-3-methyl-cyclohexane    | 126 | C9H18   | 019489-10-2 | 53   |
| 5    |    | Cyclohexane, 1-methyl-3-propyl-     | 140 | C10H20  | 004291-80-9 | 52   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
 Data File : VU034840.D  
 Acq On : 26 Sep 2019 12:48  
 Operator : JC/SP  
 Sample : K4984-08  
 Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
 Quant Title : VOC Analysis

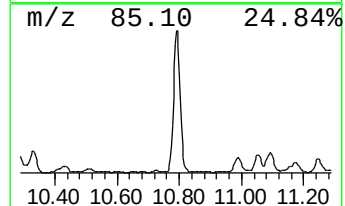
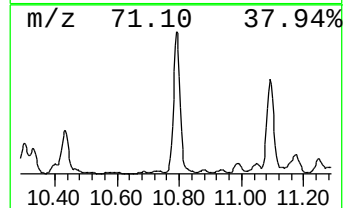
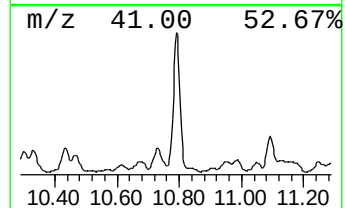
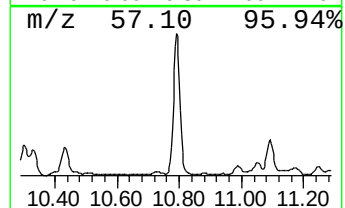
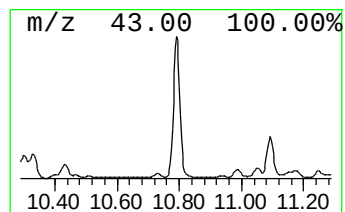
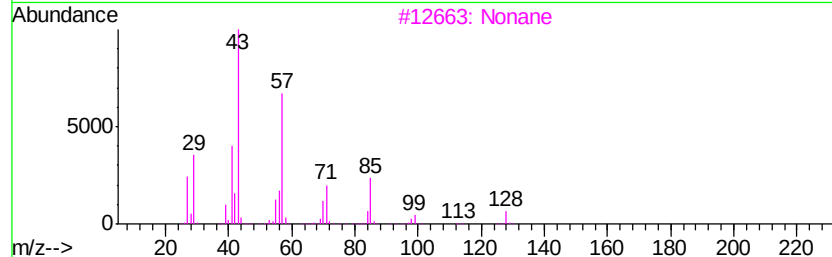
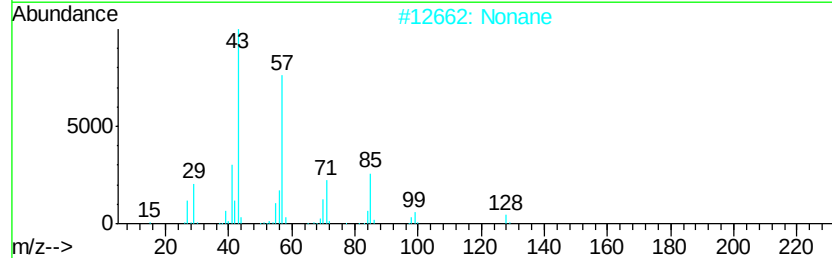
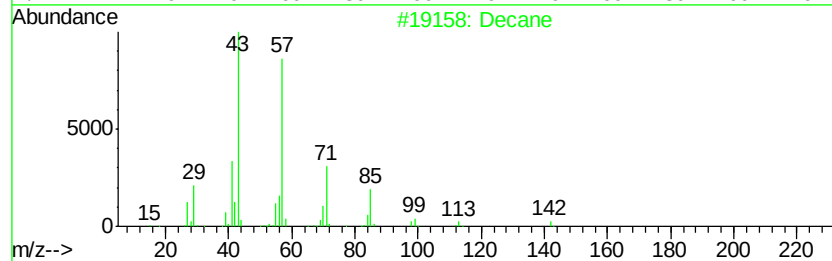
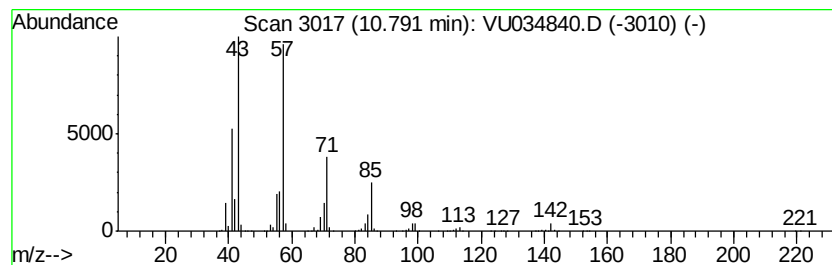
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 10.79 | 127.64 ug/L | 5223540 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Decane       | 142 | C10H22  | 000124-18-5 | 94   |
| 2    |    | Nonane       | 128 | C9H20   | 000111-84-2 | 91   |
| 3    |    | Nonane       | 128 | C9H20   | 000111-84-2 | 91   |
| 4    |    | Tetradecane  | 198 | C14H30  | 000629-59-4 | 80   |
| 5    |    | Nonane       | 128 | C9H20   | 000111-84-2 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

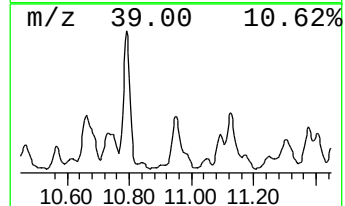
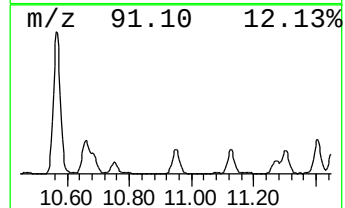
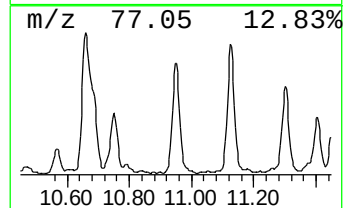
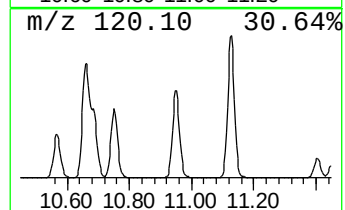
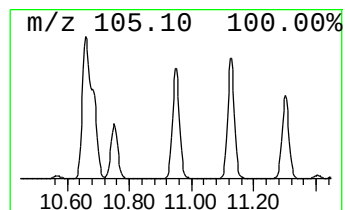
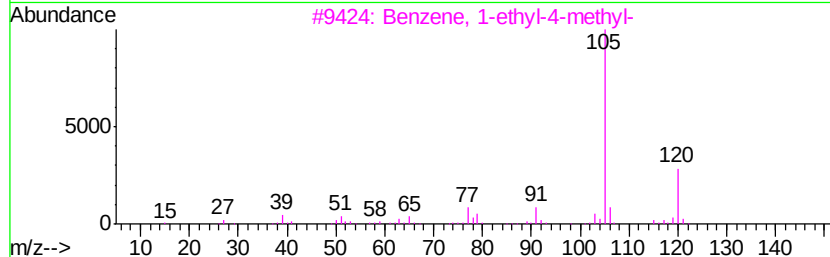
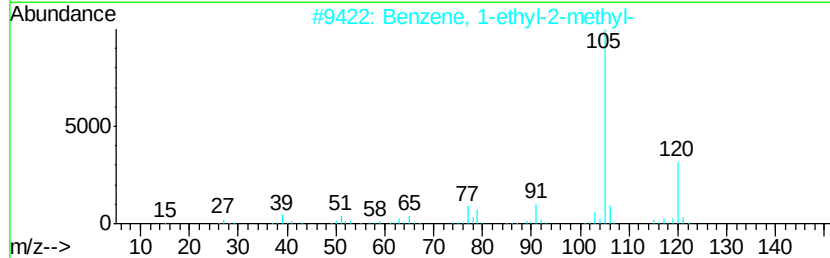
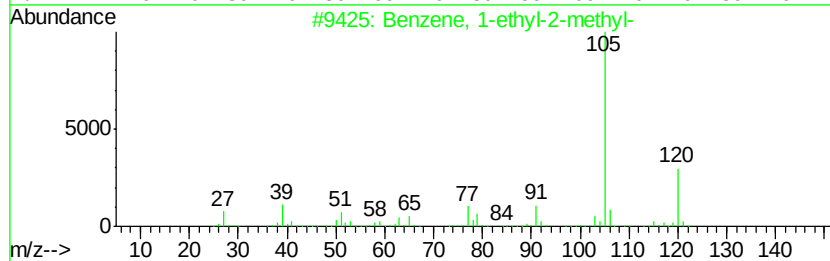
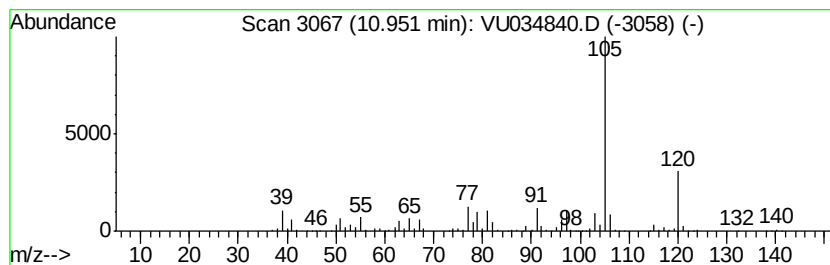
Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

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Peak Number 31 Benzene, 1-ethyl-2-methyl- Concentration Rank 17

| R.T.    | EstConc    | Area                       | Relative to ISTD       | R.T.           |
|---------|------------|----------------------------|------------------------|----------------|
| 10.95   | 52.07 ug/L | 2131000                    | 1,4-Dichlorobenzene-d4 | 11.46          |
| Hit# of | 5          | Tentative ID               | MW MolForm             | CAS# Qual      |
| 1       |            | Benzene, 1-ethyl-2-methyl- | 120 C9H12              | 000611-14-3 93 |
| 2       |            | Benzene, 1-ethyl-2-methyl- | 120 C9H12              | 000611-14-3 76 |
| 3       |            | Benzene, 1-ethyl-4-methyl- | 120 C9H12              | 000622-96-8 76 |
| 4       |            | Benzene, 1-ethyl-4-methyl- | 120 C9H12              | 000622-96-8 76 |
| 5       |            | Mesitylene                 | 120 C9H12              | 000108-67-8 76 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

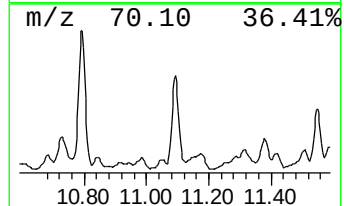
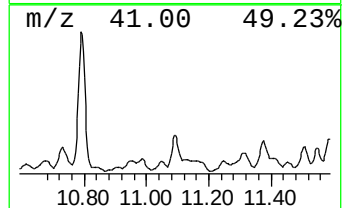
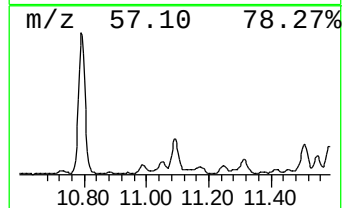
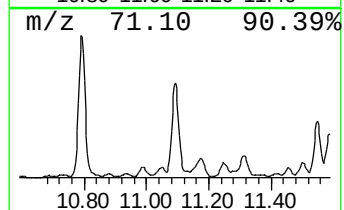
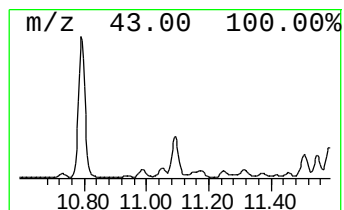
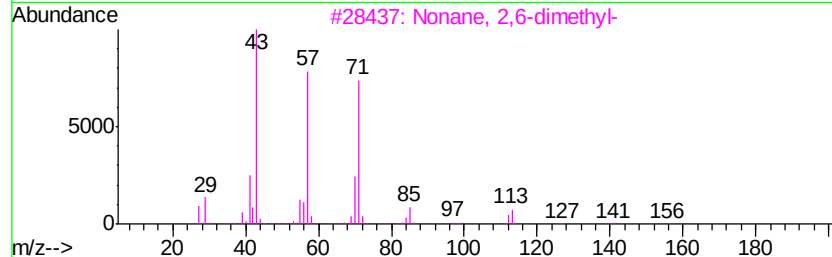
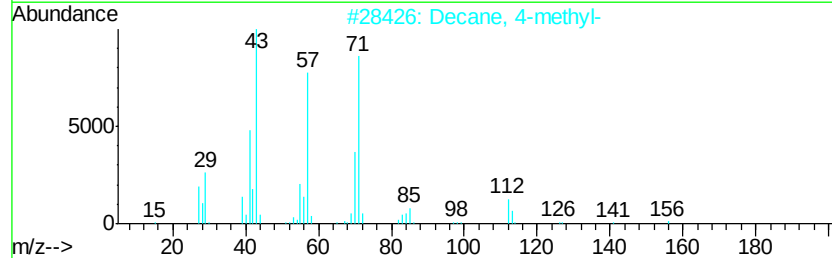
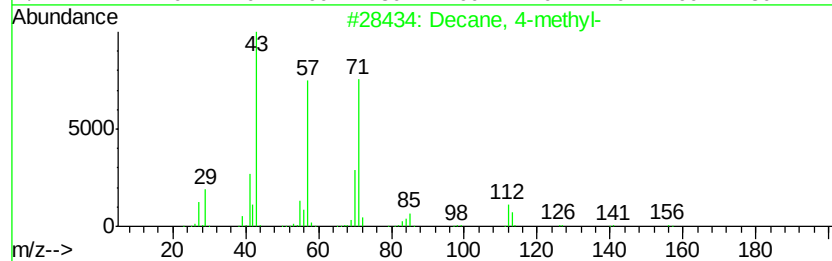
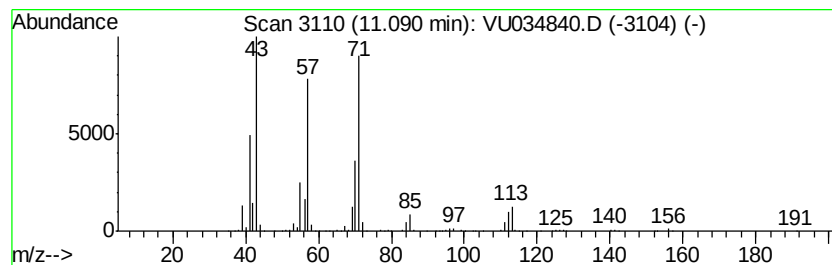
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 11.09 | 35.84 ug/L | 1466830 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 4-methyl-     | 156 | C11H24  | 002847-72-5 | 94   |
| 2    |    | Decane, 4-methyl-     | 156 | C11H24  | 002847-72-5 | 90   |
| 3    |    | Nonane, 2,6-dimethyl- | 156 | C11H24  | 017302-28-2 | 87   |
| 4    |    | Decane, 4-methyl-     | 156 | C11H24  | 002847-72-5 | 78   |
| 5    |    | Octane, 3,3-dimethyl- | 142 | C10H22  | 004110-44-5 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

