

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
 Data File : VX026043.D  
 Acq On : 25 Dec 2021 18:51  
 Operator : JC/MD  
 Sample : M5213-05 10X  
 Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 BGLD7

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
 Title : VOC Analysis

Signal : TIC: VX026043.D\data.ms

| 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         | 1.368       | 43            | 46          | 54           | rVB      | 117026         | 120680        | 0.27%           | 0.042%        |
| 2         | 1.654       | 90            | 93          | 101          | rVB      | 102648         | 120168        | 0.27%           | 0.042%        |
| 3         | 2.307       | 192           | 200         | 214          | rBV      | 201363         | 332971        | 0.74%           | 0.116%        |
| 4         | 4.465       | 544           | 554         | 570          | rBV2     | 79982          | 267320        | 0.59%           | 0.093%        |
| 5         | 5.062       | 639           | 652         | 670          | rBV3     | 110012         | 394202        | 0.87%           | 0.137%        |
| 6         | 5.495       | 709           | 723         | 737          | rVB7     | 5315           | 21555         | 0.05%           | 0.008%        |
| 7         | 5.970       | 788           | 801         | 810          | rBV3     | 340333         | 1025256       | 2.26%           | 0.357%        |
| 8         | 6.611       | 898           | 906         | 916          | rVB3     | 14801          | 36845         | 0.08%           | 0.013%        |
| 9         | 6.763       | 921           | 931         | 944          | rBV      | 247966         | 601471        | 1.33%           | 0.210%        |
| 10        | 7.312       | 1012          | 1021        | 1028         | rVV      | 210796         | 459843        | 1.02%           | 0.160%        |
| 11        | 7.379       | 1028          | 1032        | 1044         | rVB2     | 30851          | 66953         | 0.15%           | 0.023%        |
| 12        | 8.275       | 1173          | 1179        | 1183         | rBV2     | 42863          | 70874         | 0.16%           | 0.025%        |
| 13        | 8.324       | 1183          | 1187        | 1194         | rVB      | 130060         | 218170        | 0.48%           | 0.076%        |
| 14        | 8.440       | 1201          | 1206        | 1215         | rBV      | 36268          | 64255         | 0.14%           | 0.022%        |
| 15        | 8.598       | 1226          | 1232        | 1235         | rBV      | 60762          | 103993        | 0.23%           | 0.036%        |
| 16        | 8.647       | 1235          | 1240        | 1247         | rVB      | 508456         | 836556        | 1.85%           | 0.292%        |
| 17        | 8.720       | 1247          | 1252        | 1257         | rBV      | 174775         | 262191        | 0.58%           | 0.091%        |
| 18        | 8.775       | 1257          | 1261        | 1265         | rVB2     | 13964          | 21075         | 0.05%           | 0.007%        |
| 19        | 8.915       | 1278          | 1284        | 1287         | rBV      | 215602         | 312026        | 0.69%           | 0.109%        |
| 20        | 8.952       | 1287          | 1290        | 1293         | rBV      | 64904          | 83951         | 0.19%           | 0.029%        |
| 21        | 9.104       | 1304          | 1315        | 1320         | rBV2     | 31173          | 52059         | 0.11%           | 0.018%        |
| 22        | 9.293       | 1339          | 1346        | 1351         | rVB2     | 42615          | 73488         | 0.16%           | 0.026%        |
| 23        | 9.385       | 1352          | 1361        | 1371         | rBV      | 563030         | 840971        | 1.86%           | 0.293%        |
| 24        | 9.500       | 1375          | 1380        | 1386         | rBV2     | 146779         | 300486        | 0.66%           | 0.105%        |
| 25        | 9.561       | 1386          | 1390        | 1395         | rVB      | 195102         | 281292        | 0.62%           | 0.098%        |
| 26        | 9.622       | 1395          | 1400        | 1404         | rBV      | 259449         | 405676        | 0.90%           | 0.141%        |
| 27        | 9.762       | 1418          | 1423        | 1427         | rVB      | 51005          | 74360         | 0.16%           | 0.026%        |
| 28        | 9.823       | 1427          | 1433        | 1438         | rBV3     | 285162         | 614082        | 1.36%           | 0.214%        |
| 29        | 9.909       | 1442          | 1447        | 1452         | rVB      | 683253         | 989047        | 2.18%           | 0.345%        |
| 30        | 10.012      | 1457          | 1464        | 1467         | rBV      | 641937         | 894853        | 1.98%           | 0.312%        |
| 31        | 10.055      | 1467          | 1471        | 1476         | rVB      | 573715         | 758529        | 1.68%           | 0.264%        |
| 32        | 10.122      | 1476          | 1482        | 1487         | rVB4     | 66561          | 150494        | 0.33%           | 0.052%        |
| 33        | 10.201      | 1487          | 1495        | 1500         | rBV      | 9790381        | 13829915      | 30.54%          | 4.819%        |
| 34        | 10.311      | 1503          | 1513        | 1518         | rBV      | 28552636       | 45280737      | 100.00%         | 15.778%       |
| 35        | 10.366      | 1518          | 1522        | 1532         | rVB      | 4466511        | 5833715       | 12.88%          | 2.033%        |
| 36        | 10.457      | 1532          | 1537        | 1543         | rBV      | 220573         | 329780        | 0.73%           | 0.115%        |

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
 Data File : VX026043.D  
 Acq On : 25 Dec 2021 18:51  
 Operator : JC/MD  
 Sample : M5213-05 10X  
 Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 BGLD7

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
 Title : VOC Analysis

|    |        |      |      |      |      |          |          |        |        |
|----|--------|------|------|------|------|----------|----------|--------|--------|
| 37 | 10.537 | 1546 | 1550 | 1552 | rBV2 | 123090   | 174409   | 0.39%  | 0.061% |
| 38 | 10.592 | 1552 | 1559 | 1563 | rBV2 | 1162433  | 1900365  | 4.20%  | 0.662% |
| 39 | 10.646 | 1563 | 1568 | 1576 | rBV  | 10064204 | 13596678 | 30.03% | 4.738% |
| 40 | 10.714 | 1577 | 1579 | 1582 | rVB  | 335075   | 369966   | 0.82%  | 0.129% |
| 41 | 10.781 | 1582 | 1590 | 1595 | rBV2 | 2385699  | 3705167  | 8.18%  | 1.291% |
| 42 | 10.860 | 1595 | 1603 | 1606 | rBV  | 2204847  | 3814067  | 8.42%  | 1.329% |
| 43 | 10.896 | 1606 | 1609 | 1614 | rVB  | 1151089  | 1566755  | 3.46%  | 0.546% |
| 44 | 10.964 | 1614 | 1620 | 1623 | rBV2 | 969787   | 1447393  | 3.20%  | 0.504% |
| 45 | 11.006 | 1623 | 1627 | 1628 | rBV2 | 445673   | 514580   | 1.14%  | 0.179% |
| 46 | 11.031 | 1628 | 1631 | 1635 | rVB  | 896269   | 1025400  | 2.26%  | 0.357% |
| 47 | 11.085 | 1635 | 1640 | 1642 | rBV2 | 1738841  | 2523561  | 5.57%  | 0.879% |
| 48 | 11.110 | 1642 | 1644 | 1650 | rVB  | 2402140  | 3391535  | 7.49%  | 1.182% |
| 49 | 11.195 | 1650 | 1658 | 1661 | rBV  | 2795249  | 4395571  | 9.71%  | 1.532% |
| 50 | 11.305 | 1672 | 1676 | 1680 | rBV  | 2840892  | 3654325  | 8.07%  | 1.273% |
| 51 | 11.341 | 1680 | 1682 | 1685 | rVV3 | 531316   | 773879   | 1.71%  | 0.270% |
| 52 | 11.384 | 1685 | 1689 | 1695 | rVV  | 9660365  | 17166810 | 37.91% | 5.982% |
| 53 | 11.457 | 1695 | 1701 | 1703 | rVV2 | 6505832  | 10761701 | 23.77% | 3.750% |
| 54 | 11.482 | 1703 | 1705 | 1710 | rVB  | 10908235 | 12407205 | 27.40% | 4.323% |
| 55 | 11.579 | 1718 | 1721 | 1722 | rBV  | 187012   | 194309   | 0.43%  | 0.068% |
| 56 | 11.616 | 1722 | 1727 | 1735 | rVB  | 4353397  | 6635258  | 14.65% | 2.312% |
| 57 | 11.689 | 1735 | 1739 | 1741 | rBV2 | 952930   | 1170188  | 2.58%  | 0.408% |
| 58 | 11.719 | 1741 | 1744 | 1746 | rBV  | 3007114  | 3056152  | 6.75%  | 1.065% |
| 59 | 11.762 | 1746 | 1751 | 1760 | rVB  | 18245797 | 25575633 | 56.48% | 8.912% |
| 60 | 11.841 | 1760 | 1764 | 1766 | rBV2 | 975517   | 1284655  | 2.84%  | 0.448% |
| 61 | 11.896 | 1770 | 1773 | 1777 | rVB3 | 1367913  | 1636038  | 3.61%  | 0.570% |
| 62 | 11.945 | 1777 | 1781 | 1784 | rBV2 | 2798630  | 4214570  | 9.31%  | 1.469% |
| 63 | 12.024 | 1789 | 1794 | 1795 | rBV3 | 786397   | 1203583  | 2.66%  | 0.419% |
| 64 | 12.091 | 1800 | 1805 | 1810 | rBV  | 6120933  | 10036989 | 22.17% | 3.497% |
| 65 | 12.177 | 1817 | 1819 | 1826 | rVB2 | 1735103  | 2124923  | 4.69%  | 0.740% |
| 66 | 12.250 | 1826 | 1831 | 1833 | rBV  | 7047632  | 9663963  | 21.34% | 3.367% |
| 67 | 12.317 | 1838 | 1842 | 1849 | rVB3 | 5032682  | 8670255  | 19.15% | 3.021% |
| 68 | 12.427 | 1856 | 1860 | 1864 | rVB  | 5944912  | 6561157  | 14.49% | 2.286% |
| 69 | 12.469 | 1864 | 1867 | 1873 | rVB2 | 1714513  | 2631057  | 5.81%  | 0.917% |
| 70 | 12.536 | 1874 | 1878 | 1879 | rBV  | 1498459  | 1757096  | 3.88%  | 0.612% |
| 71 | 12.561 | 1879 | 1882 | 1888 | rVV3 | 1717414  | 2512338  | 5.55%  | 0.875% |
| 72 | 12.616 | 1888 | 1891 | 1894 | rVB  | 1748465  | 1824018  | 4.03%  | 0.636% |
| 73 | 12.719 | 1900 | 1908 | 1913 | rVB5 | 943093   | 2152531  | 4.75%  | 0.750% |
| 74 | 12.890 | 1932 | 1936 | 1939 | rBV2 | 965578   | 1437293  | 3.17%  | 0.501% |
| 75 | 12.920 | 1939 | 1941 | 1945 | rVV  | 1022145  | 1319969  | 2.92%  | 0.460% |
| 76 | 12.969 | 1945 | 1949 | 1956 | rVB  | 1747688  | 2739047  | 6.05%  | 0.954% |

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
 Data File : VX026043.D  
 Acq On : 25 Dec 2021 18:51  
 Operator : JC/MD  
 Sample : M5213-05 10X  
 Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 BGLD7

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXML122321WMA.M  
 Title : VOC Analysis