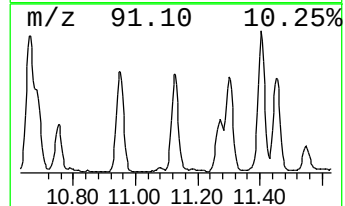
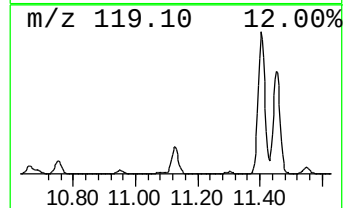
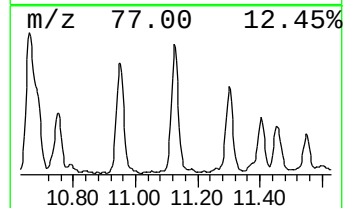
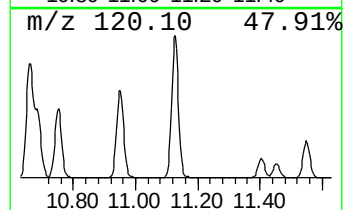
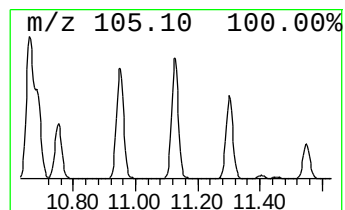
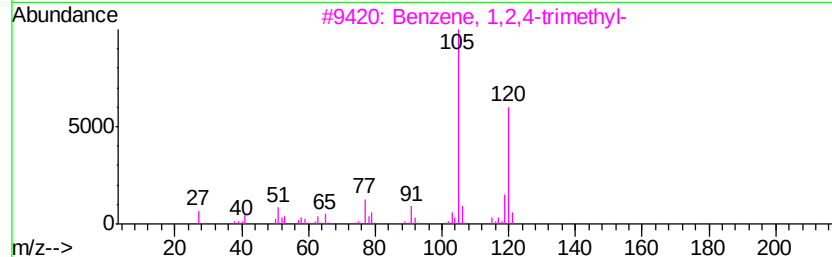
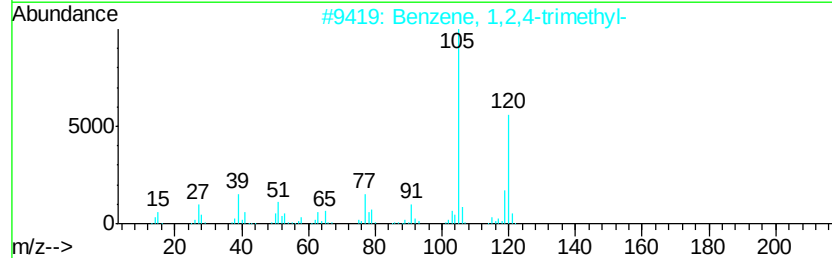
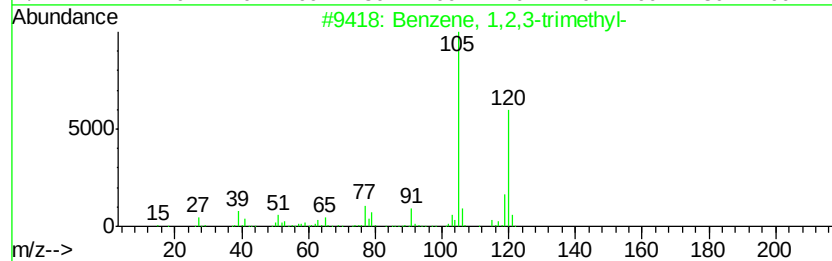
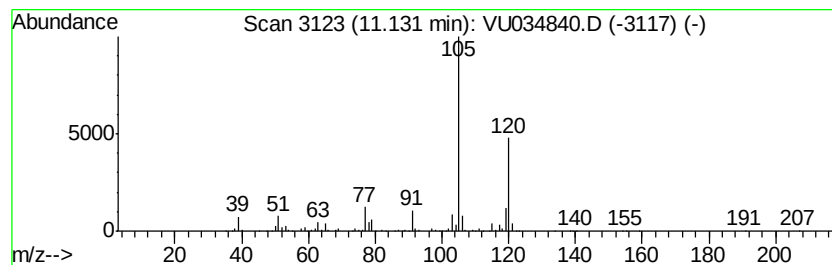
Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 33 Benzene, 1,2,3-trimethyl- Concentration Rank 21

| R.T.    | EstConc    | Area                      | Relative to ISTD       | R.T.           |
|---------|------------|---------------------------|------------------------|----------------|
| 11.13   | 42.92 ug/L | 1756550                   | 1,4-Dichlorobenzene-d4 | 11.46          |
| Hit# of | 5          | Tentative ID              | MW MolForm             | CAS# Qual      |
| 1       |            | Benzene, 1,2,3-trimethyl- | 120 C9H12              | 000526-73-8 97 |
| 2       |            | Benzene, 1,2,4-trimethyl- | 120 C9H12              | 000095-63-6 95 |
| 3       |            | Benzene, 1,2,4-trimethyl- | 120 C9H12              | 000095-63-6 95 |
| 4       |            | Mesitylene                | 120 C9H12              | 000108-67-8 94 |
| 5       |            | Benzene, 1,2,3-trimethyl- | 120 C9H12              | 000526-73-8 94 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

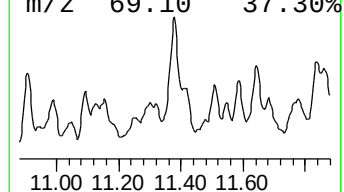
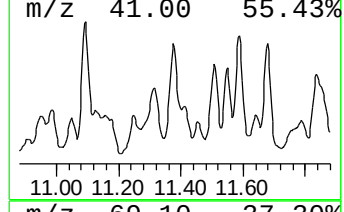
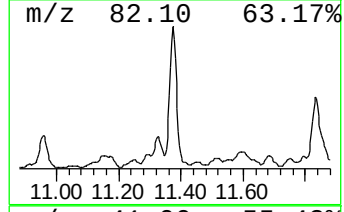
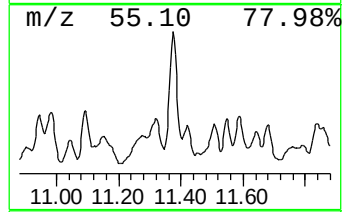
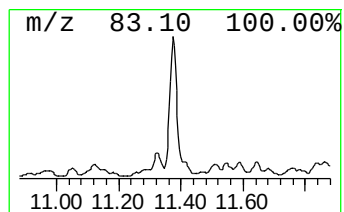
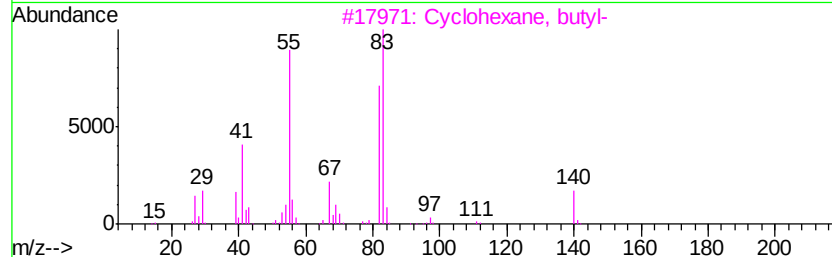
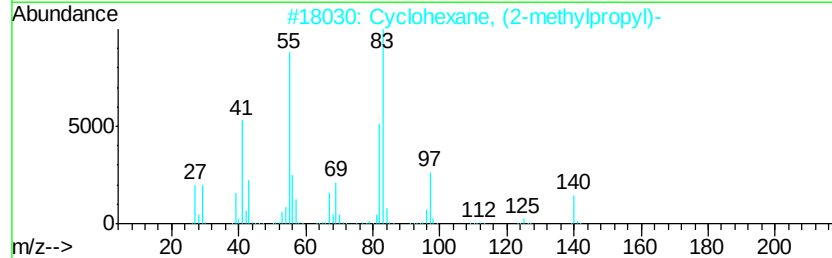
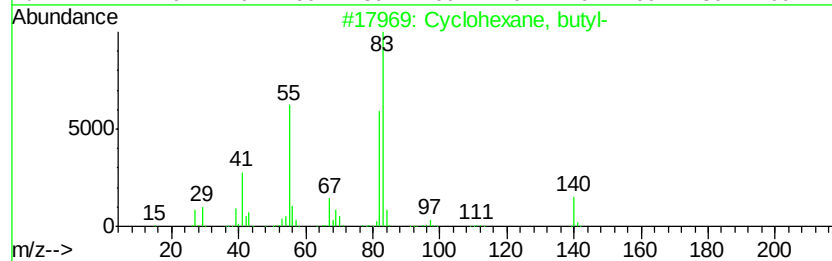
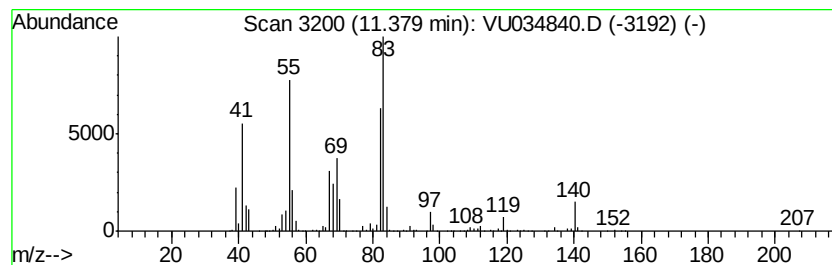
Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 34 (DEL) Alkane: Cyclic11.38 Concentration Rank 34

| R.T.    | EstConc    | Area                           | Relative to ISTD       | R.T.           |
|---------|------------|--------------------------------|------------------------|----------------|
| 11.38   | 29.03 ug/L | 1187980                        | 1,4-Dichlorobenzene-d4 | 11.46          |
| Hit# of | 5          | Tentative ID                   | MW MolForm             | CAS# Qual      |
| 1       |            | Cyclohexane, butyl-            | 140 C10H20             | 001678-93-9 72 |
| 2       |            | Cyclohexane, (2-methylpropyl)- | 140 C10H20             | 001678-98-4 72 |
| 3       |            | Cyclohexane, butyl-            | 140 C10H20             | 001678-93-9 72 |
| 4       |            | Cyclohexane, (2-methylpropyl)- | 140 C10H20             | 001678-98-4 72 |
| 5       |            | Cyclohexane, butyl-            | 140 C10H20             | 001678-93-9 64 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

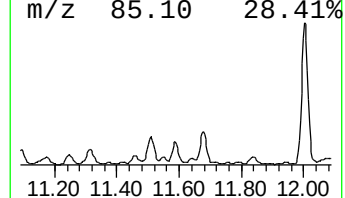
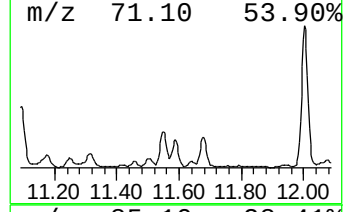
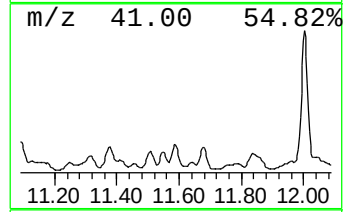
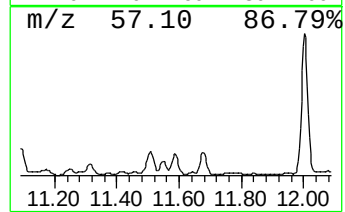
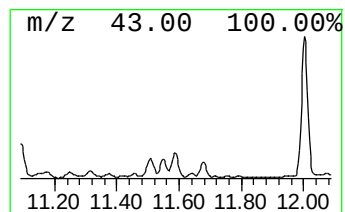
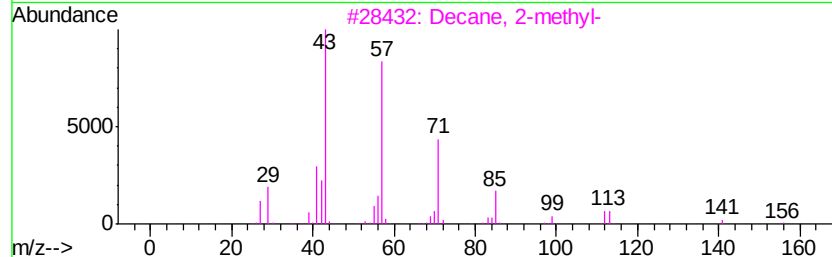
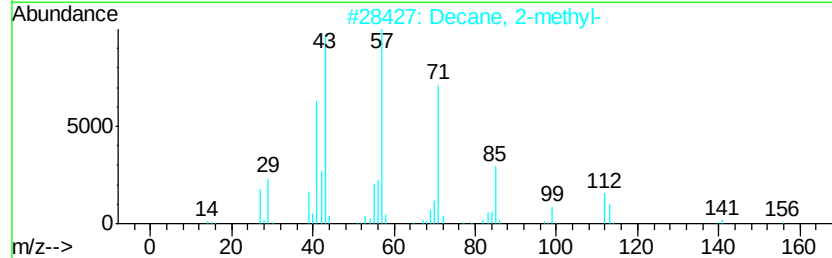
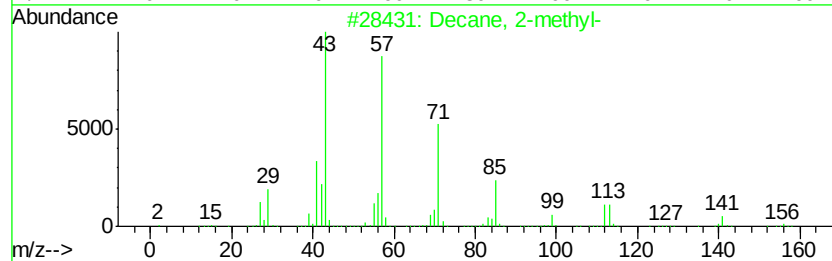
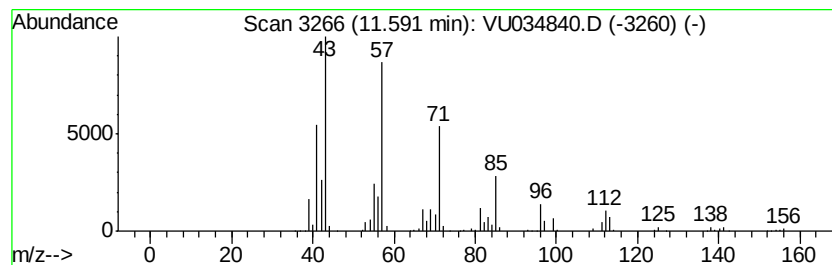
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 37

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 11.59 | 25.88 ug/L | 1058970 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2-methyl-     | 156 | C11H24  | 006975-98-0 | 87   |
| 2    |    |   | Decane, 2-methyl-     | 156 | C11H24  | 006975-98-0 | 76   |
| 3    |    |   | Decane, 2-methyl-     | 156 | C11H24  | 006975-98-0 | 72   |
| 4    |    |   | Decane, 2,9-dimethyl- | 170 | C12H26  | 001002-17-1 | 59   |
| 5    |    |   | Undecane, 5-methyl-   | 170 | C12H26  | 001632-70-8 | 59   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