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 77  | 13.042 | 1956 | 1961 | 1964 | rBV3 | 384973  | 729037  | 1.61%  | 0.254% |
| 78  | 13.164 | 1975 | 1981 | 1986 | rVB4 | 477781  | 845600  | 1.87%  | 0.295% |
| 79  | 13.213 | 1986 | 1989 | 1992 | rBV2 | 237195  | 241757  | 0.53%  | 0.084% |
| 80  | 13.268 | 1992 | 1998 | 1999 | rBV3 | 826707  | 1094003 | 2.42%  | 0.381% |
| 81  | 13.292 | 2000 | 2002 | 2007 | rVB2 | 1182080 | 1442631 | 3.19%  | 0.503% |
| 82  | 13.372 | 2013 | 2015 | 2018 | rBV2 | 128602  | 145812  | 0.32%  | 0.051% |
| 83  | 13.426 | 2020 | 2024 | 2033 | rVB3 | 700682  | 1283268 | 2.83%  | 0.447% |
| 84  | 13.506 | 2034 | 2037 | 2041 | rBV4 | 112021  | 185473  | 0.41%  | 0.065% |
| 85  | 13.585 | 2046 | 2050 | 2055 | rBV4 | 204342  | 375398  | 0.83%  | 0.131% |
| 86  | 13.676 | 2062 | 2065 | 2069 | rVB5 | 107134  | 138083  | 0.30%  | 0.048% |
| 87  | 13.780 | 2077 | 2082 | 2089 | rVB  | 6146217 | 7894070 | 17.43% | 2.751% |
| 88  | 13.872 | 2089 | 2097 | 2100 | rBV2 | 232282  | 404243  | 0.89%  | 0.141% |
| 89  | 14.042 | 2121 | 2125 | 2128 | rBV  | 168034  | 174946  | 0.39%  | 0.061% |
| 90  | 14.219 | 2150 | 2154 | 2163 | rVB2 | 118923  | 207570  | 0.46%  | 0.072% |
| 91  | 14.487 | 2193 | 2198 | 2202 | rBV4 | 46042   | 107710  | 0.24%  | 0.038% |
| 92  | 14.603 | 2214 | 2217 | 2218 | rBV  | 65582   | 75489   | 0.17%  | 0.026% |
| 93  | 14.640 | 2218 | 2223 | 2233 | rVB  | 7483250 | 9182139 | 20.28% | 3.200% |
| 94  | 14.725 | 2233 | 2237 | 2240 | rBV  | 113817  | 147389  | 0.33%  | 0.051% |
| 95  | 14.780 | 2240 | 2246 | 2255 | rVB  | 2560657 | 3336072 | 7.37%  | 1.162% |
| 96  | 15.207 | 2312 | 2316 | 2322 | rVB  | 309548  | 426102  | 0.94%  | 0.148% |
| 97  | 15.377 | 2340 | 2344 | 2348 | rBV  | 117161  | 159536  | 0.35%  | 0.056% |
| 98  | 15.481 | 2356 | 2361 | 2371 | rBV  | 182242  | 302207  | 0.67%  | 0.105% |
| 99  | 15.615 | 2378 | 2383 | 2387 | rBV  | 145230  | 235365  | 0.52%  | 0.082% |
| 100 | 15.652 | 2387 | 2389 | 2398 | rVB2 | 82832   | 124946  | 0.28%  | 0.044% |

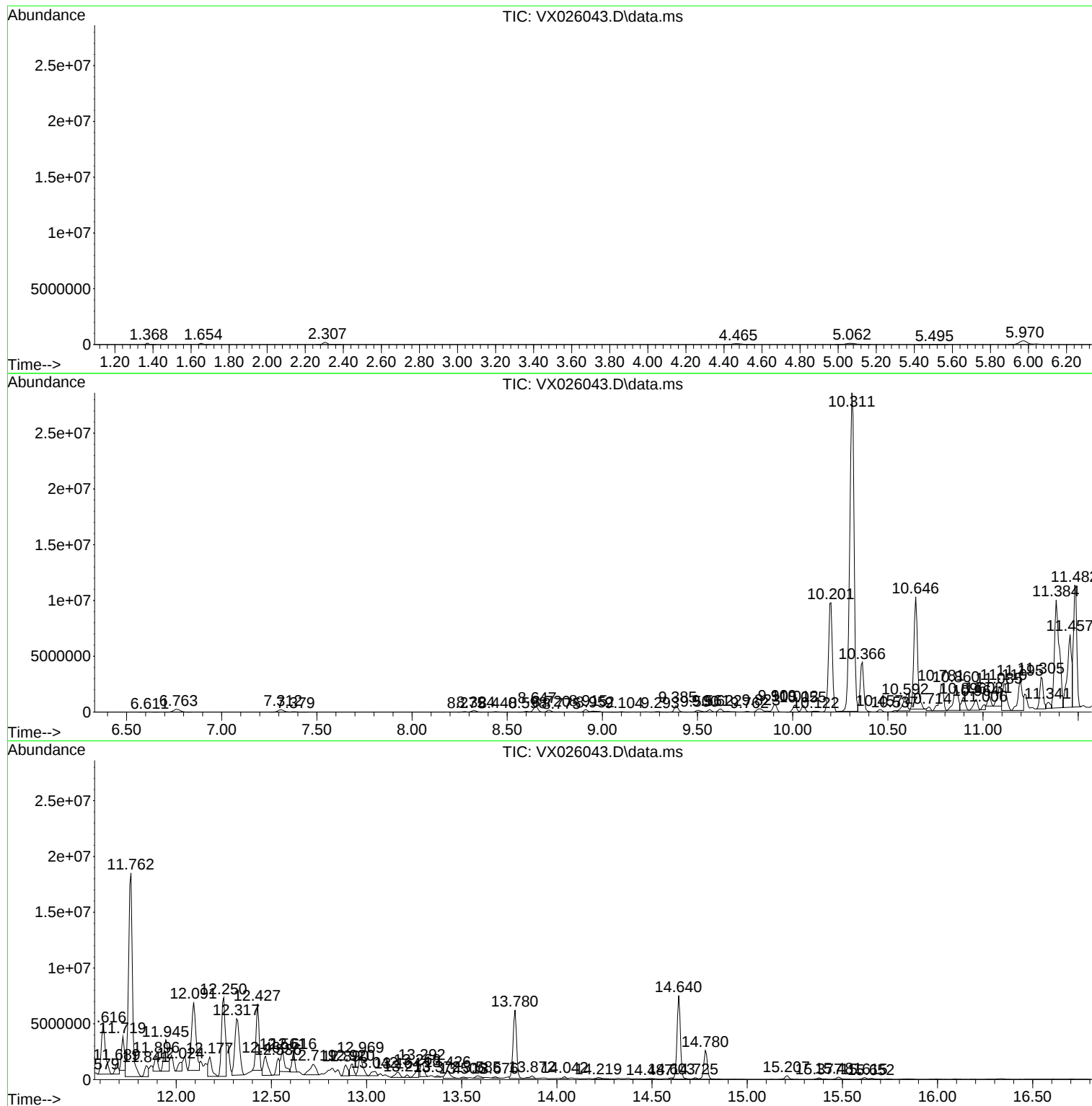
Sum of corrected areas: 286979074

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

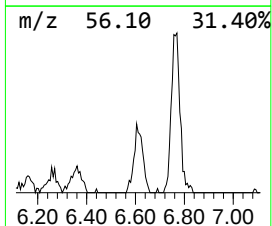
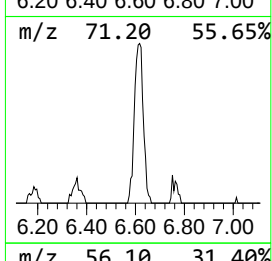
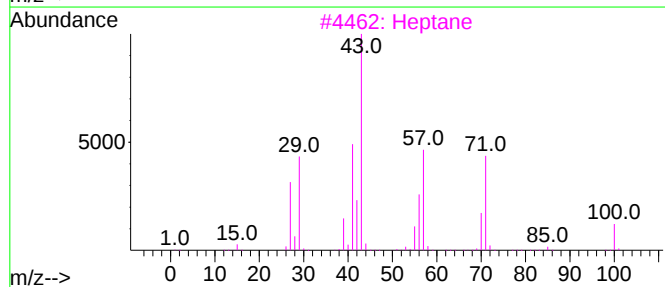
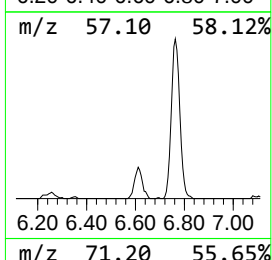
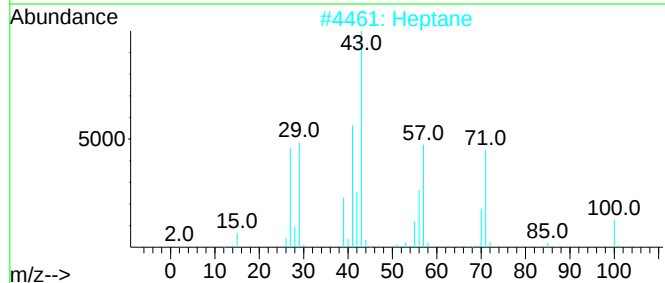
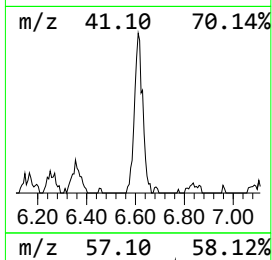
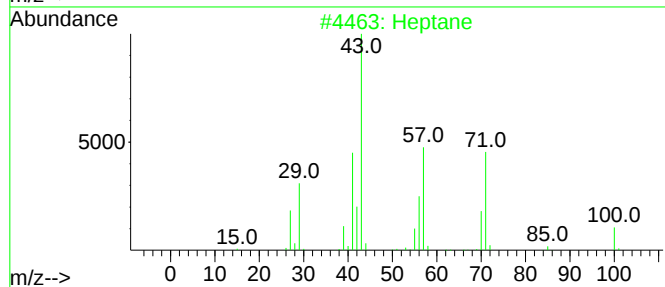
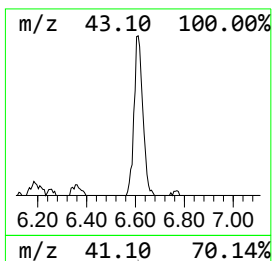
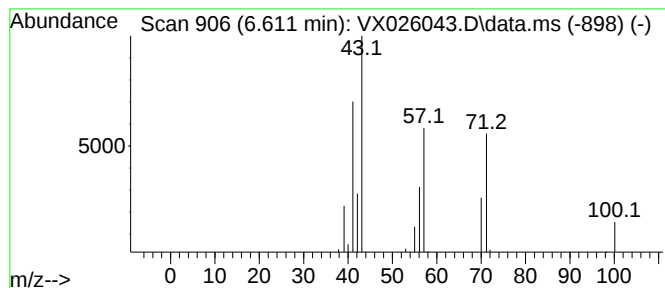
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 70

| R.T.  | EstConc   | Area  | Relative to ISTD    | R.T.  |
|-------|-----------|-------|---------------------|-------|
| 6.611 | 3.06 ug/L | 36845 | 1,4-Difluorobenzene | 6.763 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|----|---|------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 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  | 86   |
| 5    |    |   | Oxalic acid, isobutyl pentyl ester | 216 | C11H20O4 | 1000309-37-0 | 45   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