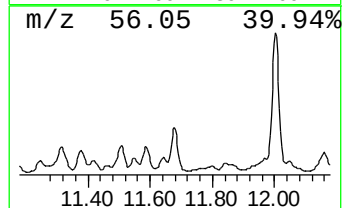
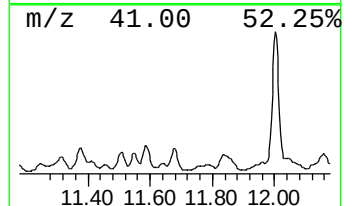
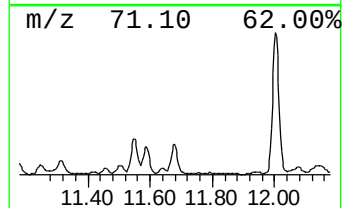
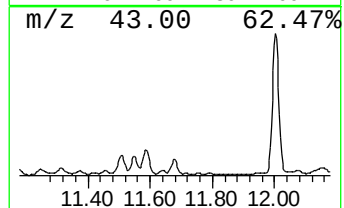
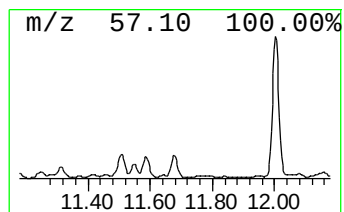
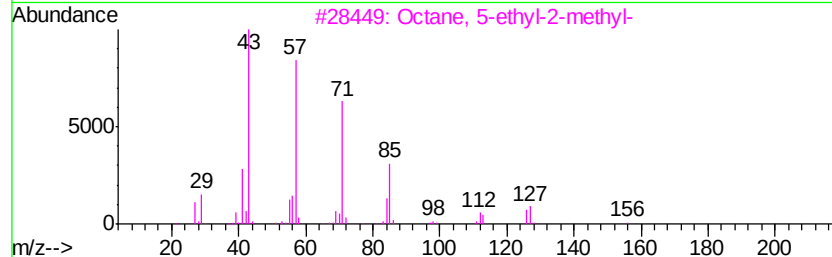
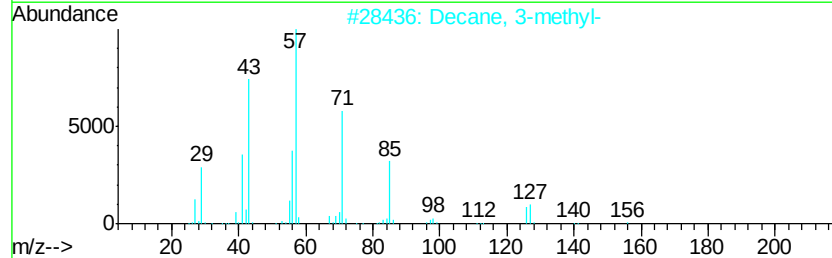
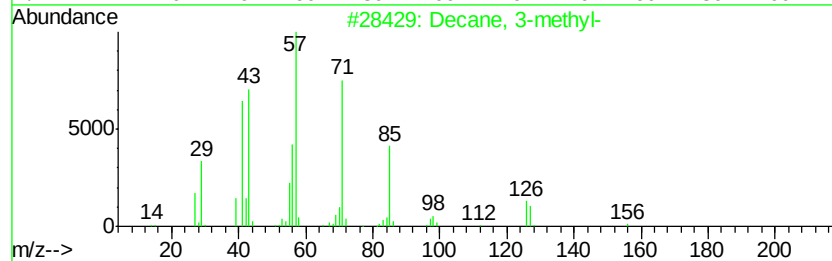
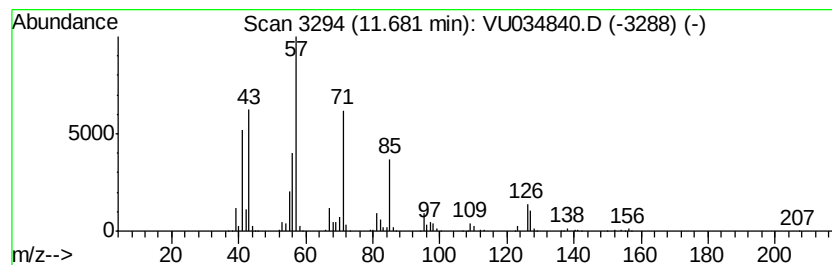
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 31

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 11.68 | 30.26 ug/L | 1238260 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                      | MW  | MolForm   | CAS#         | Qual |
|------|----|-----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Decane, 3-methyl-                 | 156 | C11H24    | 013151-34-3  | 94   |
| 2    |    | Decane, 3-methyl-                 | 156 | C11H24    | 013151-34-3  | 78   |
| 3    |    | Octane, 5-ethyl-2-methyl-         | 156 | C11H24    | 062016-18-6  | 53   |
| 4    |    | Sulfurous acid, butyl nonyl ester | 264 | C13H28O3S | 1000309-17-6 | 53   |
| 5    |    | Heptadecane, 8-methyl-            | 254 | C18H38    | 013287-23-5  | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

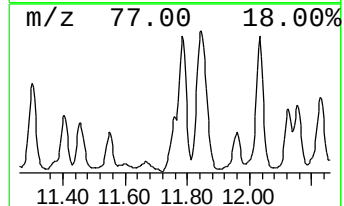
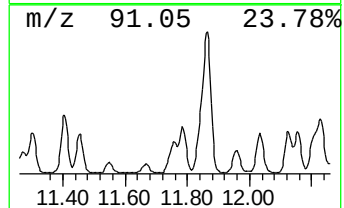
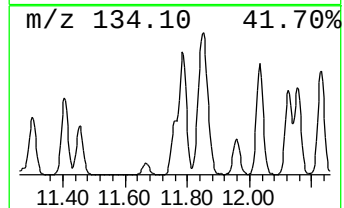
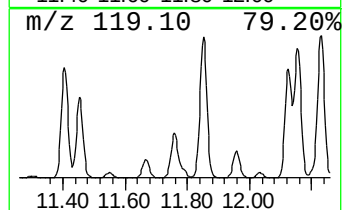
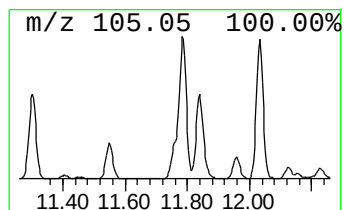
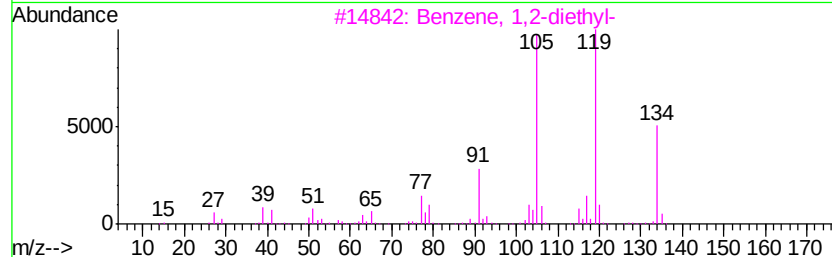
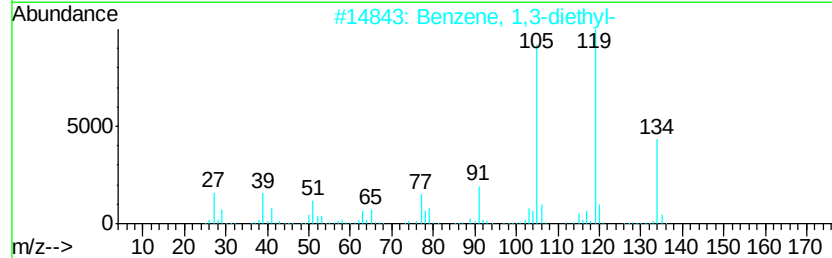
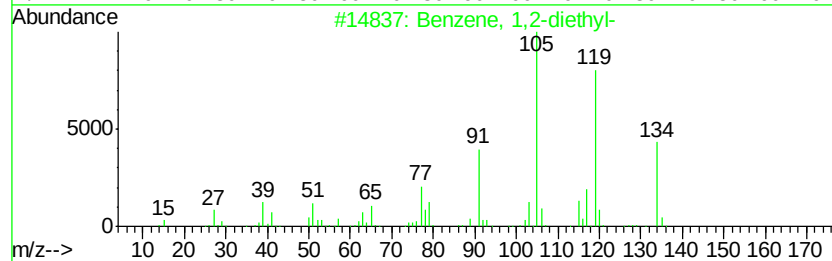
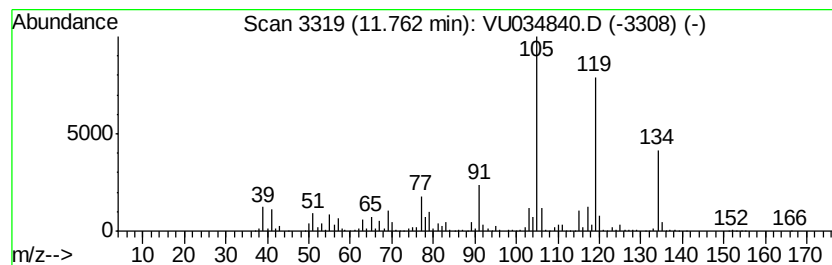
Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 38 Benzene, 1,2-diethyl- Concentration Rank 32

| R.T.    | EstConc    | Area                  | Relative to ISTD       | R.T.           |
|---------|------------|-----------------------|------------------------|----------------|
| 11.76   | 30.10 ug/L | 1231590               | 1,4-Dichlorobenzene-d4 | 11.46          |
| Hit# of | 5          | Tentative ID          | MW MolForm             | CAS# Qual      |
| 1       |            | Benzene, 1,2-diethyl- | 134 C10H14             | 000135-01-3 91 |
| 2       |            | Benzene, 1,3-diethyl- | 134 C10H14             | 000141-93-5 76 |
| 3       |            | Benzene, 1,2-diethyl- | 134 C10H14             | 000135-01-3 76 |
| 4       |            | Benzene, 1,4-diethyl- | 134 C10H14             | 000105-05-5 76 |
| 5       |            | Benzene, 1,4-diethyl- | 134 C10H14             | 000105-05-5 76 |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

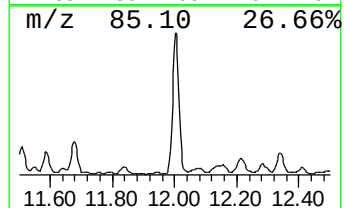
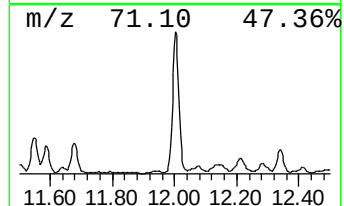
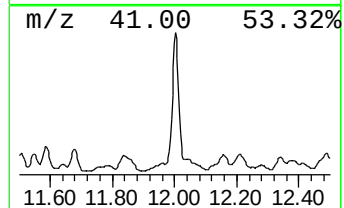
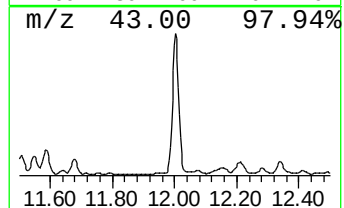
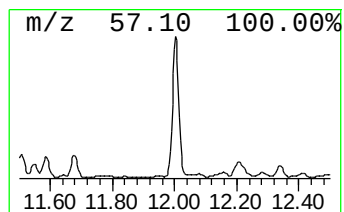
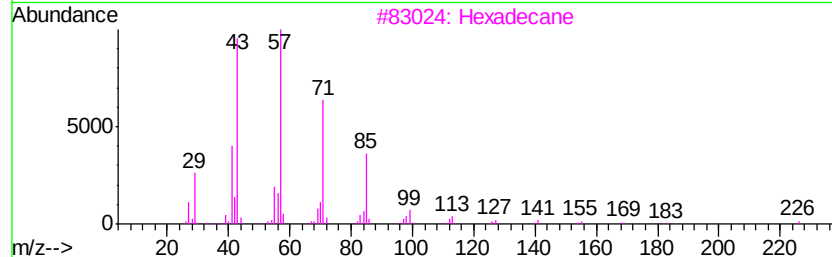
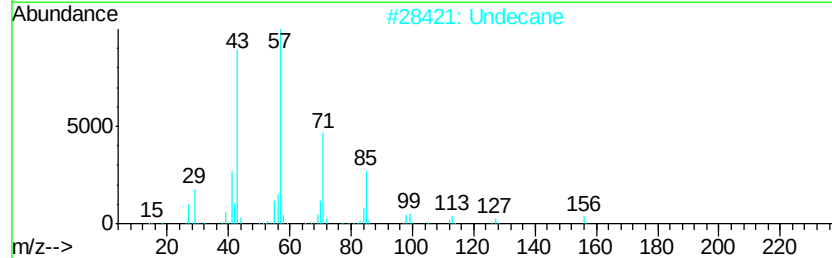
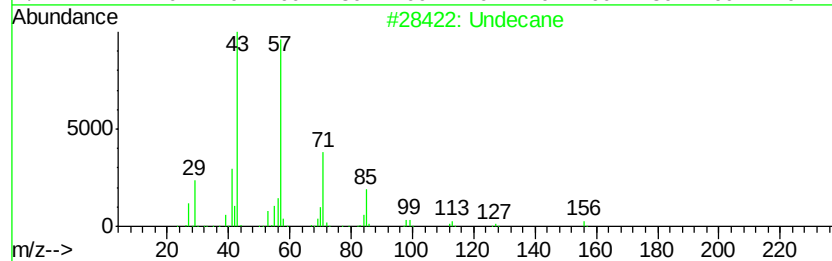
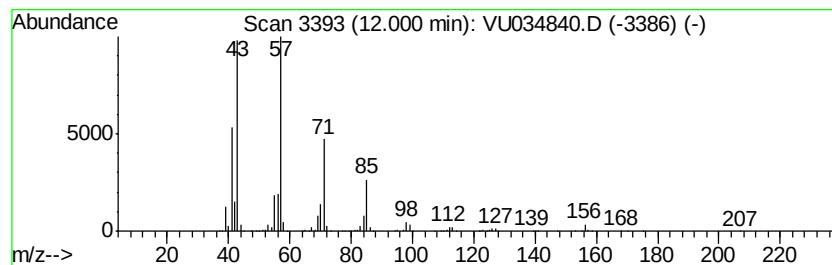
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TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 12.00 | 149.99 ug/L | 6138190 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Undecane                 | 156 | C11H24  | 001120-21-4 | 91   |
| 2    |    | Undecane                 | 156 | C11H24  | 001120-21-4 | 91   |
| 3    |    | Hexadecane               | 226 | C16H34  | 000544-76-3 | 83   |
| 4    |    | Undecane, 2,3-dimethyl-  | 184 | C13H28  | 017312-77-5 | 83   |
| 5    |    | Decane, 2,3,6-trimethyl- | 184 | C13H28  | 062238-12-4 | 78   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22a/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

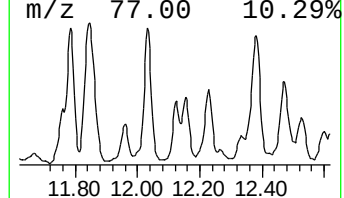
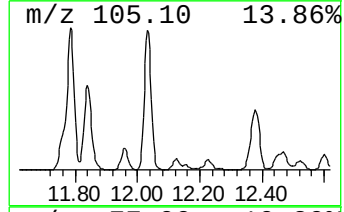
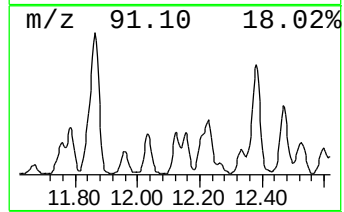
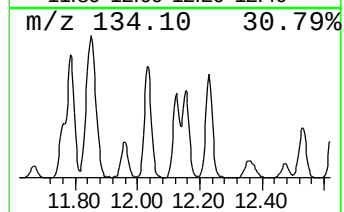
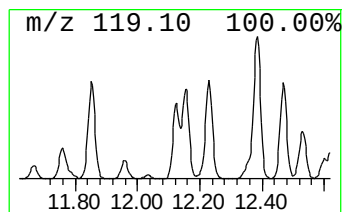
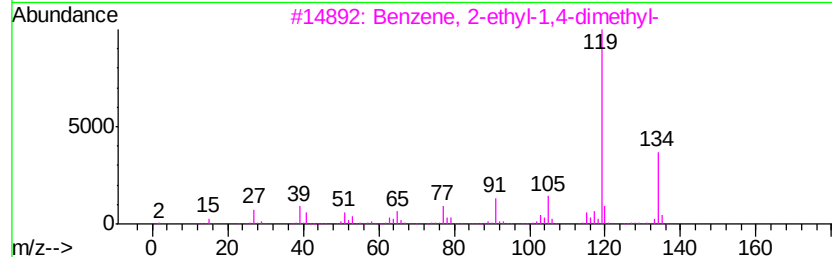
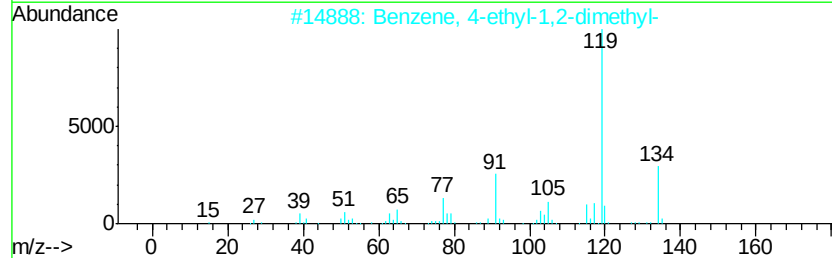
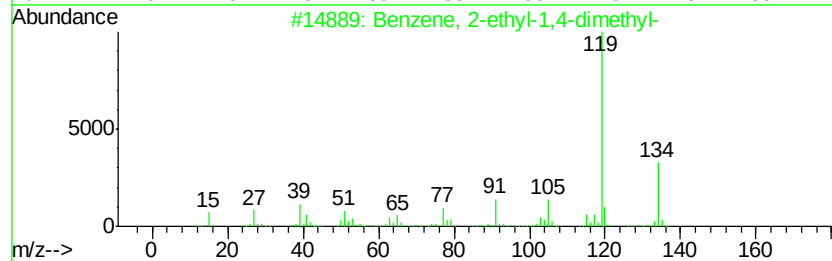
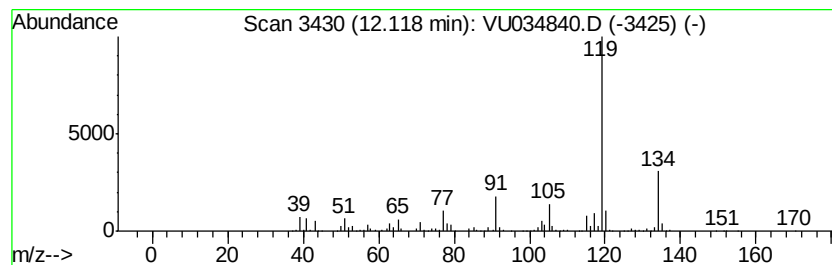
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 40 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 35