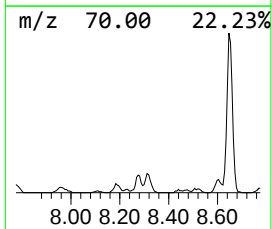
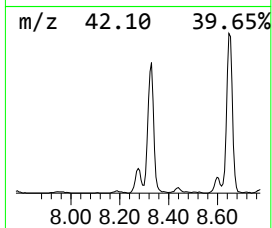
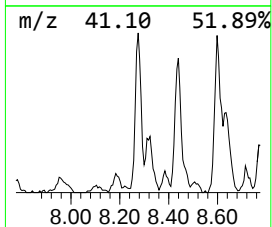
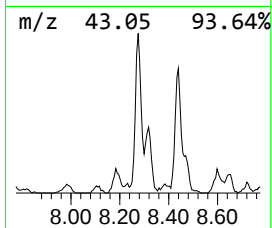
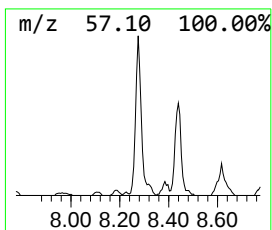
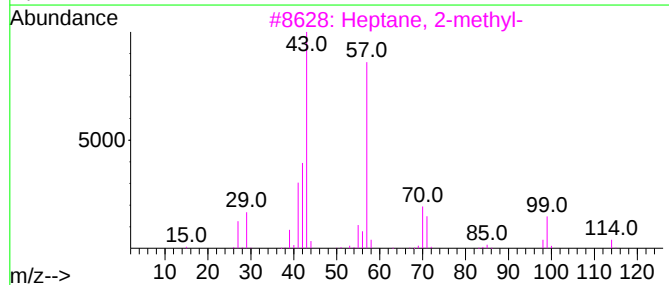
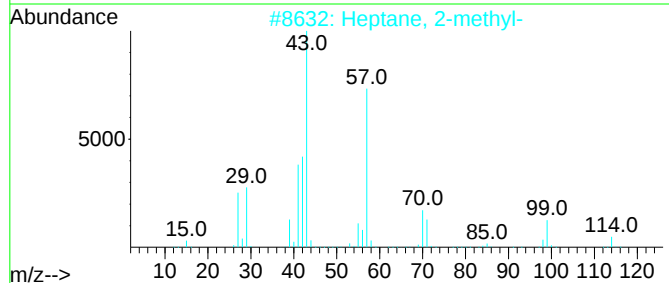
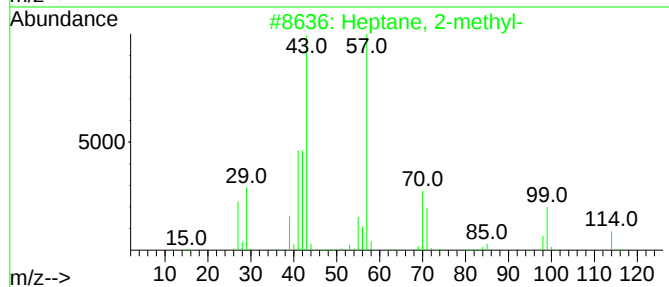
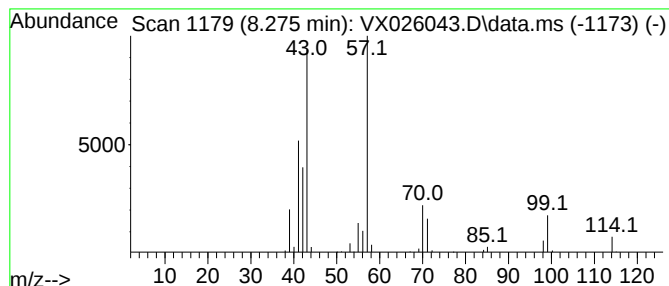
\*\*\*\*\*

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 61

| R.T.  | EstConc   | Area  | Relative to ISTD    | R.T.  |
|-------|-----------|-------|---------------------|-------|
| 8.275 | 5.89 ug/L | 70874 | 1,4-Difluorobenzene | 6.763 |

| Hit# | of 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------|-----|---------|-------------|------|
| 1    |      | Heptane, 2-methyl-    | 114 | C8H18   | 000592-27-8 | 97   |
| 2    |      | Heptane, 2-methyl-    | 114 | C8H18   | 000592-27-8 | 95   |
| 3    |      | Heptane, 2-methyl-    | 114 | C8H18   | 000592-27-8 | 95   |
| 4    |      | Hexane, 2,5-dimethyl- | 114 | C8H18   | 000592-13-2 | 58   |
| 5    |      | Hexane, 2,5-dimethyl- | 114 | C8H18   | 000592-13-2 | 58   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

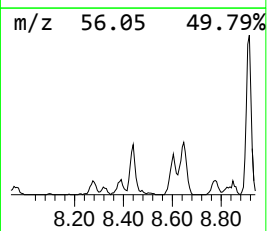
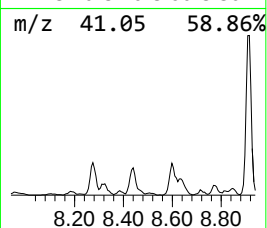
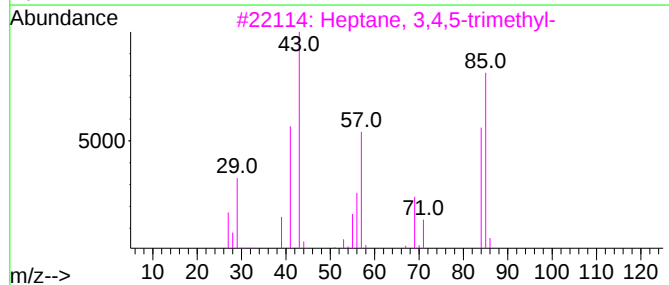
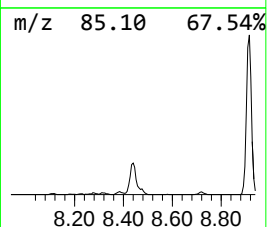
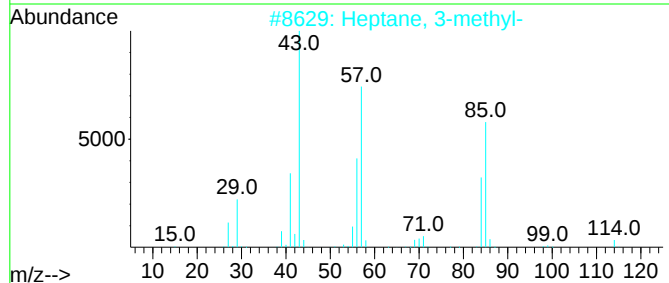
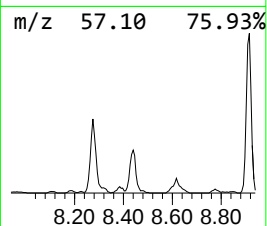
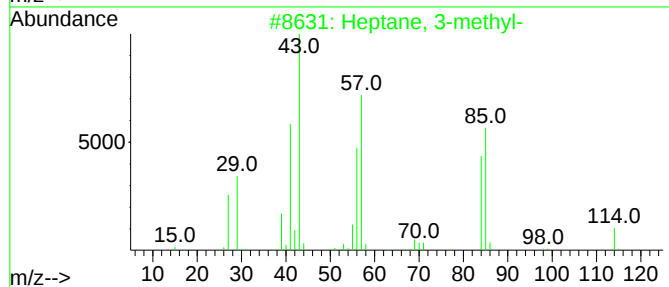
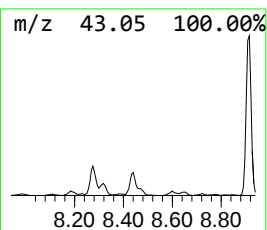
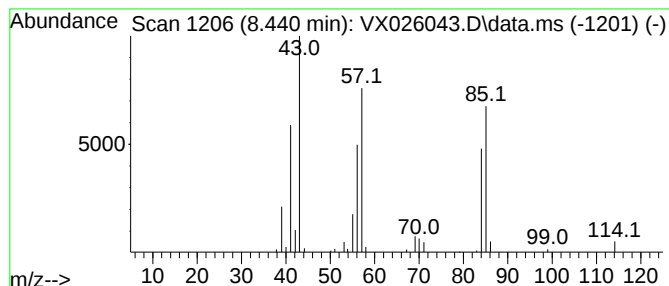
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 67

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.   |
|-------|-----------|-------|------------------|--------|
| 8.440 | 4.24 ug/L | 64255 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | 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 | 59   |
| 4    |    |   | Hexane, 2-methyl-         | 100 | C7H16   | 000591-76-4 | 53   |
| 5    |    |   | Hexane, 3-ethyl-          | 114 | C8H18   | 000619-99-8 | 53   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

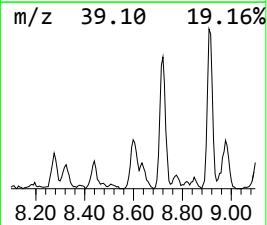
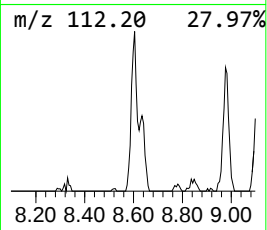
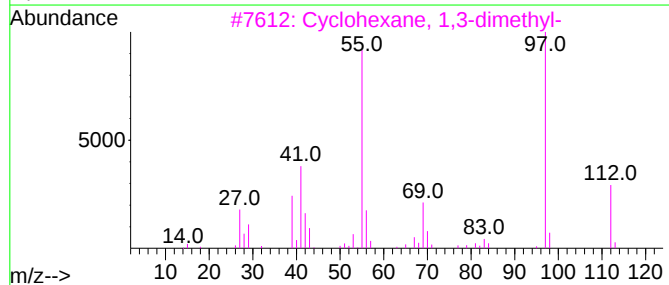
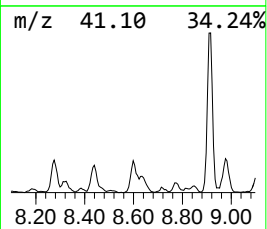
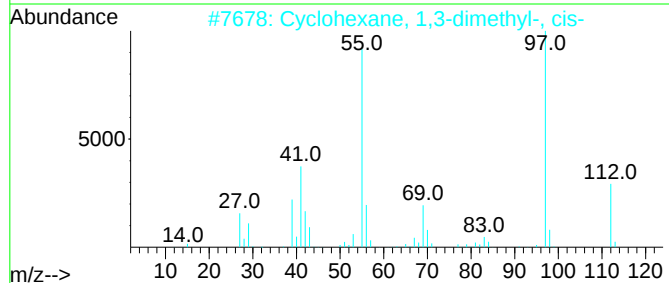
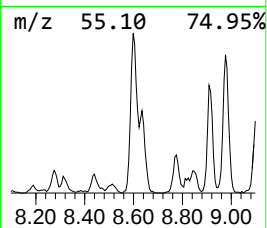
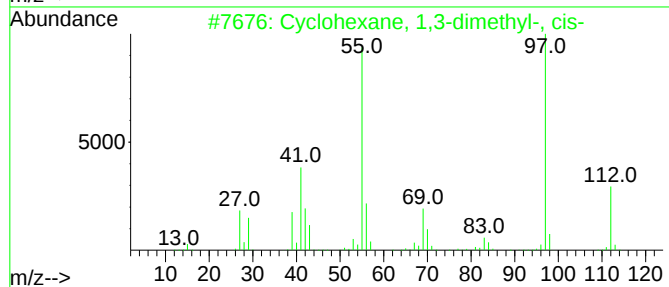
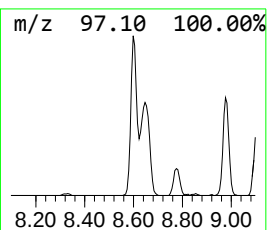
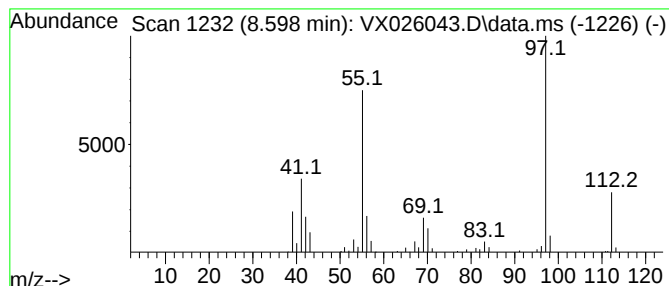
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Cyclic8.598 Concentration Rank 57

| R.T.  | EstConc   | Area   | Relative to ISTD | R.T.   |
|-------|-----------|--------|------------------|--------|
| 8.598 | 6.85 ug/L | 103993 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,3-dimethyl-, cis- | 112 | C8H16   | 000638-04-0 | 97   |
| 2    |    |   | Cyclohexane, 1,3-dimethyl-, cis- | 112 | C8H16   | 000638-04-0 | 95   |
| 3    |    |   | Cyclohexane, 1,3-dimethyl-       | 112 | C8H16   | 000591-21-9 | 95   |
| 4    |    |   | Cyclohexane, 1,3-dimethyl-, cis- | 112 | C8H16   | 000638-04-0 | 94   |
| 5    |    |   | Cyclohexane, 1,3-dimethyl-, cis- | 112 | C8H16   | 000638-04-0 | 94   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