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.12 | 27.84 ug/L | 1139220 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 97   |
| 2    |    | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 96   |
| 3    |    | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 95   |
| 4    |    | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 95   |
| 5    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

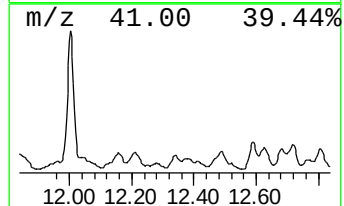
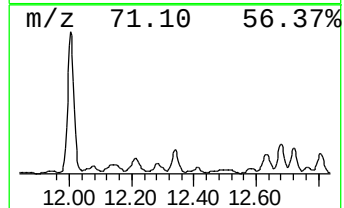
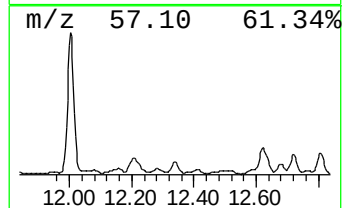
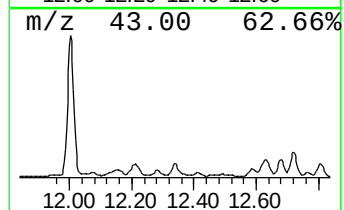
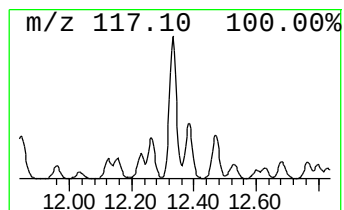
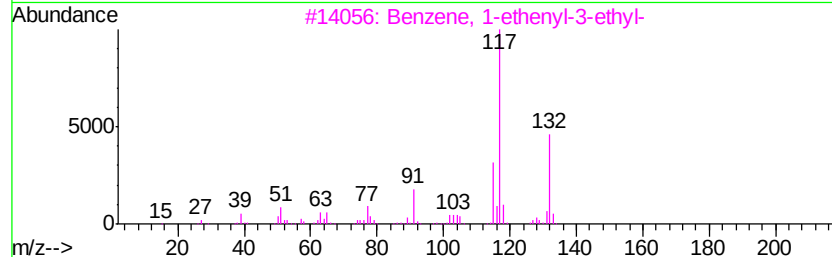
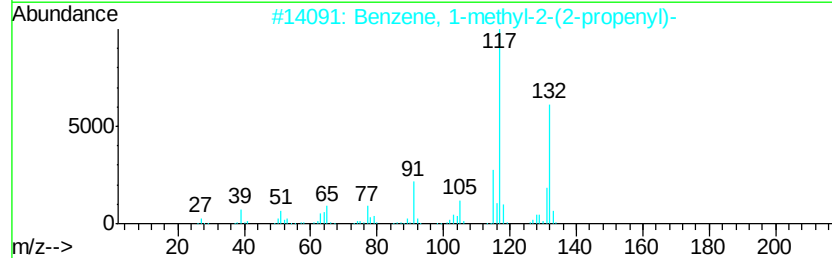
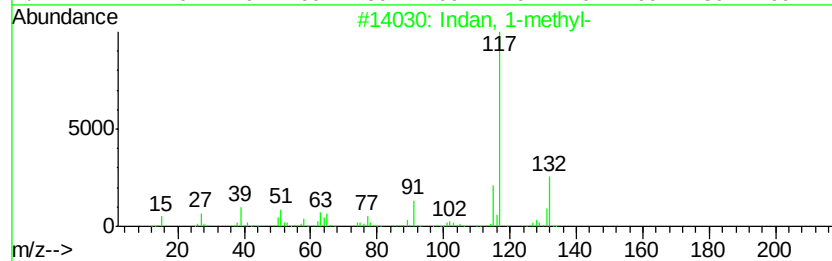
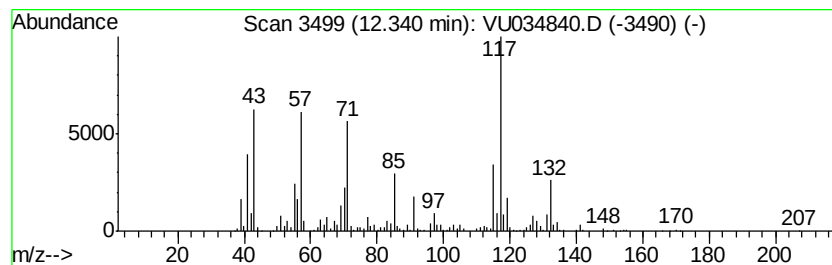
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 41 Indan, 1-methyl- Concentration Rank 25

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.34 | 37.42 ug/L | 1531410 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Indan, 1-methyl-                    | 132 | C10H12  | 000767-58-8 | 87   |
| 2    |    | Benzene, 1-methyl-2-(2-propenyl)-   | 132 | C10H12  | 001587-04-8 | 60   |
| 3    |    | Benzene, 1-ethenyl-3-ethyl-         | 132 | C10H12  | 007525-62-4 | 60   |
| 4    |    | Indan, 1-methyl-                    | 132 | C10H12  | 000767-58-8 | 55   |
| 5    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 49   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

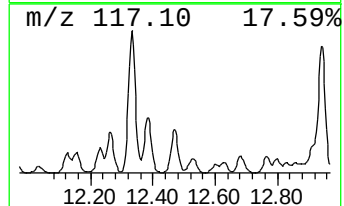
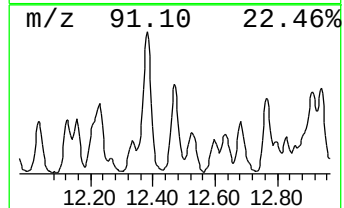
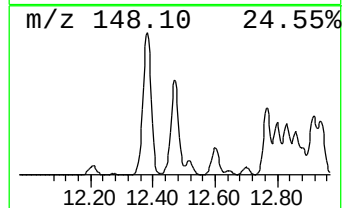
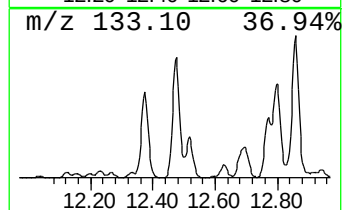
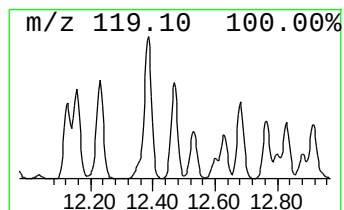
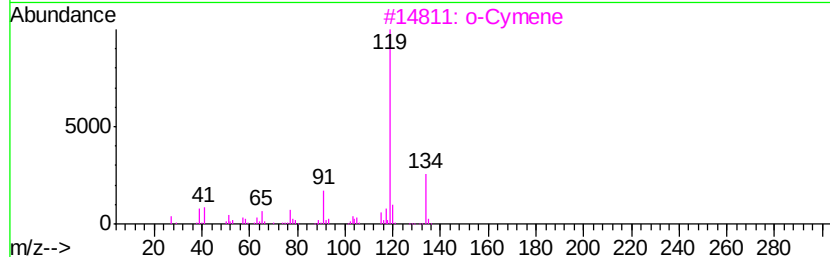
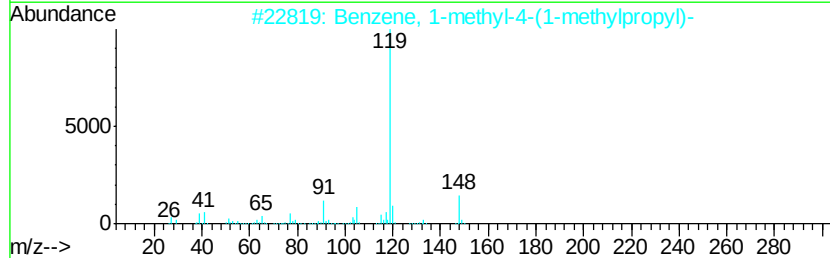
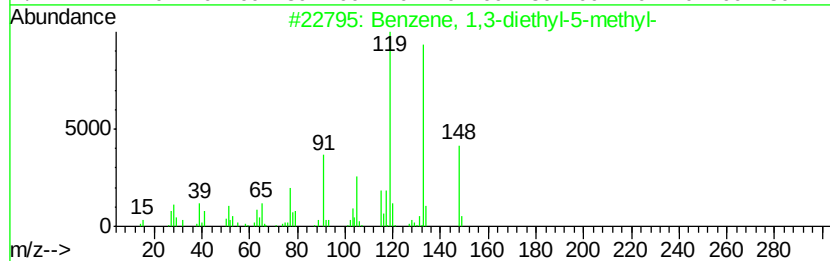
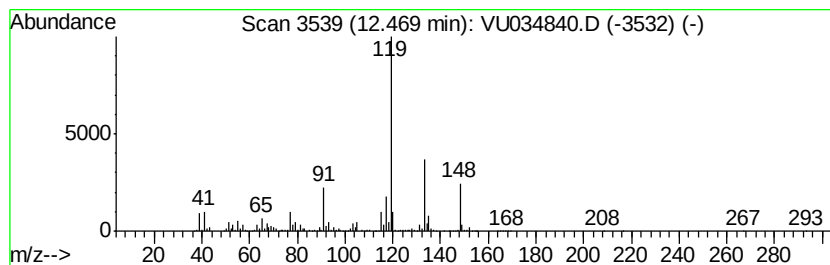
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 42 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 13

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.47 | 67.90 ug/L | 2778490 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,3-diethyl-5-methyl-      | 148 | C11H16  | 002050-24-0 | 70   |
| 2    |    | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16  | 001595-16-0 | 58   |
| 3    |    | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 55   |
| 4    |    | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8 | 52   |
| 5    |    | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8 | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

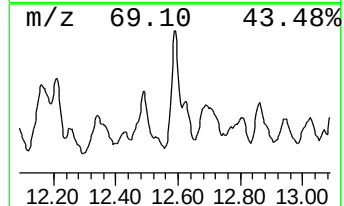
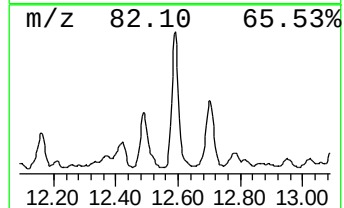
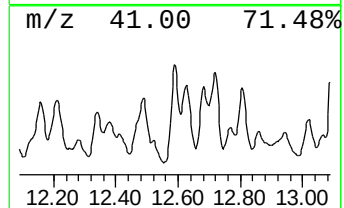
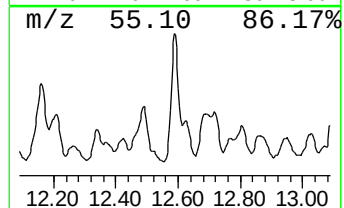
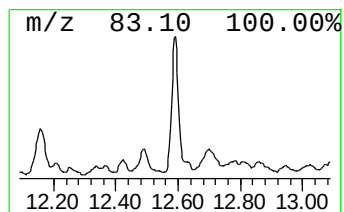
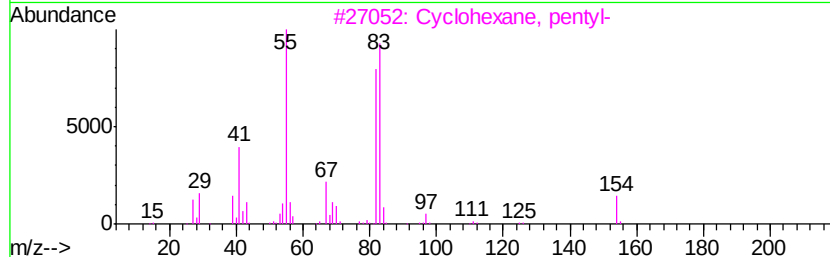
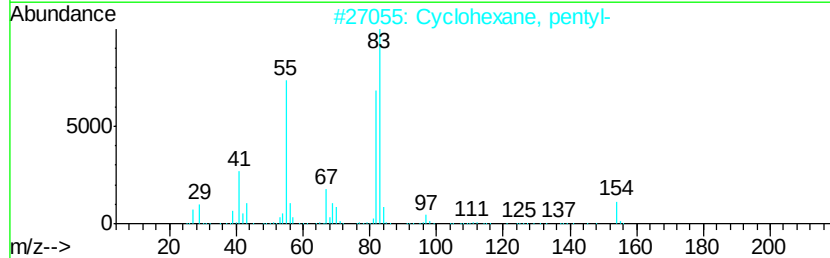
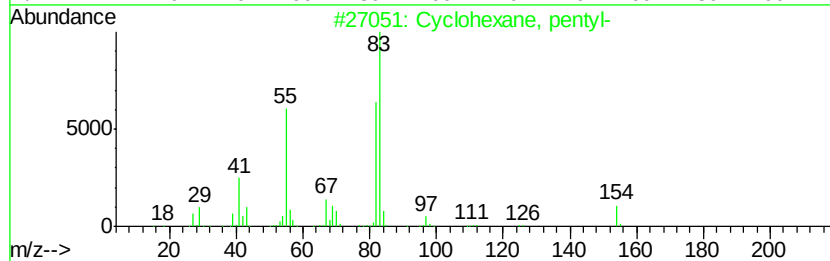
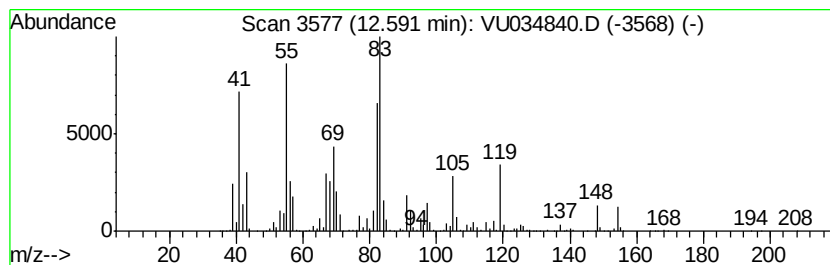
Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 43 (DEL) Alkane: Cyclic12.59 Concentration Rank 18

| R.T.    | EstConc    | Area                    | Relative to ISTD       | R.T.           |
|---------|------------|-------------------------|------------------------|----------------|
| 12.59   | 48.73 ug/L | 1994060                 | 1,4-Dichlorobenzene-d4 | 11.46          |
| Hit# of | 5          | Tentative ID            | MW MolForm             | CAS# Qual      |
| 1       |            | Cyclohexane, pentyl-    | 154 C11H22             | 004292-92-6 52 |
| 2       |            | Cyclohexane, pentyl-    | 154 C11H22             | 004292-92-6 52 |
| 3       |            | Cyclohexane, pentyl-    | 154 C11H22             | 004292-92-6 49 |
| 4       |            | Cyclohexane, nonadecyl- | 350 C25H50             | 022349-03-7 47 |
| 5       |            | Cyclohexane, propyl-    | 126 C9H18              | 001678-92-8 47 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