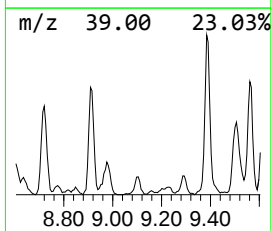
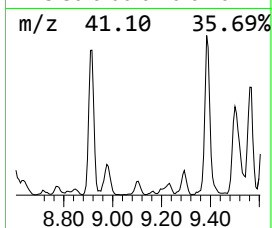
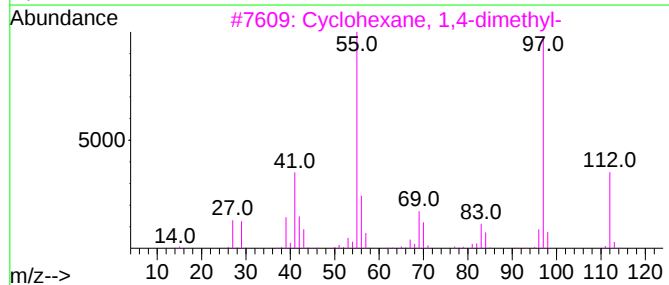
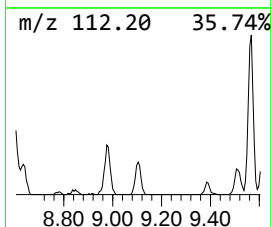
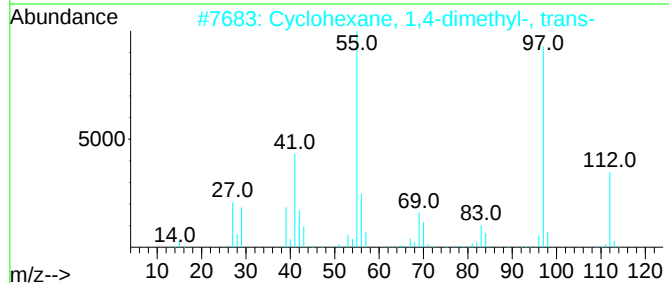
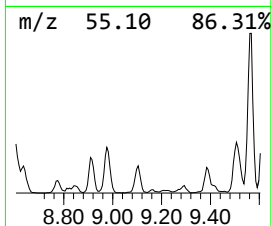
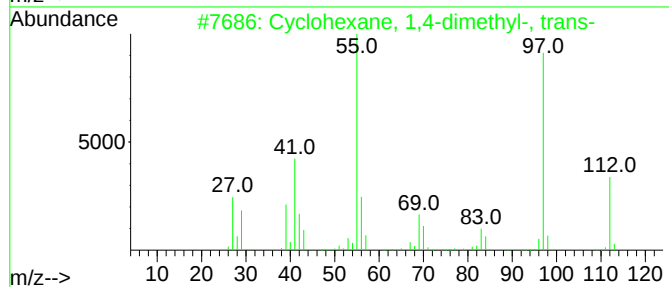
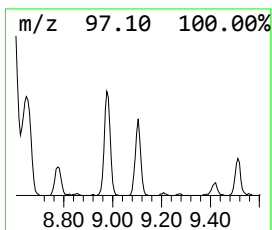
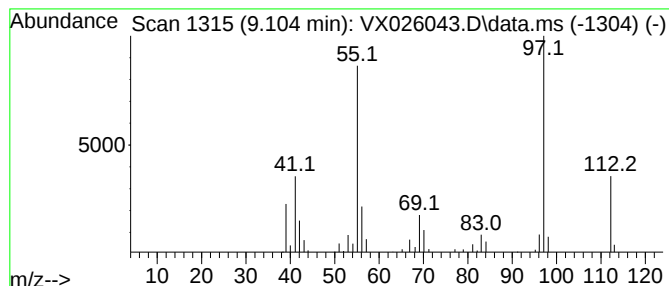
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Cyclic9.104 Concentration Rank 68

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.   |
|-------|-----------|-------|------------------|--------|
| 9.104 | 3.43 ug/L | 52059 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,4-dimethyl-, trans- | 112 | C8H16   | 002207-04-7 | 97   |
| 2    |    |   | Cyclohexane, 1,4-dimethyl-, trans- | 112 | C8H16   | 002207-04-7 | 97   |
| 3    |    |   | Cyclohexane, 1,4-dimethyl-         | 112 | C8H16   | 000589-90-2 | 97   |
| 4    |    |   | Cyclohexane, 1,4-dimethyl-, trans- | 112 | C8H16   | 002207-04-7 | 96   |
| 5    |    |   | Cyclohexane, 1,4-dimethyl-, cis-   | 112 | C8H16   | 000624-29-3 | 96   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

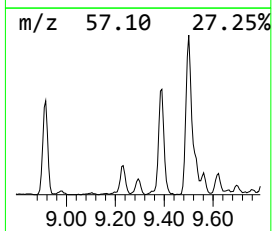
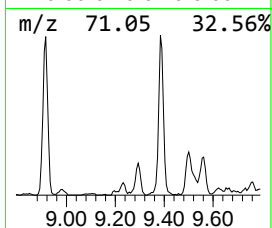
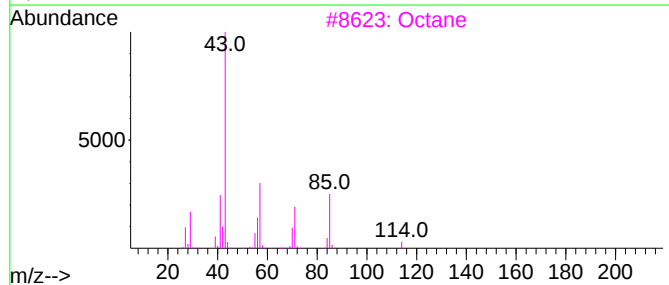
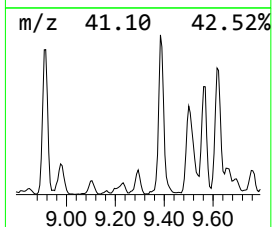
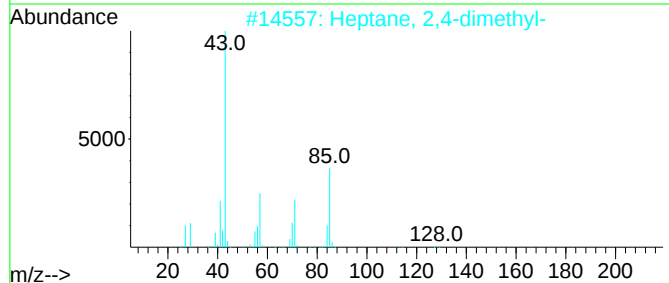
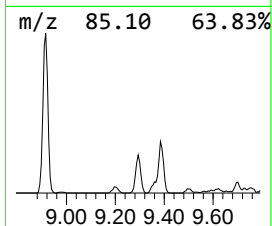
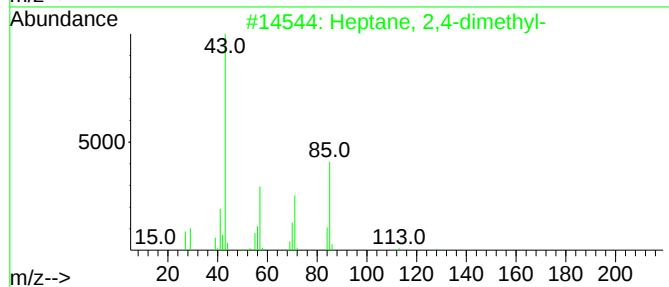
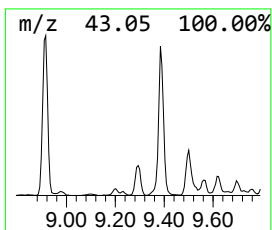
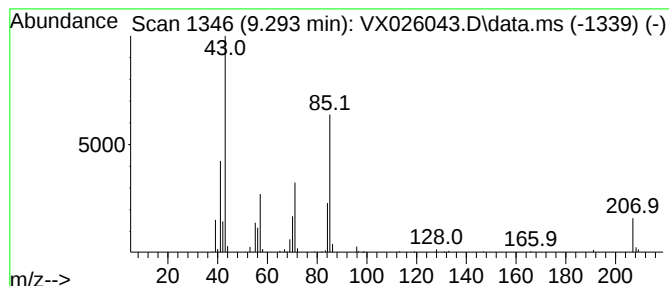
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 65

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.   |
|-------|-----------|-------|------------------|--------|
| 9.293 | 4.84 ug/L | 73488 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm   | CAS#         | Qual |
|------|----|---|------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Heptane, 2,4-dimethyl-             | 128 | C9H20     | 002213-23-2  | 74   |
| 2    |    |   | Heptane, 2,4-dimethyl-             | 128 | C9H20     | 002213-23-2  | 74   |
| 3    |    |   | Octane                             | 114 | C8H18     | 000111-65-9  | 64   |
| 4    |    |   | Sulfurous acid, hexyl pentyl ester | 236 | C11H24O3S | 1000309-14-1 | 59   |
| 5    |    |   | Hexane, 2,3,4-trimethyl-           | 128 | C9H20     | 000921-47-1  | 58   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

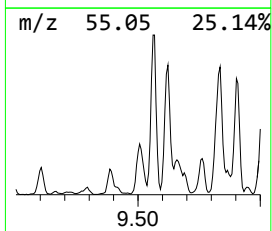
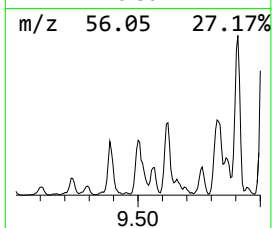
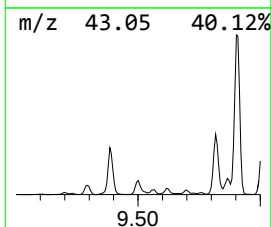
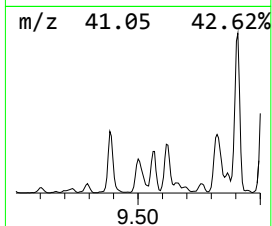
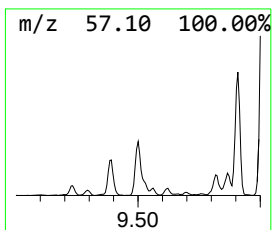
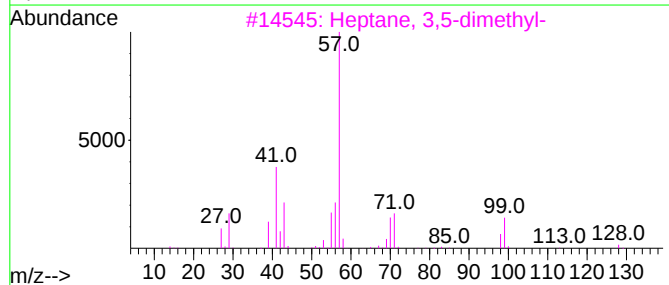
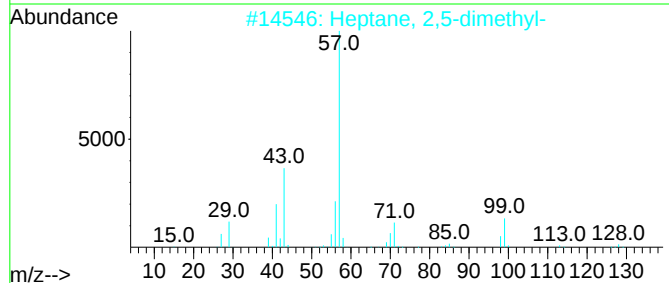
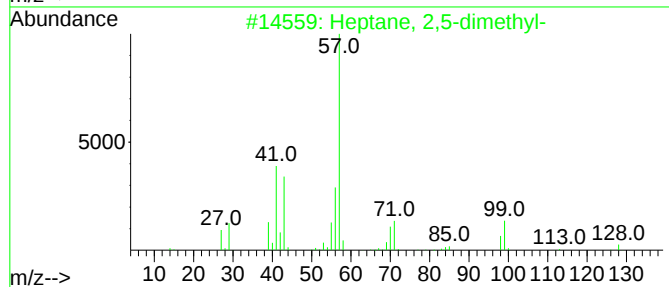
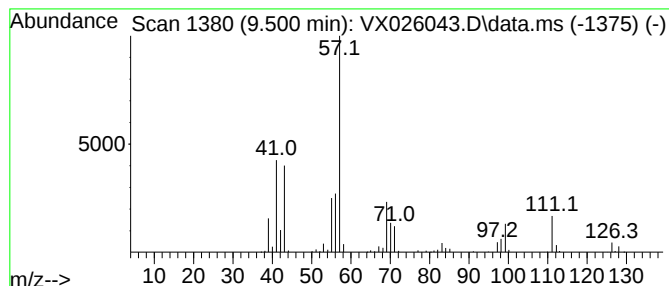
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 42

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.   |
|-------|------------|--------|------------------|--------|
| 9.500 | 19.81 ug/L | 300486 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 76   |
| 2    |    |   | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 64   |
| 3    |    |   | Heptane, 3,5-dimethyl- | 128 | C9H20   | 000926-82-9 | 52   |
| 4    |    |   | Octane, 3-methyl-      | 128 | C9H20   | 002216-33-3 | 52   |
| 5    |    |   | Heptane, 3,5-dimethyl- | 128 | C9H20   | 000926-82-9 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

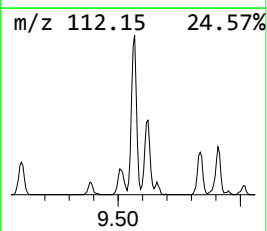
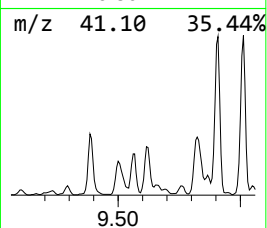
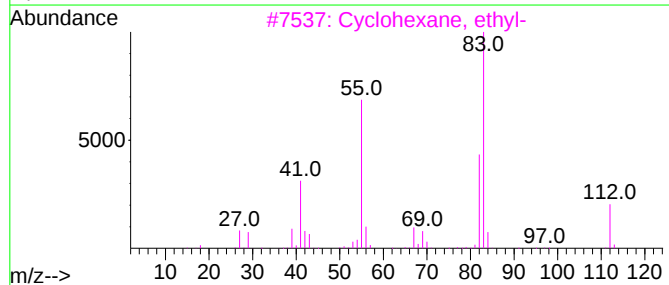
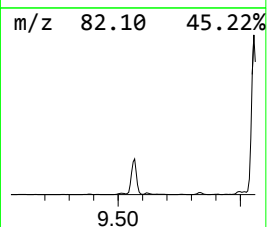
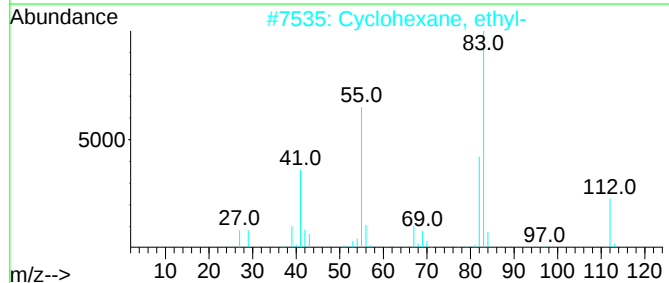
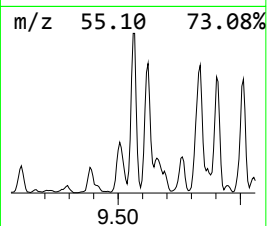
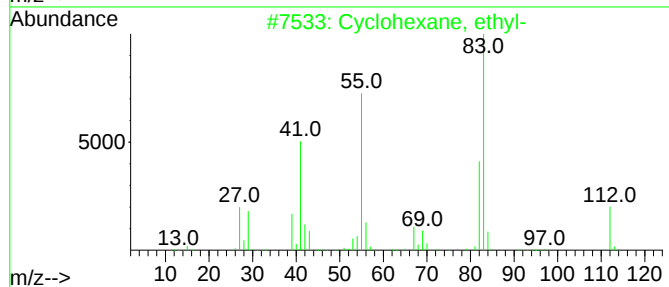
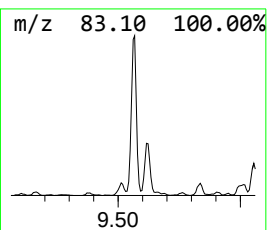
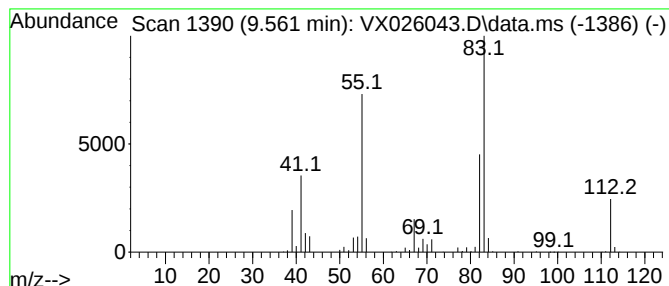
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 (DEL) Alkane: Cyclic9.561 Concentration Rank 43

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.   |
|-------|------------|--------|------------------|--------|
| 9.561 | 18.54 ug/L | 281292 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Cyclohexane, ethyl-                 | 112 | C8H16    | 001678-91-7  | 87   |
| 2    |    |   | Cyclohexane, ethyl-                 | 112 | C8H16    | 001678-91-7  | 87   |
| 3    |    |   | Cyclohexane, ethyl-                 | 112 | C8H16    | 001678-91-7  | 87   |
| 4    |    |   | Cyclohexane, ethyl-                 | 112 | C8H16    | 001678-91-7  | 58   |
| 5    |    |   | Oxalic acid, cyclohexyl dodecyl ... | 340 | C20H36O4 | 1000309-31-4 | 53   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

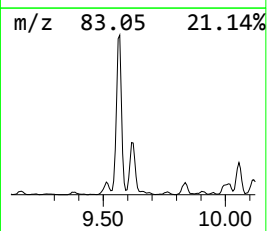
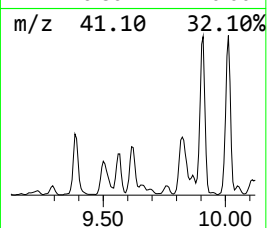
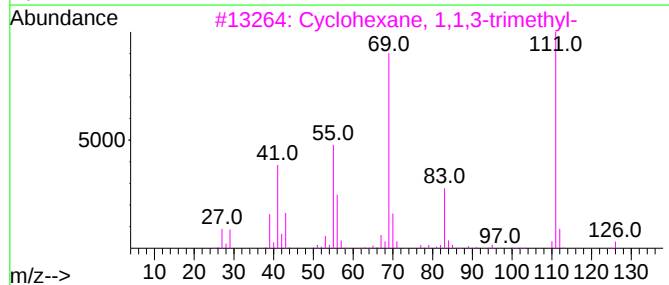
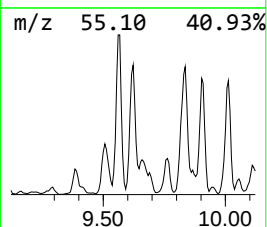
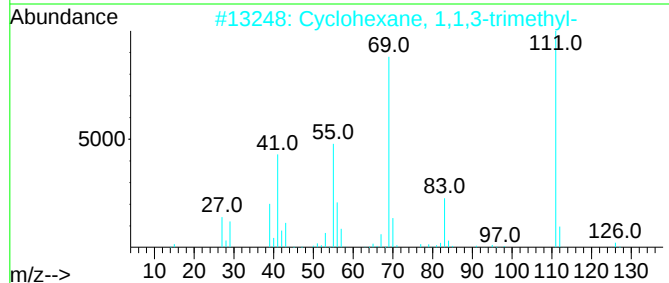
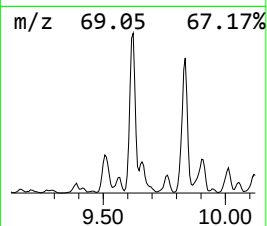
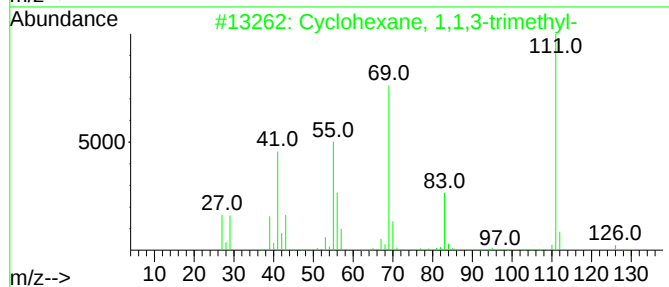
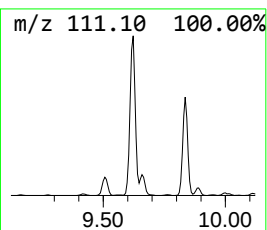
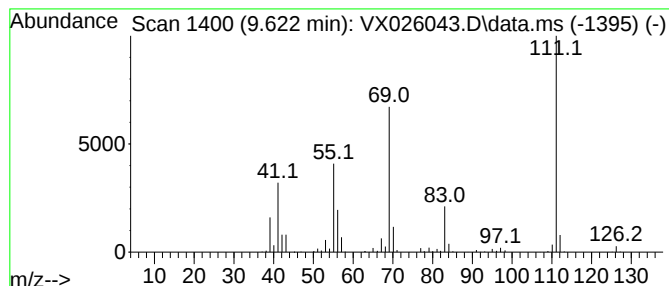
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Cyclic9.622 Concentration Rank 39

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.   |
|-------|------------|--------|------------------|--------|
| 9.622 | 26.74 ug/L | 405676 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | 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, 1,1,3-trimethyl- | 126 | C9H18   | 003073-66-3 | 92   |
| 5    |    |   | 1,1,4-Trimethylcyclohexane    | 126 | C9H18   | 007094-27-1 | 72   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

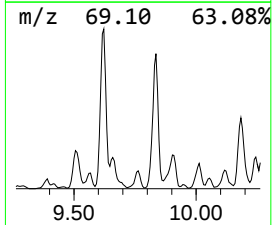
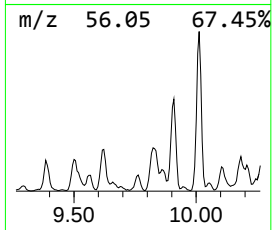
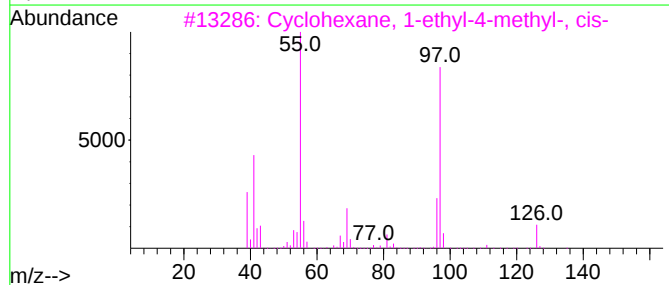
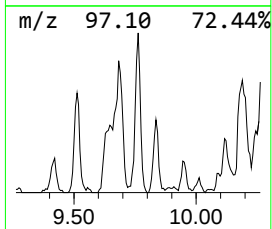
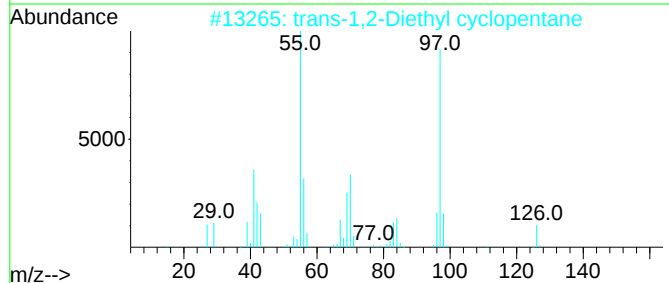
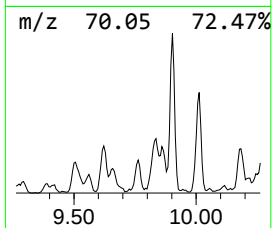
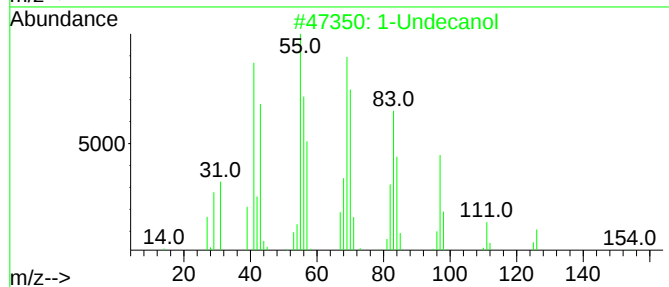
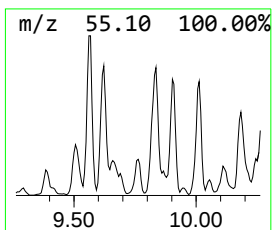
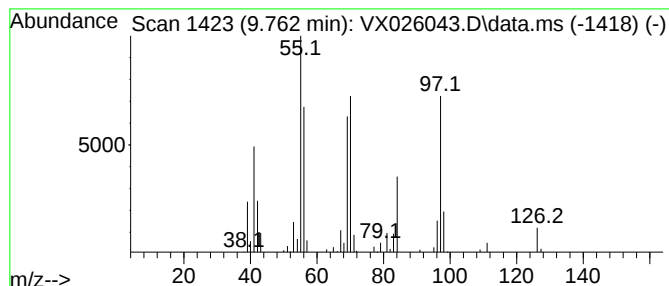
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 1-Undecanol Concentration Rank 64