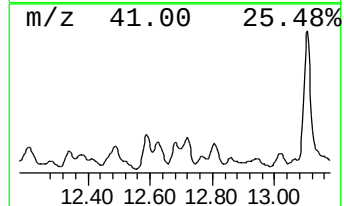
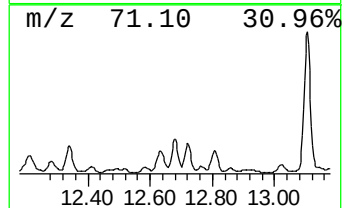
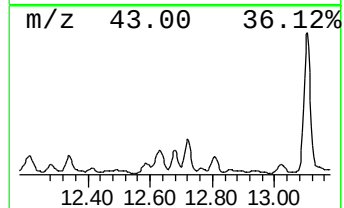
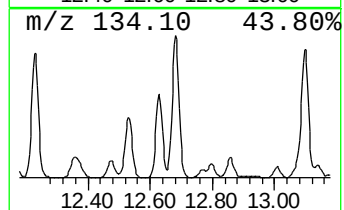
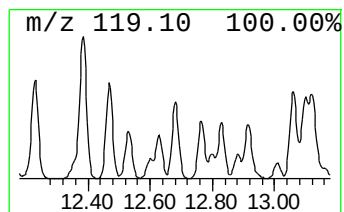
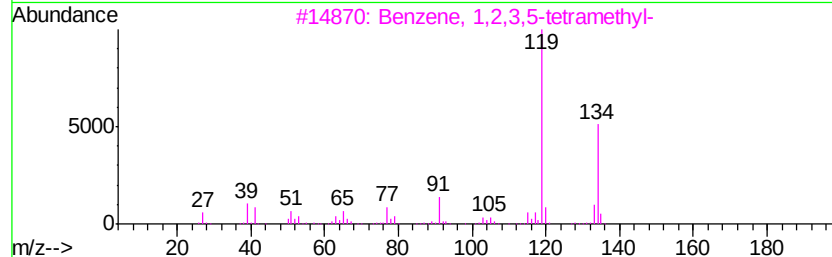
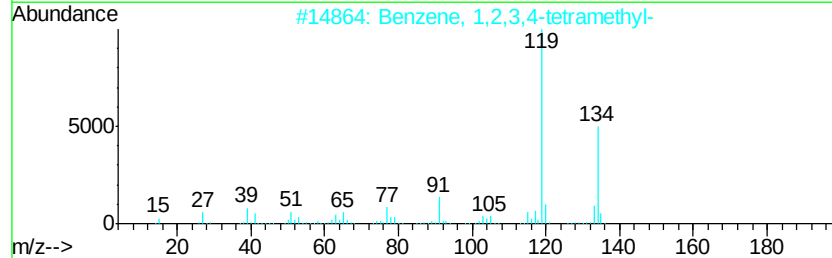
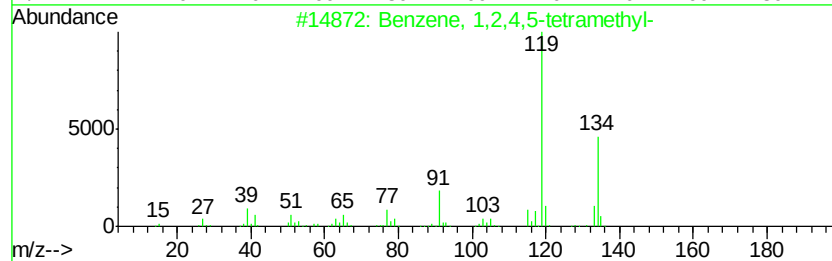
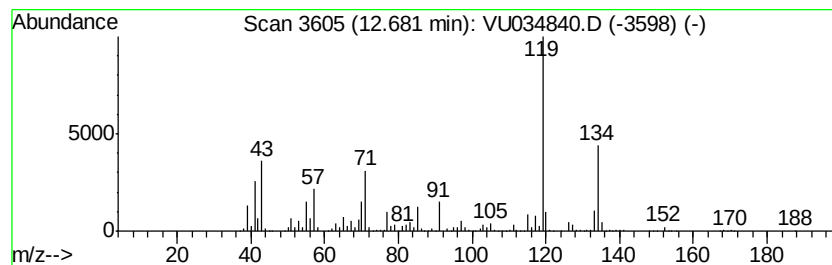
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 44 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 15

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.68 | 58.98 ug/L | 2413690 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 94   |
| 2    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 94   |
| 3    |    | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 93   |
| 4    |    | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 93   |
| 5    |    | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

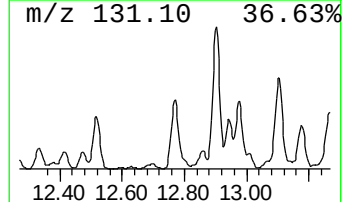
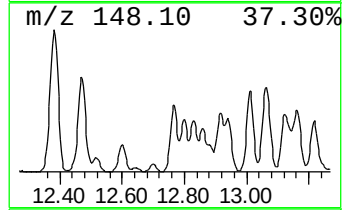
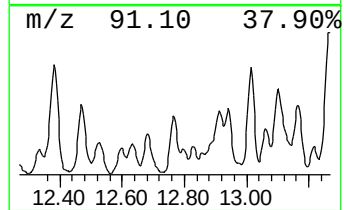
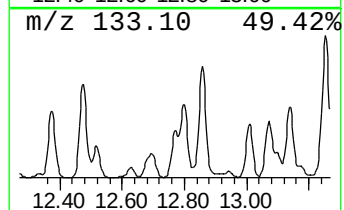
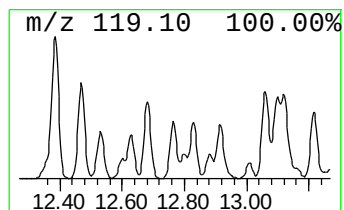
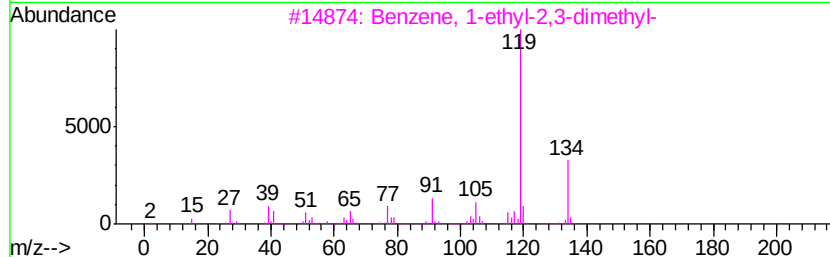
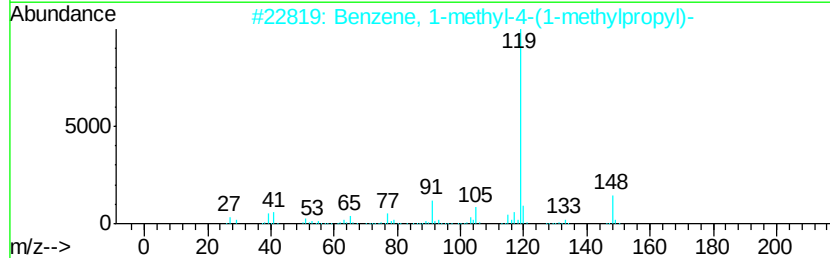
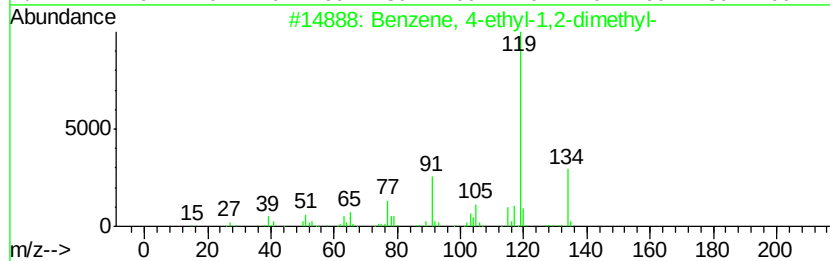
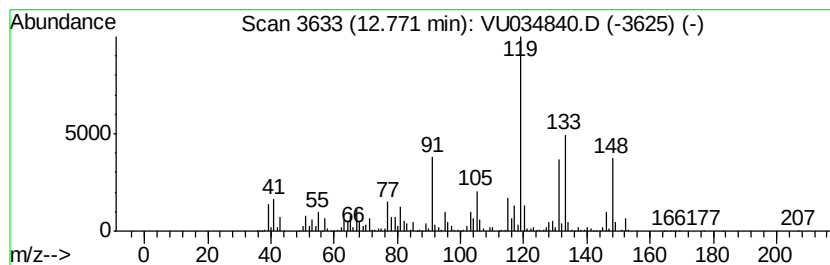
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 45 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 29

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 12.77 | 34.24 ug/L | 1401300 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 64   |
| 2    |    | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16  | 001595-16-0 | 58   |
| 3    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 50   |
| 4    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 50   |
| 5    |    | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9 | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

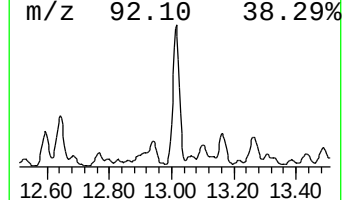
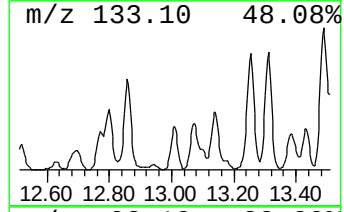
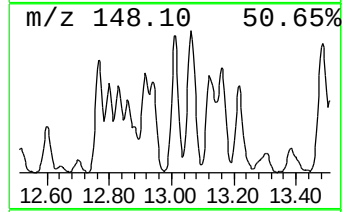
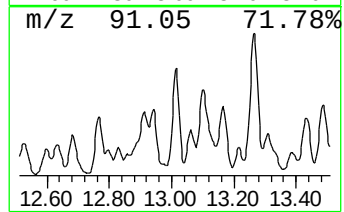
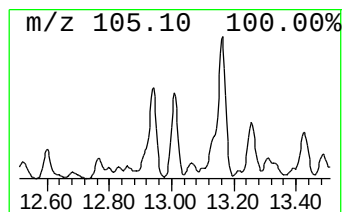
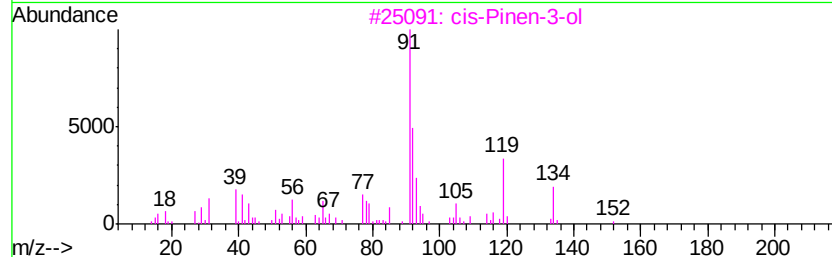
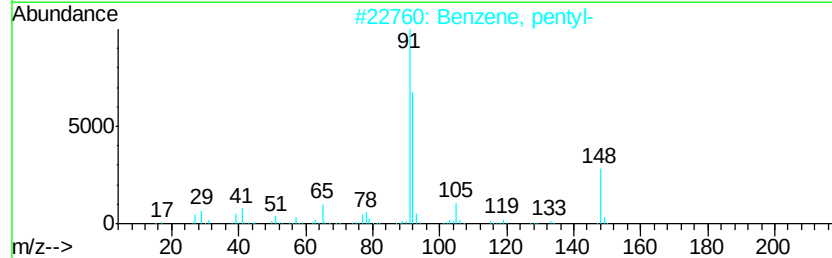
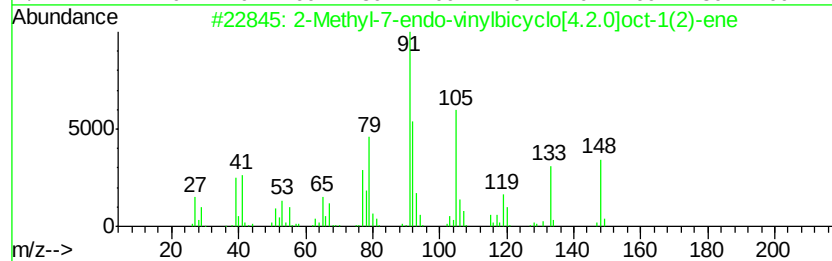
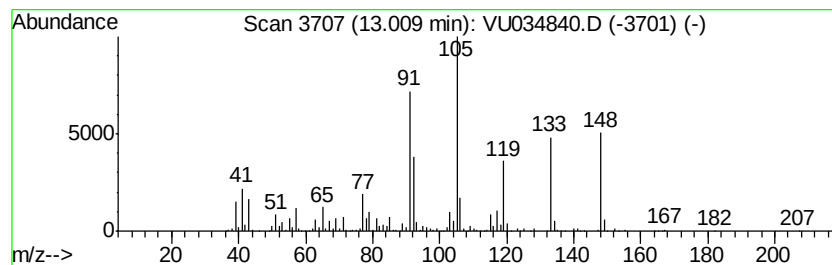
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 46 2-Methyl-7-endo-vinylbicycl... Concentration Rank 28

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 13.01 | 35.46 ug/L | 1451020 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                                    | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 2-Methyl-7-endo-vinylbicyclo[4.2.0]oct-1(2)-ene | 148 | C11H16    | 1000200-99-1 | 91   |
| 2    |    | Benzene, pentyl-                                | 148 | C11H16    | 000538-68-1  | 43   |
| 3    |    | cis-Pinen-3-ol                                  | 152 | C10H16O   | 1000292-85-2 | 41   |
| 4    |    | Benzene, (1-ethylpropyl)-                       | 148 | C11H16    | 001196-58-3  | 38   |
| 5    |    | 2-Nitrophenyl-.beta.-phenylpropi...             | 271 | C15H13NO4 | 040123-51-1  | 38   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

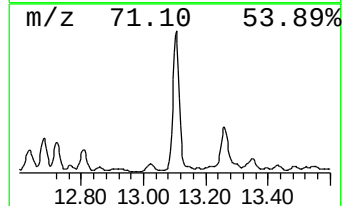
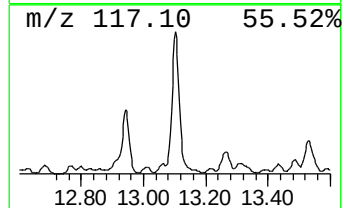
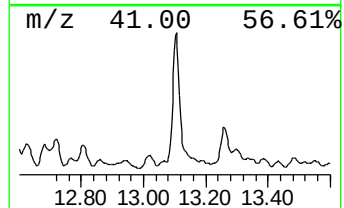
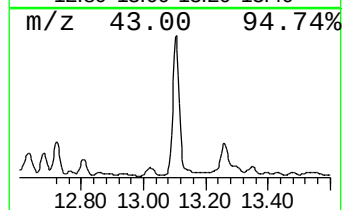
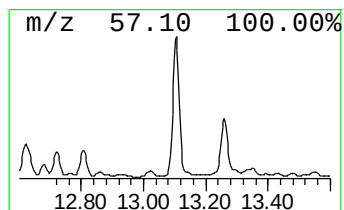
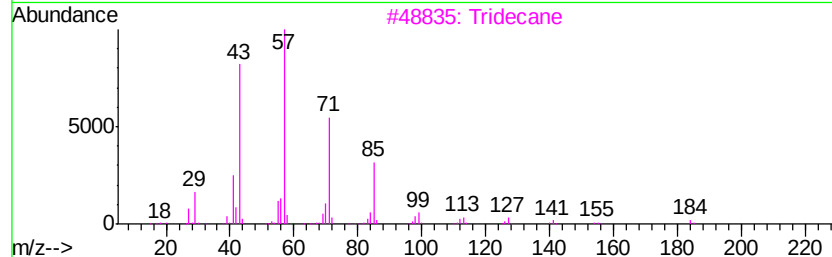
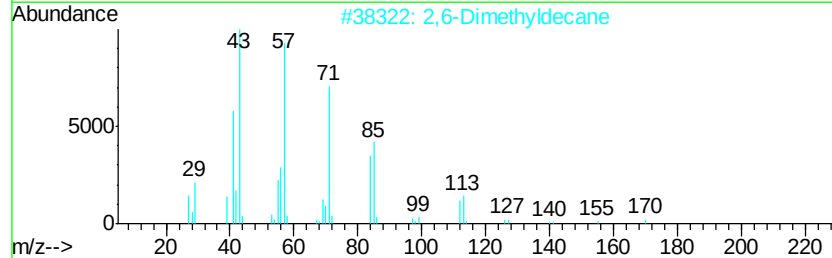
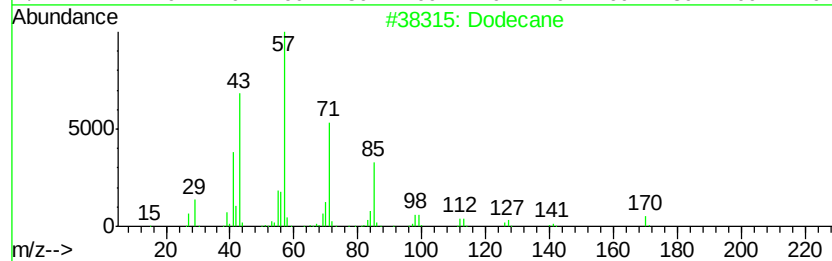
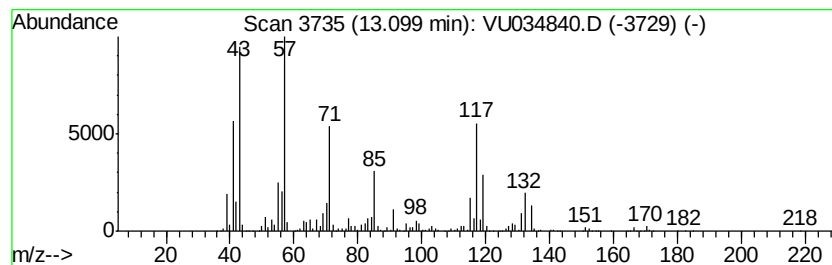
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 13.10 | 186.08 ug/L | 7614820 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane           | 170 | C12H26  | 000112-40-3 | 70   |
| 2    |    | 2,6-Dimethyldecane | 170 | C12H26  | 013150-81-7 | 43   |
| 3    |    | Tridecane          | 184 | C13H28  | 000629-50-5 | 43   |
| 4    |    | Tetradecane        | 198 | C14H30  | 000629-59-4 | 43   |
| 5    |    | Tridecane          | 184 | C13H28  | 000629-50-5 | 43   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