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.   |
|-------|-----------|-------|------------------|--------|
| 9.762 | 4.90 ug/L | 74360 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Undecanol                         | 172 | C11H24O | 000112-42-5 | 72   |
| 2    |    |   | trans-1,2-Diethyl cyclopentane      | 126 | C9H18   | 000932-40-1 | 62   |
| 3    |    |   | Cyclohexane, 1-ethyl-4-methyl-, ... | 126 | C9H18   | 004926-78-7 | 30   |
| 4    |    |   | Cyclohexane, 1-ethyl-2-methyl-      | 126 | C9H18   | 003728-54-9 | 27   |
| 5    |    |   | trans--4-Nonene                     | 126 | C9H18   | 010405-85-3 | 27   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

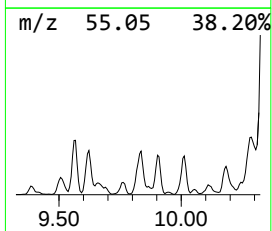
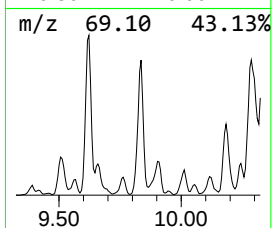
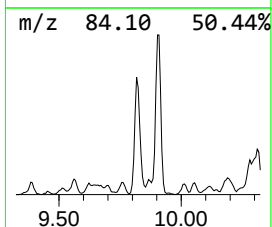
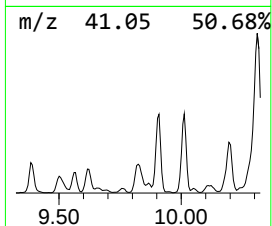
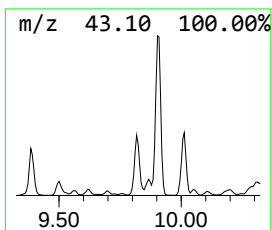
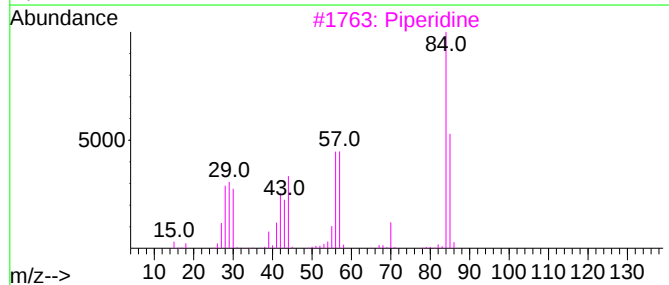
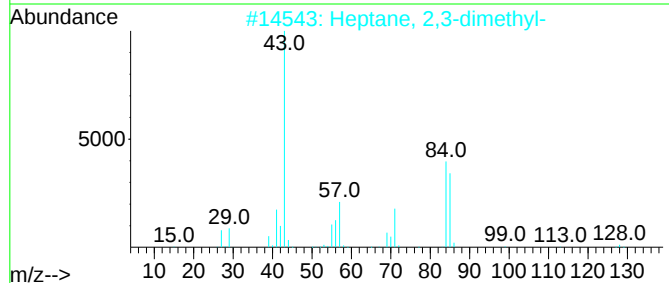
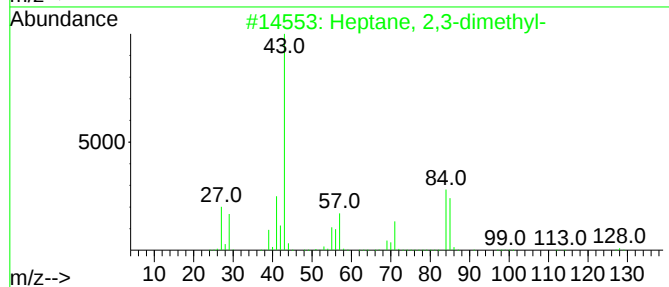
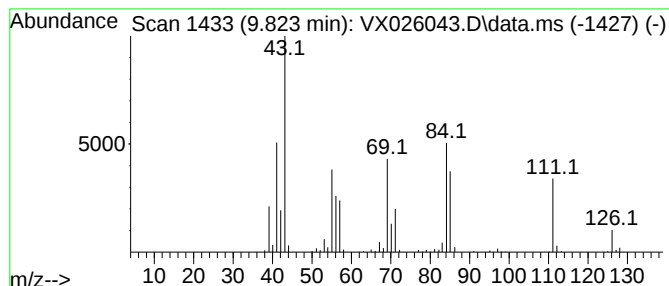
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 35

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.   |
|-------|------------|--------|------------------|--------|
| 9.823 | 40.48 ug/L | 614082 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptane, 2,3-dimethyl-    | 128 | C9H20   | 003074-71-3 | 50   |
| 2    |    |   | Heptane, 2,3-dimethyl-    | 128 | C9H20   | 003074-71-3 | 49   |
| 3    |    |   | Piperidine                | 85  | C5H11N  | 000110-89-4 | 46   |
| 4    |    |   | Nonane, 4-ethyl-5-methyl- | 170 | C12H26  | 001632-71-9 | 45   |
| 5    |    |   | Hexanal, 3,3-dimethyl-    | 128 | C8H16O  | 055320-57-5 | 43   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

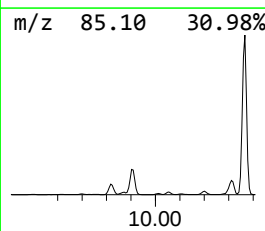
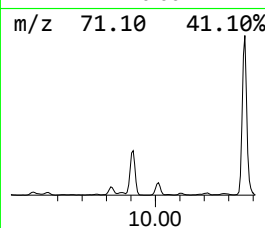
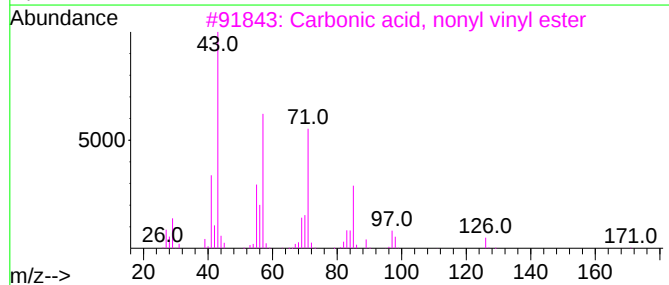
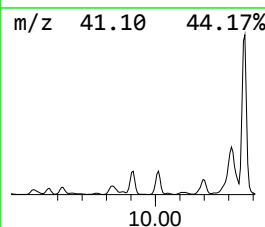
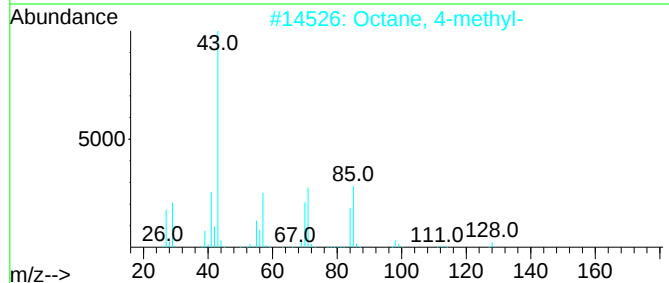
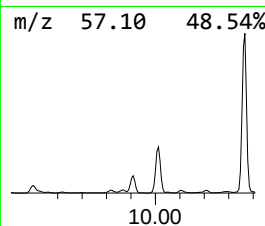
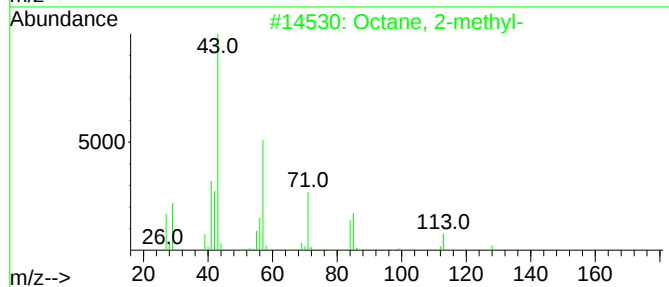
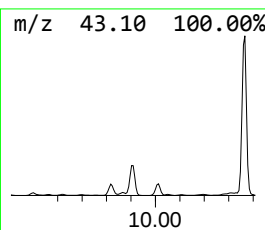
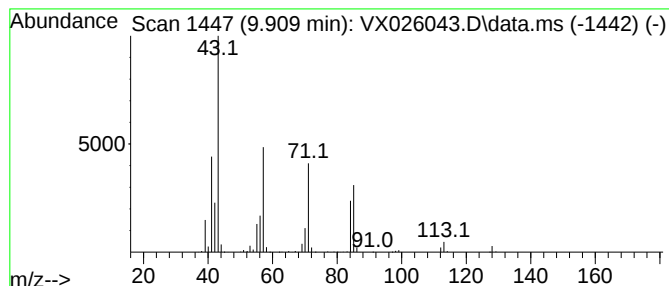
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.   |
|-------|------------|--------|------------------|--------|
| 9.909 | 65.20 ug/L | 989047 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm  | CAS#         | Qual |
|------|----|---|----------------------------------|-----|----------|--------------|------|
| 1    |    |   | Octane, 2-methyl-                | 128 | C9H20    | 003221-61-2  | 90   |
| 2    |    |   | Octane, 4-methyl-                | 128 | C9H20    | 002216-34-4  | 72   |
| 3    |    |   | Carbonic acid, nonyl vinyl ester | 214 | C12H22O3 | 1000383-25-6 | 64   |
| 4    |    |   | Dodecane, 5-methyl-              | 184 | C13H28   | 017453-93-9  | 59   |
| 5    |    |   | Heptane, 2,6-dimethyl-           | 128 | C9H20    | 001072-05-5  | 58   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

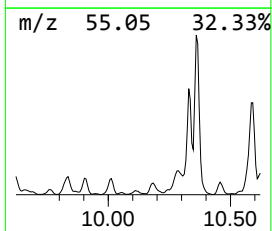
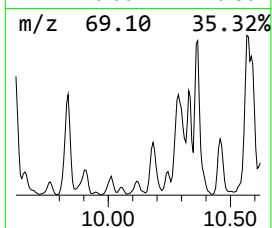
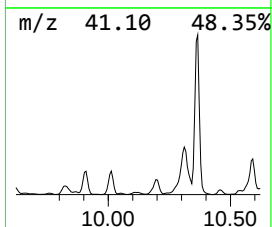
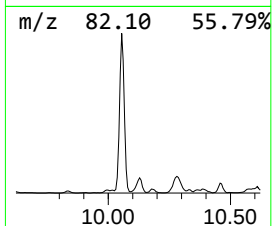
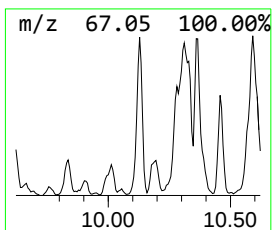
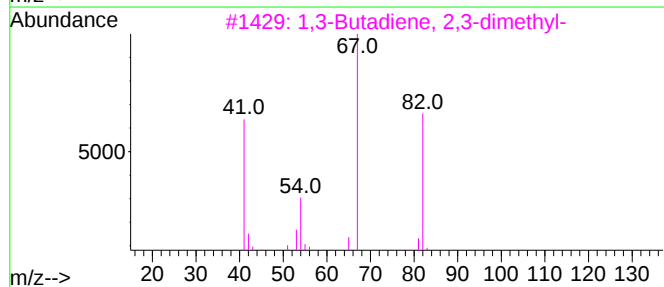
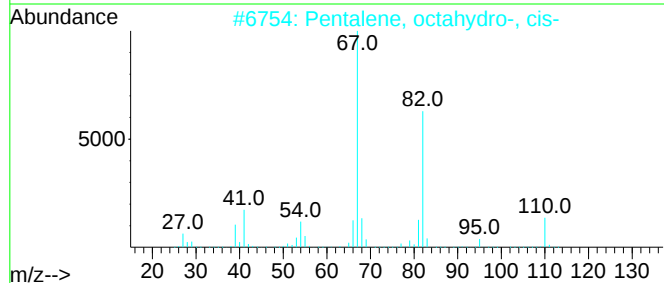
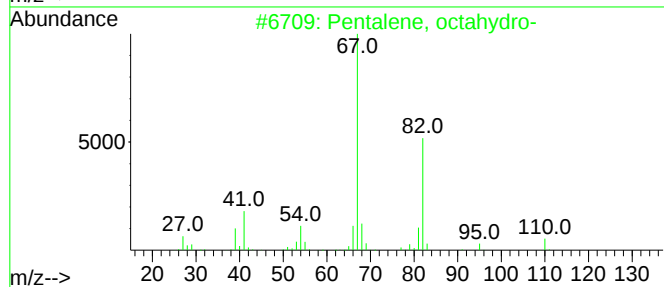
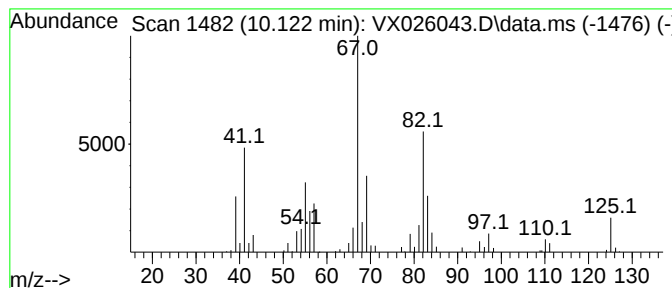
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Pentalene, octahydro- Concentration Rank 51

| R.T.   | EstConc   | Area   | Relative to ISTD | R.T.   |
|--------|-----------|--------|------------------|--------|
| 10.122 | 9.92 ug/L | 150494 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentalene, octahydro-        | 110 | C8H14   | 000694-72-4 | 58   |
| 2    |    |   | Pentalene, octahydro-, cis-  | 110 | C8H14   | 001755-05-1 | 50   |
| 3    |    |   | 1,3-Butadiene, 2,3-dimethyl- | 82  | C6H10   | 000513-81-5 | 50   |
| 4    |    |   | 4-Methyl-1,3-pentadiene      | 82  | C6H10   | 000926-56-7 | 49   |
| 5    |    |   | 4-Hexen-1-ol, (E)-           | 100 | C6H12O  | 000928-92-7 | 47   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

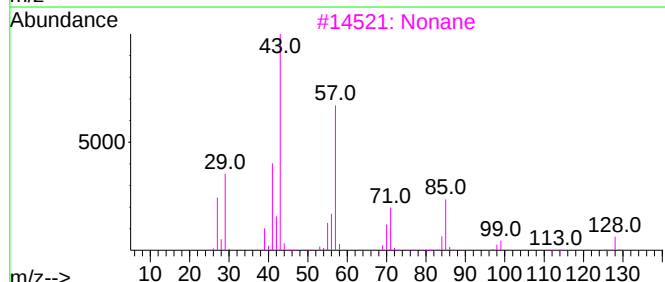
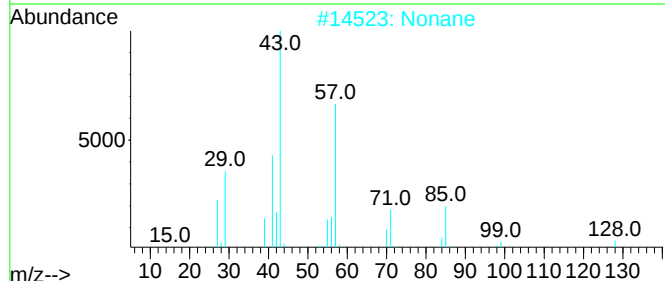
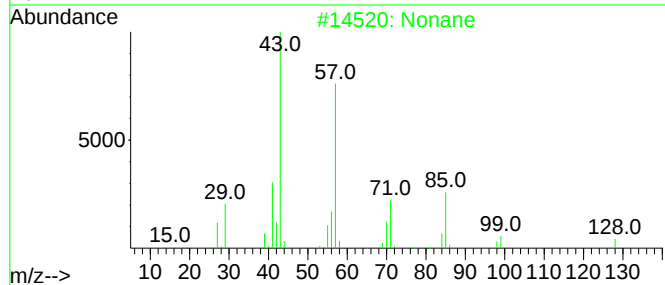
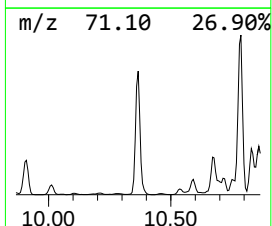
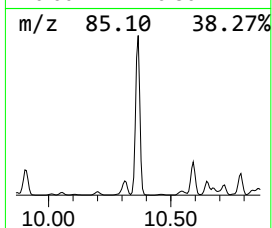
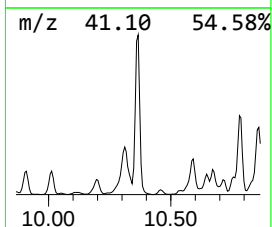
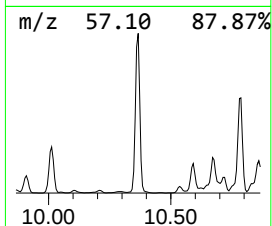
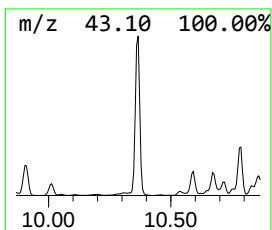
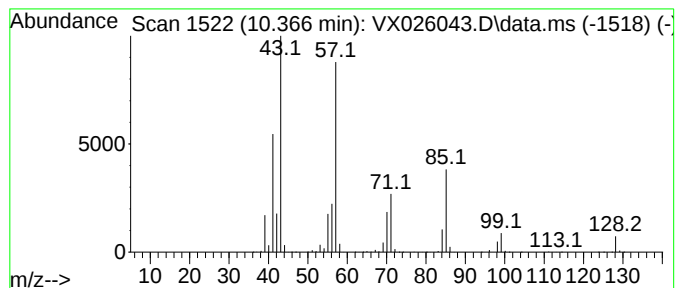
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 10.366 | 384.54 ug/L | 5833720 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Nonane                              | 128 | C9H20    | 000111-84-2  | 94   |
| 2    |    |   | Nonane                              | 128 | C9H20    | 000111-84-2  | 94   |
| 3    |    |   | Nonane                              | 128 | C9H20    | 000111-84-2  | 91   |
| 4    |    |   | Nonane                              | 128 | C9H20    | 000111-84-2  | 87   |
| 5    |    |   | Carbonic acid, decyl prop-1-en-2... | 242 | C14H26O3 | 1000382-90-5 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

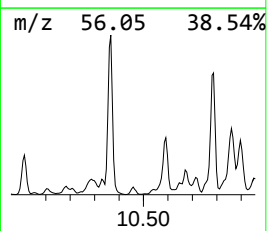
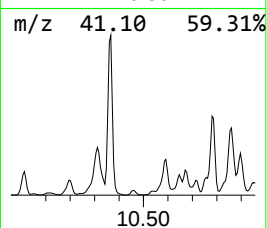
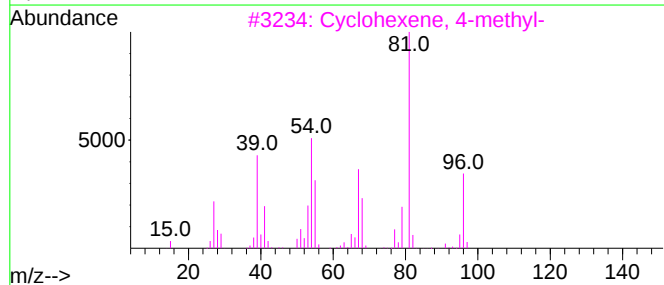
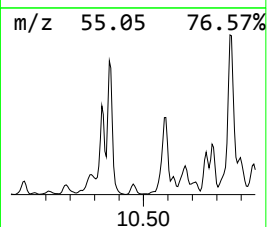
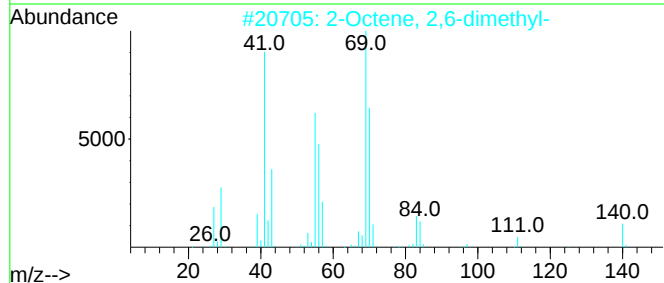
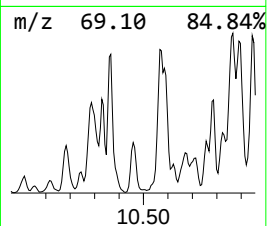
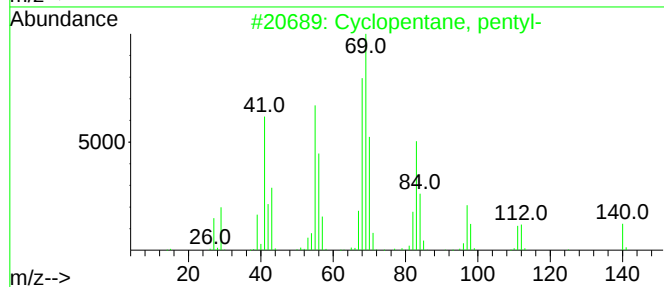
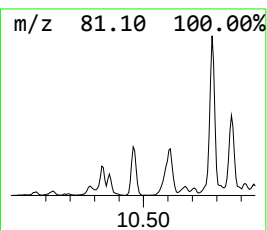
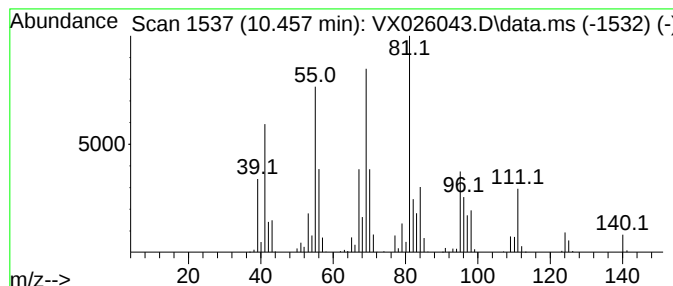
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Cyclic10.458 Concentration Rank 41