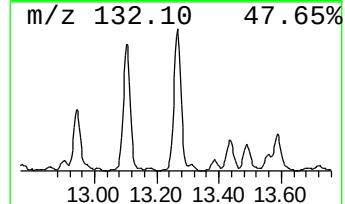
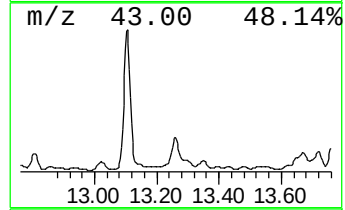
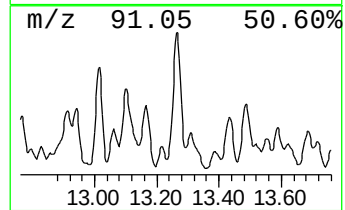
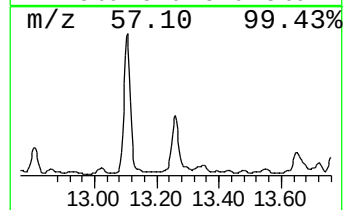
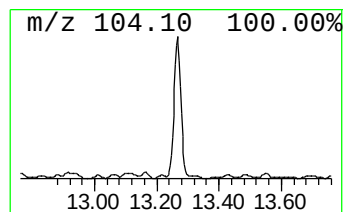
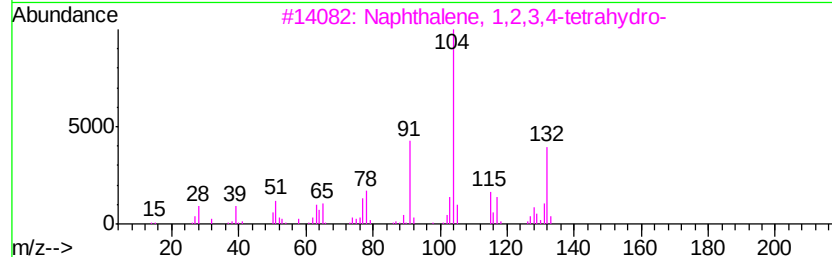
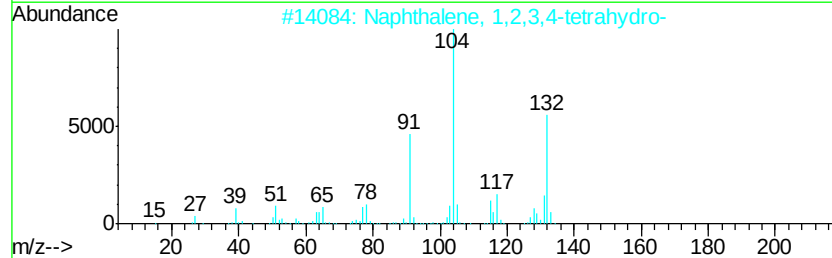
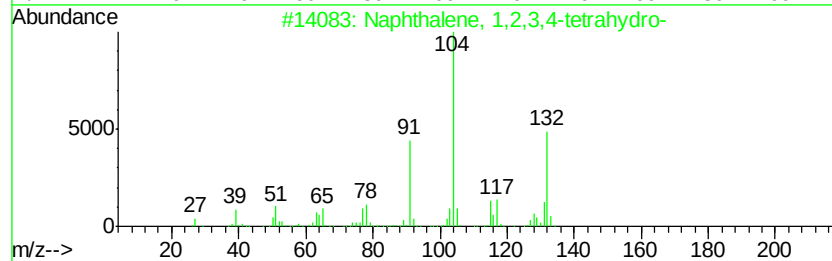
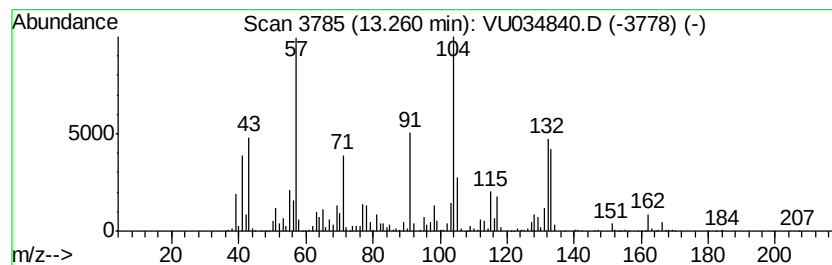
Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 48 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 10

| R.T.      | EstConc                             | Area    | Relative to ISTD       | R.T.        |      |
|-----------|-------------------------------------|---------|------------------------|-------------|------|
| 13.26     | 96.19 ug/L                          | 3936330 | 1,4-Dichlorobenzene-d4 | 11.46       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm                | CAS#        | Qual |
| 1         | Naphthalene, 1,2,3,4-tetrahydro-    | 132     | C10H12                 | 000119-64-2 | 93   |
| 2         | Naphthalene, 1,2,3,4-tetrahydro-    | 132     | C10H12                 | 000119-64-2 | 93   |
| 3         | Naphthalene, 1,2,3,4-tetrahydro-    | 132     | C10H12                 | 000119-64-2 | 89   |
| 4         | Naphthalene, 1,2,3,4-tetrahydro-    | 132     | C10H12                 | 000119-64-2 | 70   |
| 5         | Isoquinoline, 1,2,3,4-tetrahydro... | 147     | C10H13N                | 029726-60-1 | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

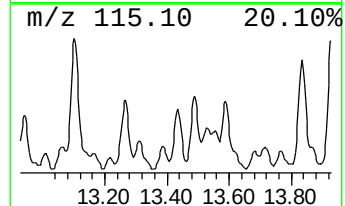
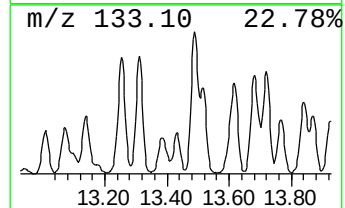
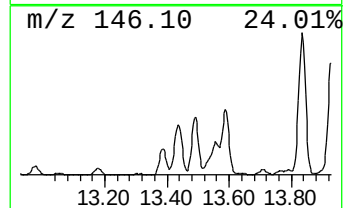
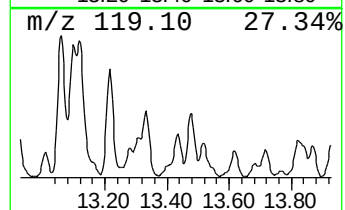
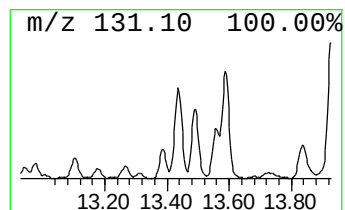
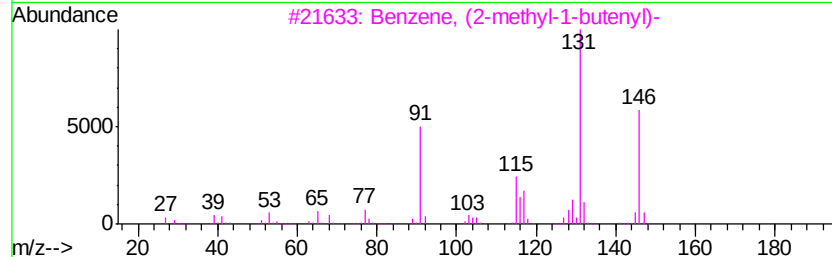
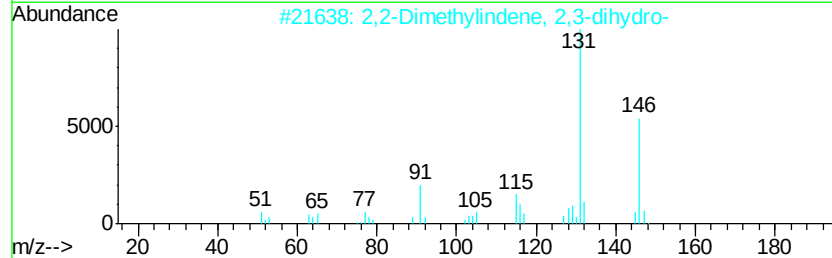
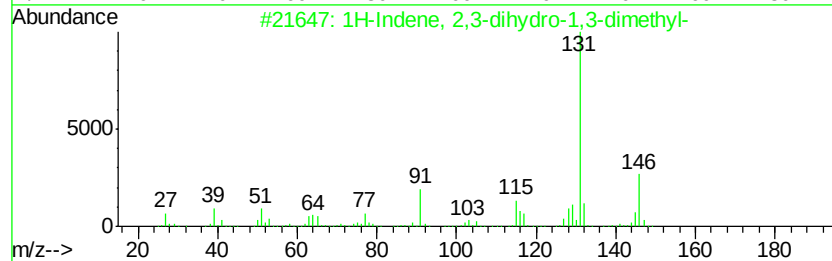
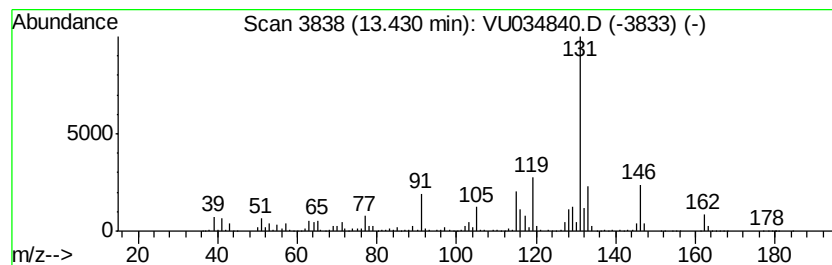
Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 49 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 23

| R.T.    | EstConc    | Area                                | Relative to ISTD       | R.T.    |             |      |
|---------|------------|-------------------------------------|------------------------|---------|-------------|------|
| 13.43   | 40.74 ug/L | 1667140                             | 1,4-Dichlorobenzene-d4 | 11.46   |             |      |
| Hit# of | 5          | Tentative ID                        | MW                     | MolForm | CAS#        | Qual |
| 1       |            | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146                    | C11H14  | 004175-53-5 | 89   |
| 2       |            | 2,2-Dimethylindene, 2,3-dihydro-    | 146                    | C11H14  | 020836-11-7 | 89   |
| 3       |            | Benzene, (2-methyl-1-butenyl)-      | 146                    | C11H14  | 056253-64-6 | 89   |
| 4       |            | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146                    | C11H14  | 006682-71-9 | 86   |
| 5       |            | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146                    | C11H14  | 017059-48-2 | 76   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

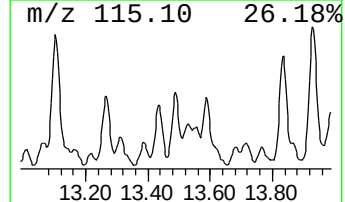
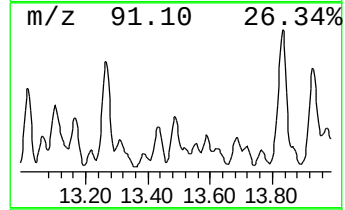
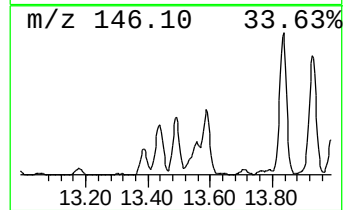
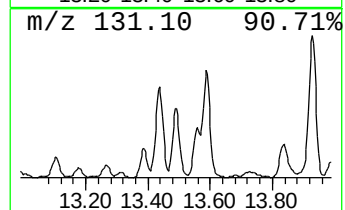
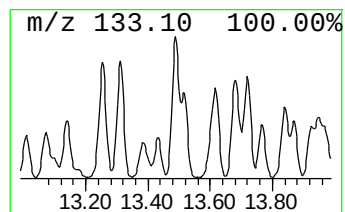
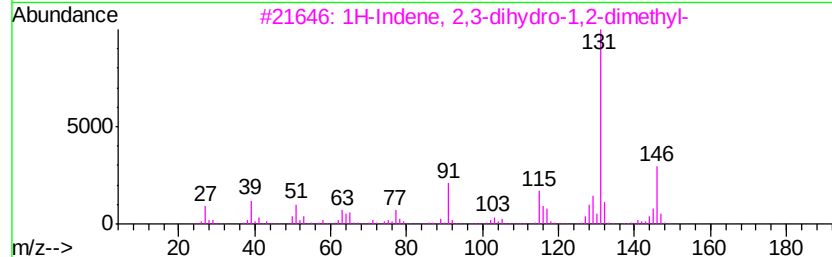
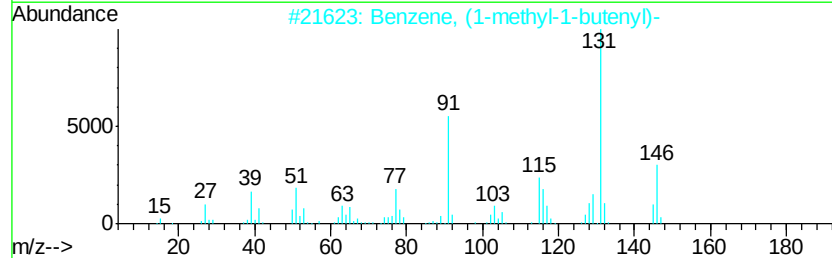
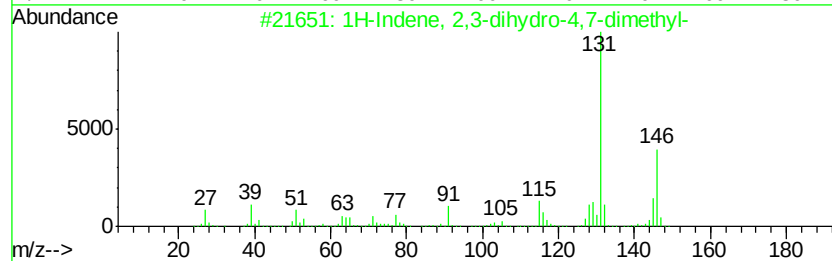
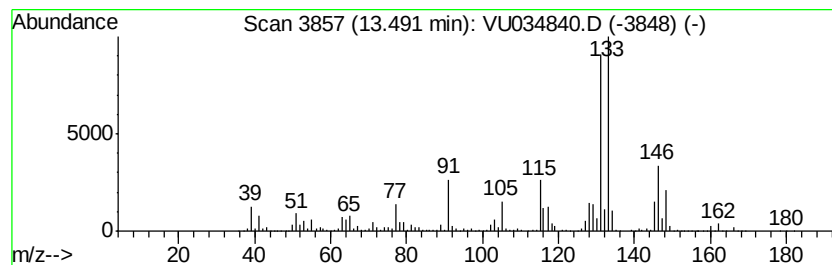
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 50 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 16

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 13.49 | 58.37 ug/L | 2388690 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 89   |
| 2    |    | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 87   |
| 3    |    | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 64   |
| 4    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001559-81-5 | 50   |
| 5    |    | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

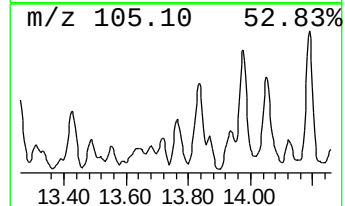
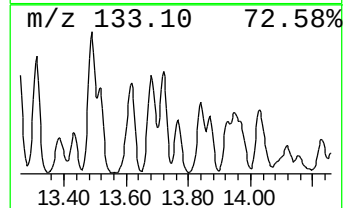
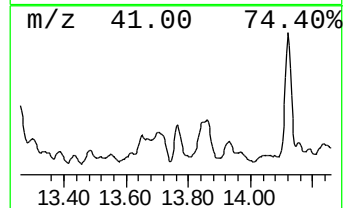
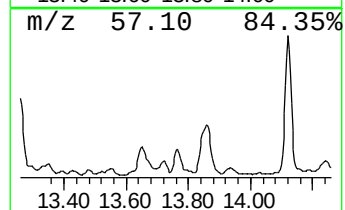
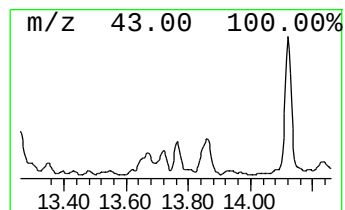
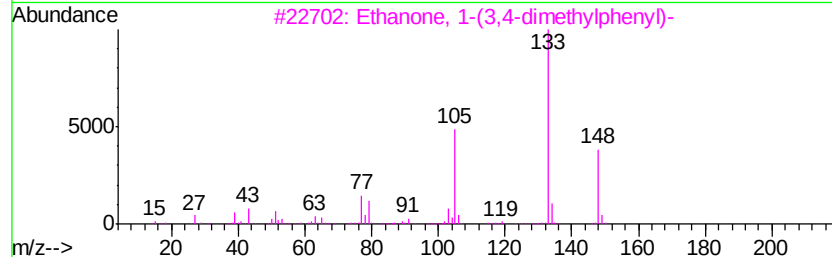
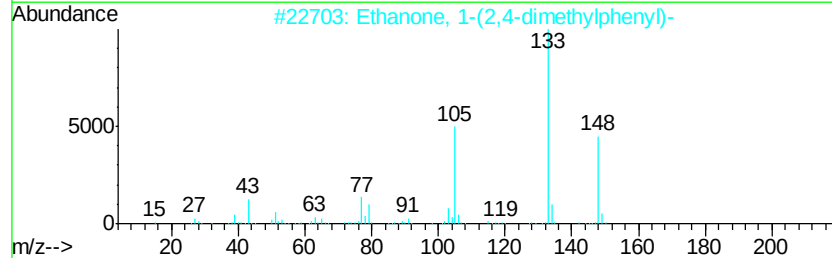
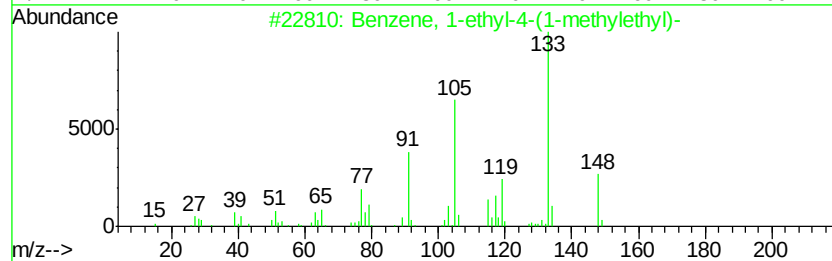
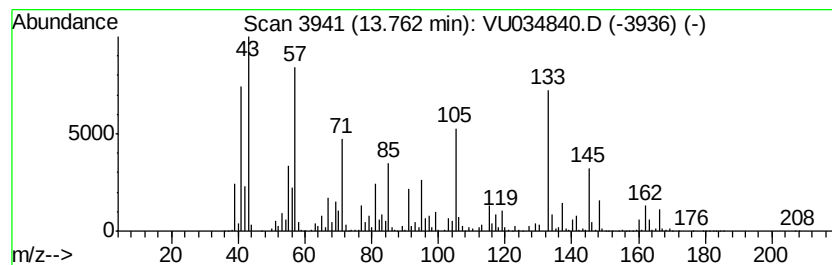
Instrument :  
MSVOA\_U  
ClientSampled :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 51 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 22

| R.T.    | EstConc    | Area                                | Relative to ISTD       | R.T.           |
|---------|------------|-------------------------------------|------------------------|----------------|
| 13.76   | 42.20 ug/L | 1726930                             | 1,4-Dichlorobenzene-d4 | 11.46          |
| Hit# of | 5          | Tentative ID                        | MW MolForm             | CAS# Qual      |
| 1       |            | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 C11H16             | 004218-48-8 50 |
| 2       |            | Ethanone, 1-(2,4-dimethylphenyl)-   | 148 C10H12O            | 000089-74-7 42 |
| 3       |            | Ethanone, 1-(3,4-dimethylphenyl)-   | 148 C10H12O            | 003637-01-2 42 |
| 4       |            | Benzene, 1-ethyl-3-(1-methylethyl)- | 148 C11H16             | 004920-99-4 38 |
| 5       |            | Benzene, 1,3-dimethyl-5-(1-methy... | 148 C11H16             | 004706-90-5 38 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

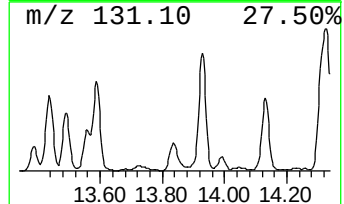
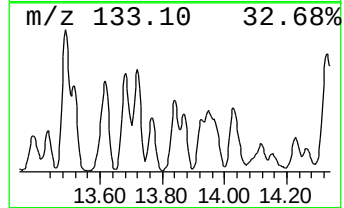
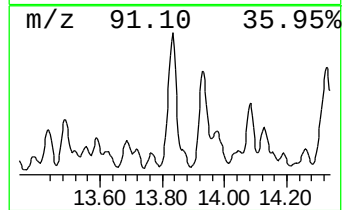
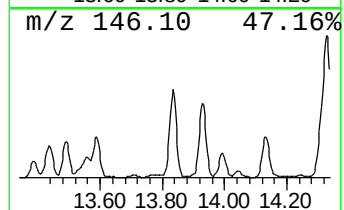
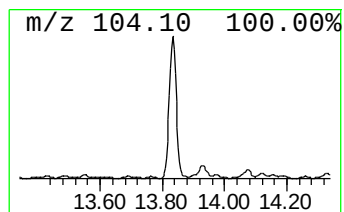
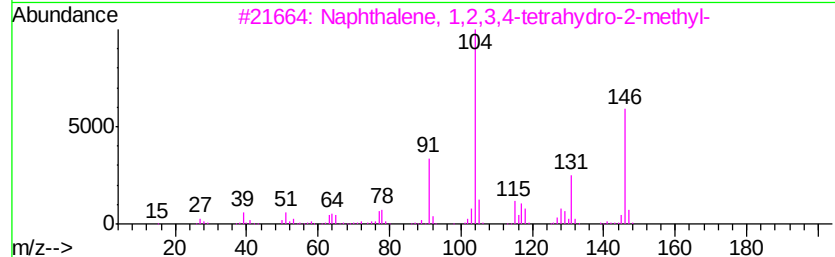
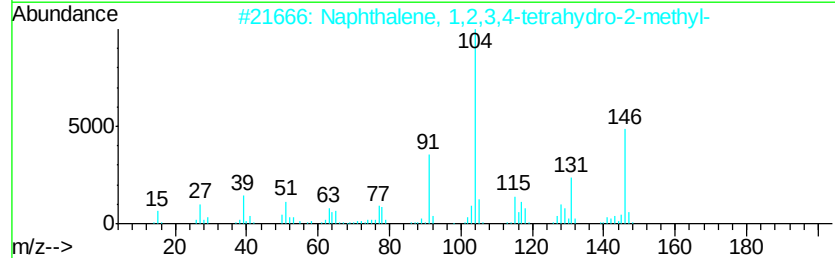
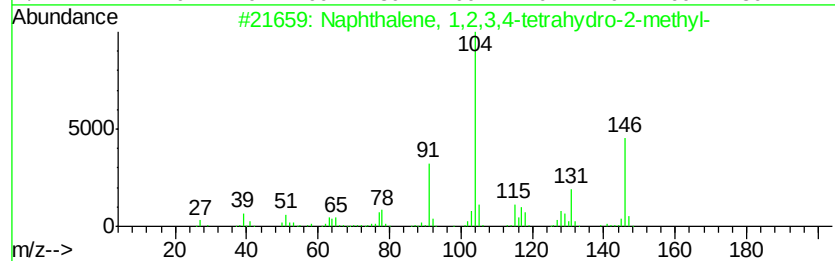
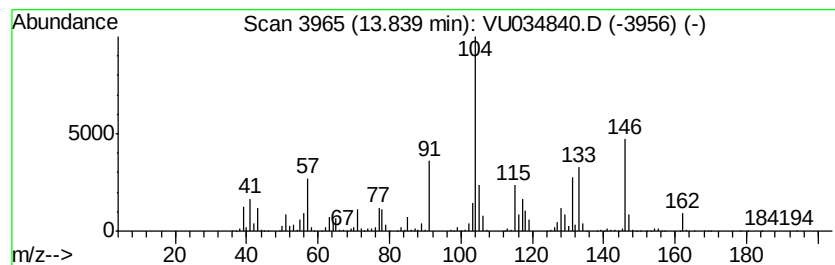
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 52 Naphthalene, 1,2,3,4-tetra... Concentration Rank 8

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 13.84 | 98.71 ug/L | 4039380 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 003877-19-8 | 81   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 003877-19-8 | 81   |
| 3    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 003877-19-8 | 76   |
| 4    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 003877-19-8 | 68   |
| 5    |    | Pyrazolo[1,5-a]pyridine, 2,3-dim... | 146 | C9H10N2 | 028537-56-6 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22uL/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

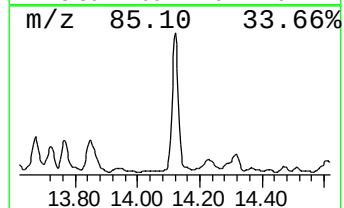
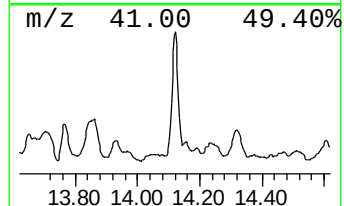
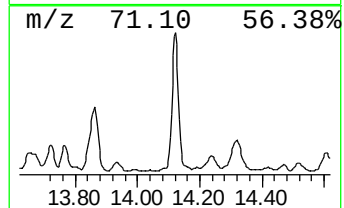
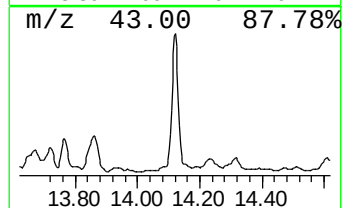
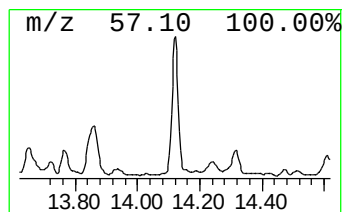
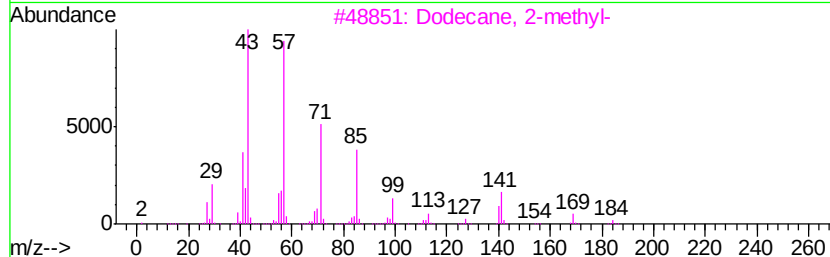
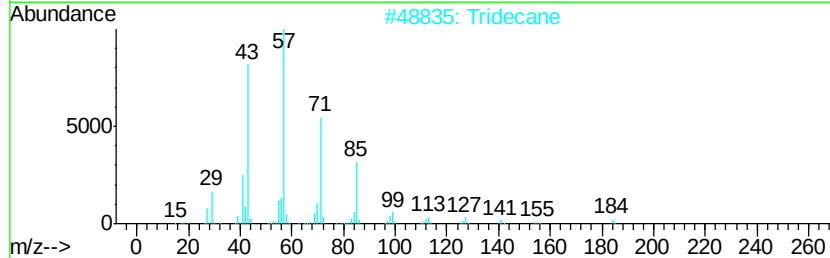
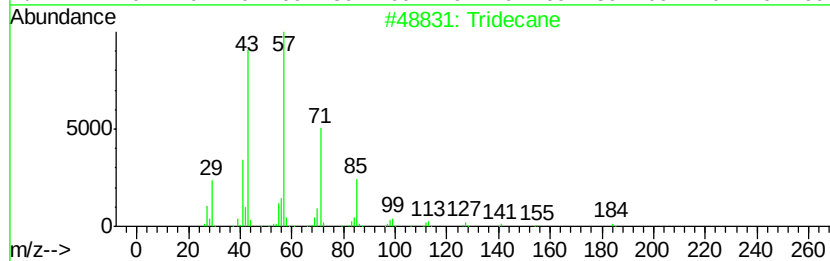
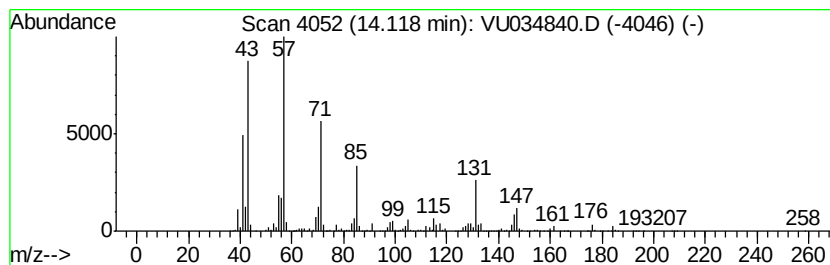
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 14.12 | 125.73 ug/L | 5145170 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane           | 184 | C13H28  | 000629-50-5 | 92   |
| 2    |    | Tridecane           | 184 | C13H28  | 000629-50-5 | 78   |
| 3    |    | Dodecane, 2-methyl- | 184 | C13H28  | 001560-97-0 | 53   |
| 4    |    | Tridecane           | 184 | C13H28  | 000629-50-5 | 52   |
| 5    |    | Hexadecane          | 226 | C16H34  | 000544-76-3 | 46   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

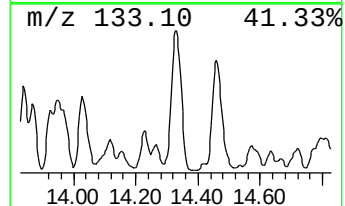
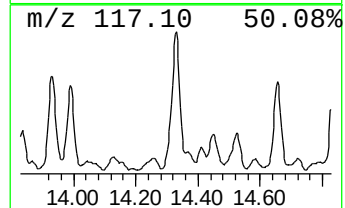
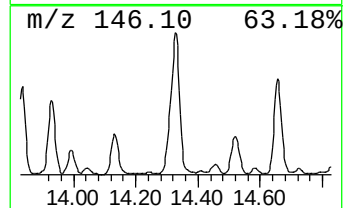
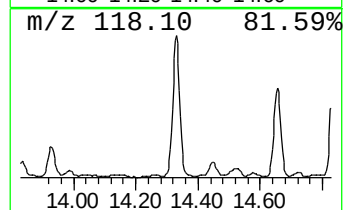
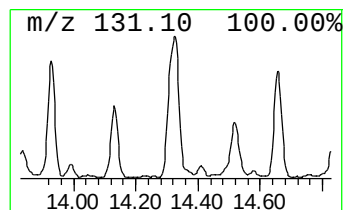
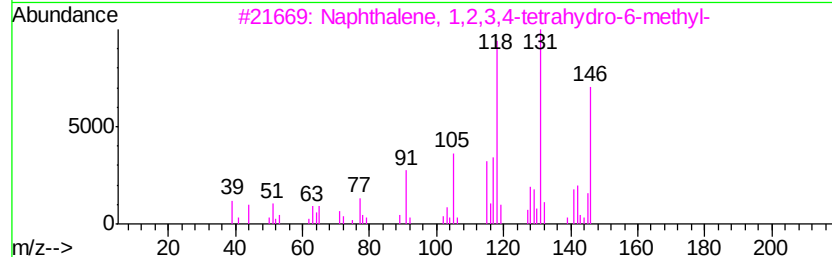
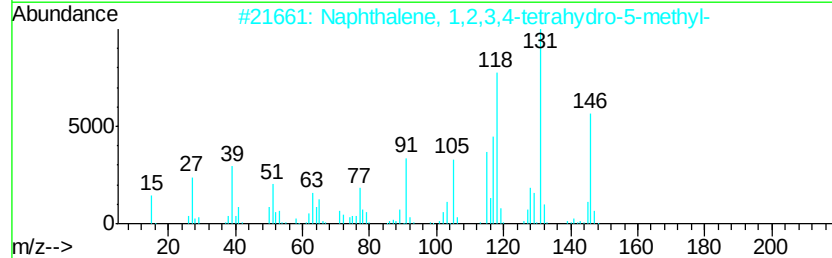
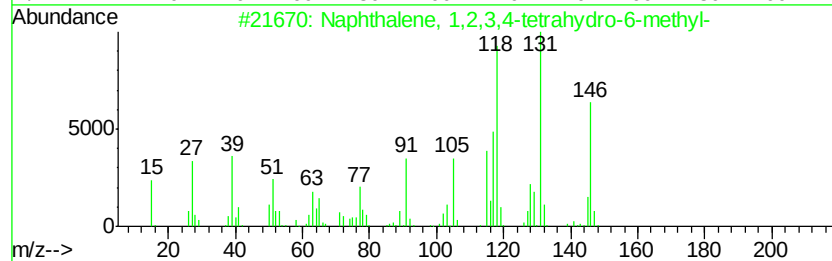
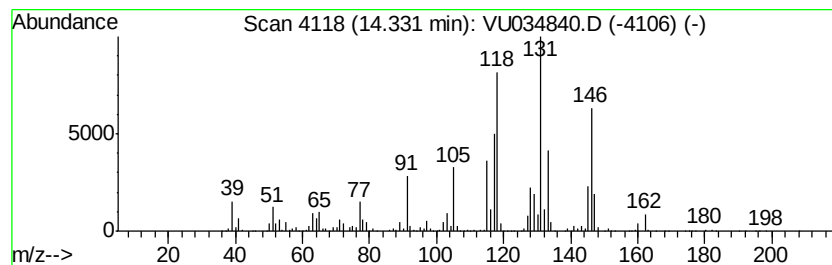
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 55 Naphthalene, 1,2,3,4-tetra... Concentration Rank 2

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 14.33 | 181.16 ug/L | 7413720 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9 | 93   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5 | 91   |
| 3    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9 | 89   |
| 4    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9 | 70   |
| 5    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9 | 70   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU092619\  
Data File : VU034840.D  
Acq On : 26 Sep 2019 12:48  
Operator : JC/SP  
Sample : K4984-08  
Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
Quant Title : VOC Analysis

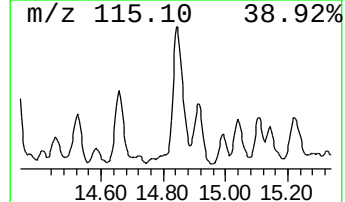
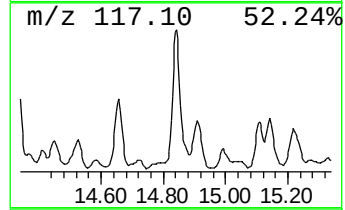
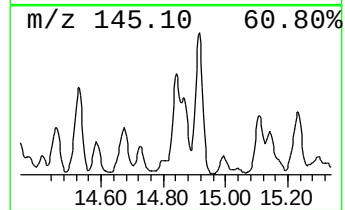
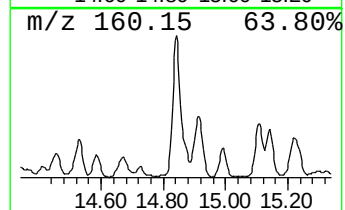
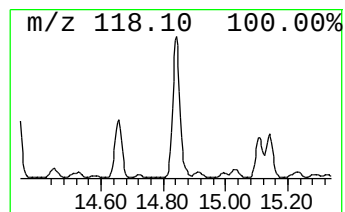
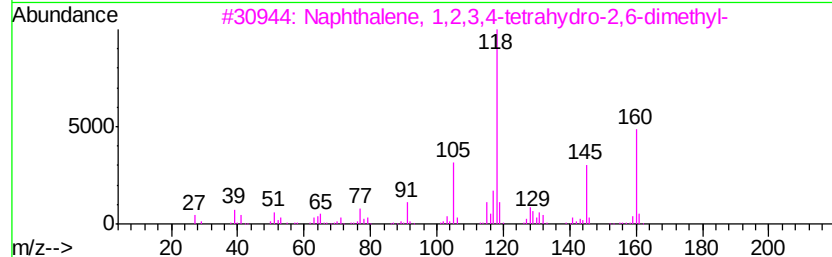
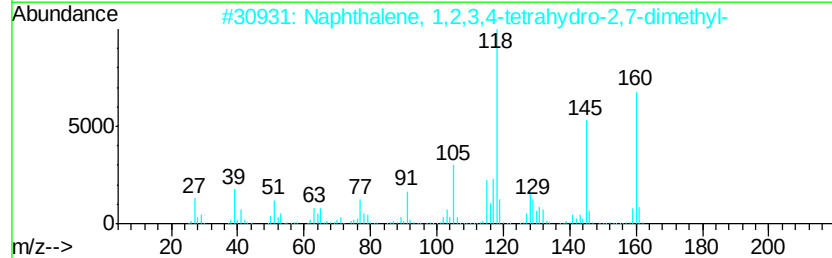
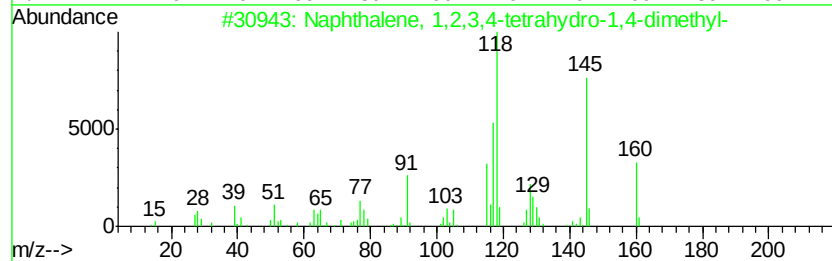
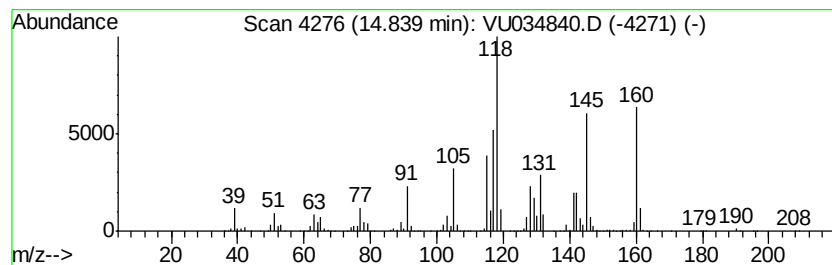
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 57 Naphthalene, 1,2,3,4-tetra... Concentration Rank 4

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 14.84 | 144.69 ug/L | 5920990 | 1,4-Dichlorobenzene-d4 | 11.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 004175-54-6 | 95   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 013065-07-1 | 86   |
| 3    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 007524-63-2 | 55   |
| 4    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 007524-63-2 | 53   |
| 5    |    | Benzene, (3-methyl-1-methylenebu... | 160 | C12H16  | 038212-14-5 | 53   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU092619\  
 Data File : VU034840.D  
 Acq On : 26 Sep 2019 12:48  
 Operator : JC/SP  
 Sample : K4984-08  
 Misc : 6.22g/5mL/100uL/5.0mL/MSVOA\_U/MEOH  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BEDL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMULM091919WMA.M  
 Quant Title : VOC Analysis

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

|                      |       |         |       |          |   | --Internal Standard-- |         |      |  |
|----------------------|-------|---------|-------|----------|---|-----------------------|---------|------|--|
| TIC Top Hit name     | RT    | EstConc | Units | Response | # | RT                    | Resp    | Conc |  |
| (DEL) Alkane: Str... | 5.75  | 4.6     | ug/L  | 111487   | 1 | 5.86                  | 1222970 | 50.0 |  |
| (DEL) Alkane: Cyc... | 6.61  | 2.5     | ug/L  | 62475    | 1 | 5.86                  | 1222970 | 50.0 |  |
| (DEL) Alkane: Cyc... | 6.71  | 4.7     | ug/L  | 116001   | 1 | 5.86                  | 1222970 | 50.0 |  |
| (DEL) Alkane: Cyc... | 6.88  | 6.0     | ug/L  | 147888   | 1 | 5.86                  | 1222970 | 50.0 |  |
| (DEL) Alkane: Str... | 7.08  | 2.9     | ug/L  | 71852    | 1 | 5.86                  | 1222970 | 50.0 |  |
| (DEL) Alkane: Str... | 7.15  | 10.6    | ug/L  | 259267   | 1 | 5.86                  | 1222970 | 50.0 |  |
| (DEL) Alkane: Str... | 7.31  | 14.8    | ug/L  | 360653   | 1 | 5.86                  | 1222970 | 50.0 |  |
| (DEL) Alkane: Cyc... | 7.42  | 2.9     | ug/L  | 70697    | 1 | 5.86                  | 1222970 | 50.0 |  |
| (DEL) Alkane: Cyc... | 7.68  | 7.8     | ug/L  | 230694   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Cyc... | 7.89  | 17.2    | ug/L  | 510322   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Cyc... | 8.02  | 13.5    | ug/L  | 400083   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Str... | 8.20  | 6.3     | ug/L  | 185790   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Str... | 8.43  | 6.4     | ug/L  | 190015   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Cyc... | 8.52  | 24.2    | ug/L  | 718363   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Cyc... | 8.58  | 29.9    | ug/L  | 886971   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Cyc... | 8.73  | 6.7     | ug/L  | 198168   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Str... | 8.79  | 12.7    | ug/L  | 376618   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Str... | 8.89  | 36.8    | ug/L  | 1092010  | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Str... | 9.01  | 32.7    | ug/L  | 971426   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Str... | 9.42  | 110.3   | ug/L  | 3274570  | 2 | 9.07                  | 1483920 | 50.0 |  |
| unknown-01           | 9.54  | 8.6     | ug/L  | 254207   | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Cyc... | 9.69  | 47.8    | ug/L  | 1418420  | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Str... | 9.93  | 62.2    | ug/L  | 1844890  | 2 | 9.07                  | 1483920 | 50.0 |  |
| (DEL) Alkane: Cyc... | 10.02 | 39.9    | ug/L  | 1184770  | 2 | 9.07                  | 1483920 | 50.0 |  |
| Benzene, propyl-     | 10.57 | 19.7    | ug/L  | 805039   | 3 | 11.46                 | 2046160 | 50.0 |  |
| Benzene, 1-ethyl-... | 10.66 | 75.0    | ug/L  | 3067180  | 3 | 11.46                 | 2046160 | 50.0 |  |
| (DEL) Alkane: Cyc... | 10.73 | 43.7    | ug/L  | 1788310  | 3 | 11.46                 | 2046160 | 50.0 |  |
| (DEL) Alkane: Str... | 10.79 | 127.6   | ug/L  | 5223540  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Benzene, 1-ethyl-... | 10.95 | 52.1    | ug/L  | 2131000  | 3 | 11.46                 | 2046160 | 50.0 |  |
| (DEL) Alkane: Str... | 11.09 | 35.8    | ug/L  | 1466830  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Benzene, 1,2,3-tr... | 11.13 | 42.9    | ug/L  | 1756550  | 3 | 11.46                 | 2046160 | 50.0 |  |
| (DEL) Alkane: Cyc... | 11.38 | 29.0    | ug/L  | 1187980  | 3 | 11.46                 | 2046160 | 50.0 |  |
| (DEL) Alkane: Str... | 11.59 | 25.9    | ug/L  | 1058970  | 3 | 11.46                 | 2046160 | 50.0 |  |
| (DEL) Alkane: Str... | 11.68 | 30.3    | ug/L  | 1238260  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Benzene, 1,2-diet... | 11.76 | 30.1    | ug/L  | 1231590  | 3 | 11.46                 | 2046160 | 50.0 |  |
| (DEL) Alkane: Str... | 12.00 | 150.0   | ug/L  | 6138190  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Benzene, 2-ethyl-... | 12.12 | 27.8    | ug/L  | 1139220  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Indan, 1-methyl-     | 12.34 | 37.4    | ug/L  | 1531410  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Benzene, 1,3-diet... | 12.47 | 67.9    | ug/L  | 2778490  | 3 | 11.46                 | 2046160 | 50.0 |  |
| (DEL) Alkane: Cyc... | 12.59 | 48.7    | ug/L  | 1994060  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Benzene, 1,2,4,5-... | 12.68 | 59.0    | ug/L  | 2413690  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Benzene, 4-ethyl-... | 12.77 | 34.2    | ug/L  | 1401300  | 3 | 11.46                 | 2046160 | 50.0 |  |
| 2-Methyl-7-endo-v... | 13.01 | 35.5    | ug/L  | 1451020  | 3 | 11.46                 | 2046160 | 50.0 |  |
| (DEL) Alkane: Str... | 13.10 | 186.1   | ug/L  | 7614820  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Naphthalene, 1,2,... | 13.26 | 96.2    | ug/L  | 3936330  | 3 | 11.46                 | 2046160 | 50.0 |  |
| 1H-Indene, 2,3-di... | 13.43 | 40.7    | ug/L  | 1667140  | 3 | 11.46                 | 2046160 | 50.0 |  |
| 1H-Indene, 2,3-di... | 13.49 | 58.4    | ug/L  | 2388690  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Benzene, 1-ethyl-... | 13.76 | 42.2    | ug/L  | 1726930  | 3 | 11.46                 | 2046160 | 50.0 |  |
| Naphthalene, 1,2,... | 13.84 | 98.7    | ug/L  | 4039380  | 3 | 11.46                 | 2046160 | 50.0 |  |

|                      |       |            |         |   |       |         |      |
|----------------------|-------|------------|---------|---|-------|---------|------|
| (DEL) Alkane: Str... | 14.12 | 125.7 ug/L | 5145170 | 3 | 11.46 | 2046160 | 50.0 |
| Naphthalene, 1,2,... | 14.33 | 181.2 ug/L | 7413720 | 3 | 11.46 | 2046160 | 50.0 |
| Naphthalene, 1,2,... | 14.84 | 144.7 ug/L | 5920990 | 3 | 11.46 | 2046160 | 50.0 |

Instrument :  
MSVOA\_U  
ClientSampleId :  
BEDL2