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.458 | 21.74 ug/L | 329780 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclopentane, pentyl-     | 140 | C10H20  | 003741-00-2 | 43   |
| 2    |    |   | 2-Octene, 2,6-dimethyl-   | 140 | C10H20  | 004057-42-5 | 38   |
| 3    |    |   | Cyclohexene, 4-methyl-    | 96  | C7H12   | 000591-47-9 | 35   |
| 4    |    |   | 3-Hexene, 2-methyl-, (E)- | 98  | C7H14   | 000692-24-0 | 30   |
| 5    |    |   | 3-Hexene, 2-methyl-, (E)- | 98  | C7H14   | 000692-24-0 | 30   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

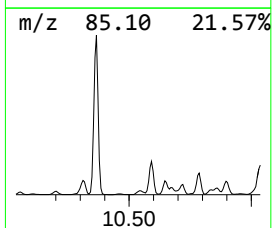
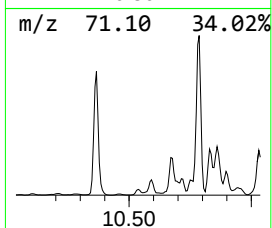
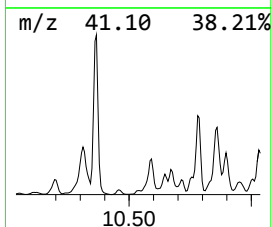
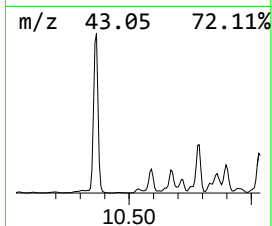
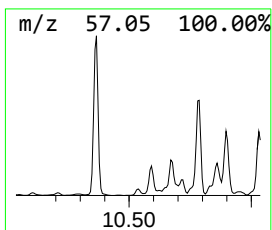
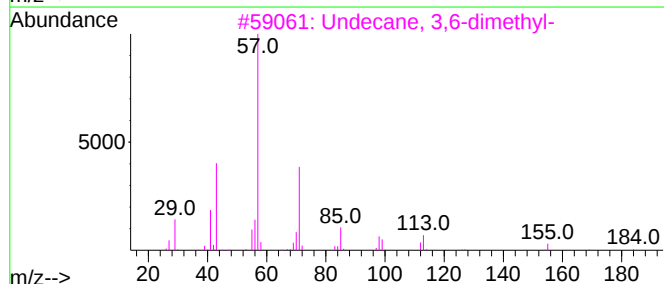
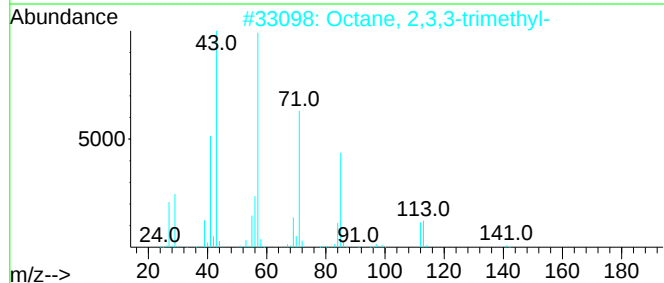
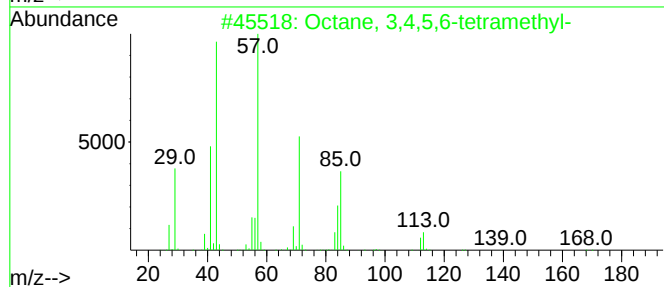
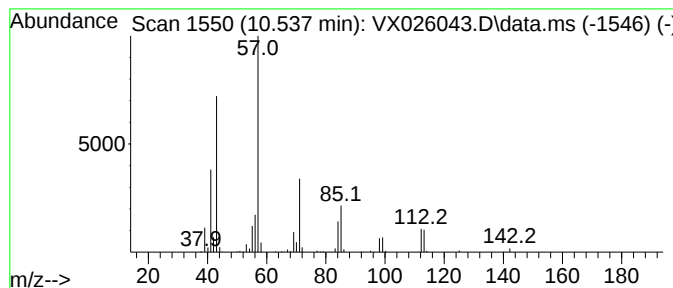
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 49

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.537 | 11.50 ug/L | 174409 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 3,4,5,6-tetramethyl- | 170 | C12H26  | 062185-21-1 | 64   |
| 2    |    |   | Octane, 2,3,3-trimethyl-     | 156 | C11H24  | 062016-30-2 | 64   |
| 3    |    |   | Undecane, 3,6-dimethyl-      | 184 | C13H28  | 017301-28-9 | 59   |
| 4    |    |   | Octane, 3,5-dimethyl-        | 142 | C10H22  | 015869-93-9 | 52   |
| 5    |    |   | Undecane, 5-methyl-          | 170 | C12H26  | 001632-70-8 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

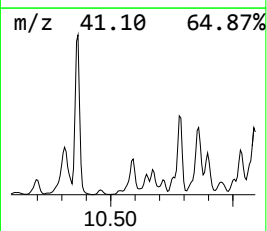
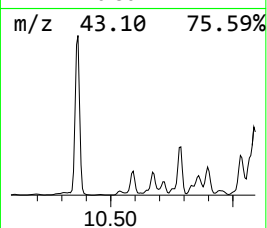
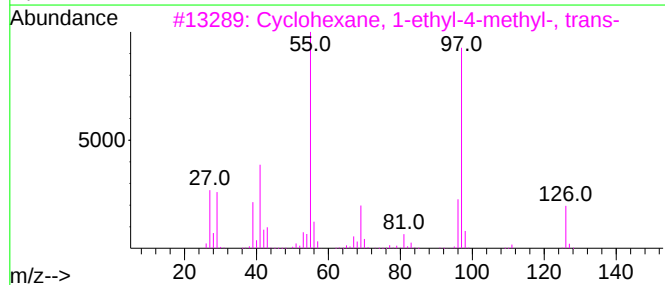
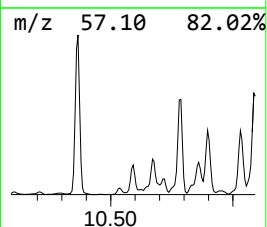
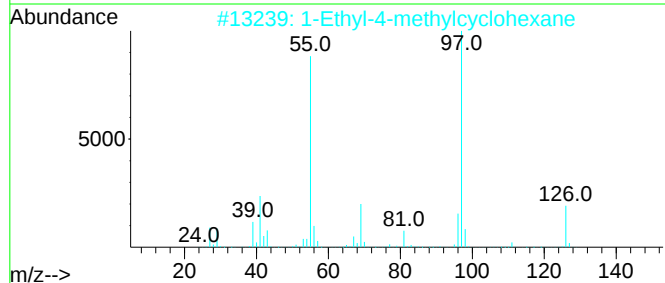
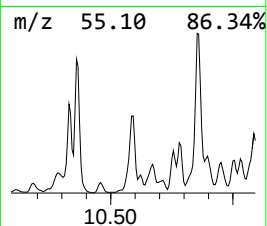
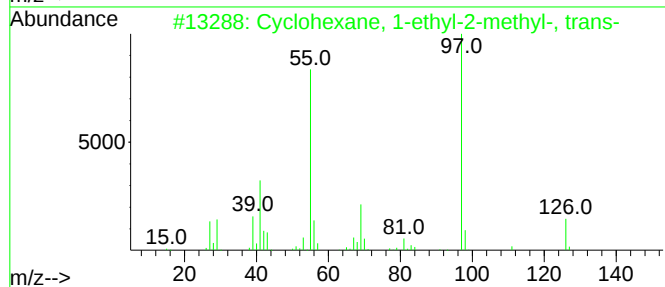
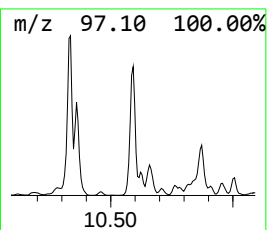
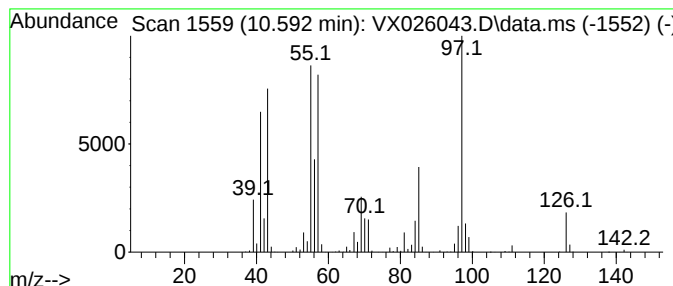
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 (DEL) Alkane: Cyclic10.592 Concentration Rank 15

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 10.592 | 125.27 ug/L | 1900370 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1-ethyl-2-methyl-, ... | 126 | C9H18   | 004923-78-8 | 55   |
| 2    |    |   | 1-Ethyl-4-methylcyclohexane         | 126 | C9H18   | 003728-56-1 | 49   |
| 3    |    |   | Cyclohexane, 1-ethyl-4-methyl-, ... | 126 | C9H18   | 006236-88-0 | 49   |
| 4    |    |   | Cyclohexane, 1-ethyl-2-methyl-      | 126 | C9H18   | 003728-54-9 | 46   |
| 5    |    |   | 1-Ethyl-3-methylcyclohexane (c,t)   | 126 | C9H18   | 003728-55-0 | 46   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

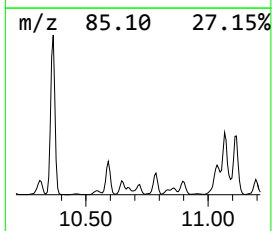
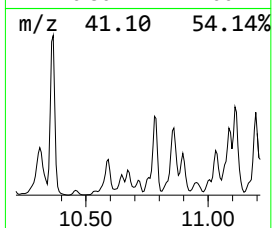
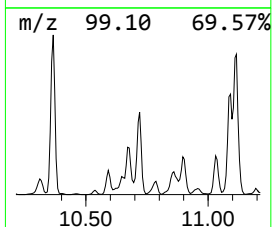
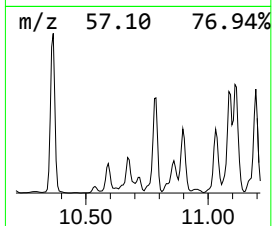
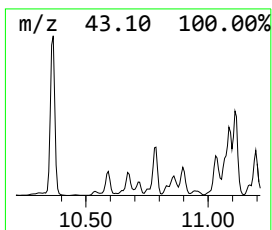
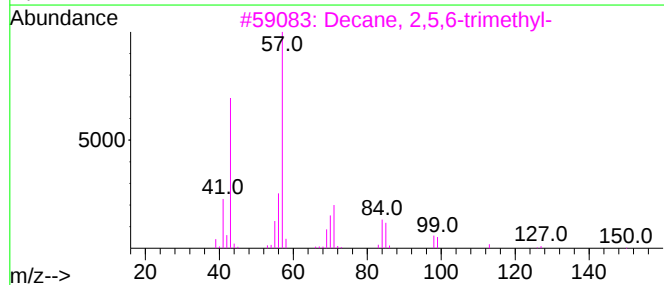
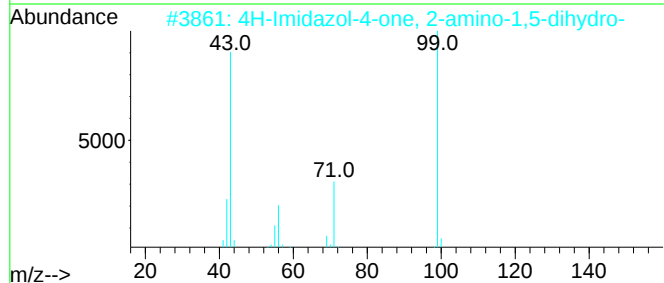
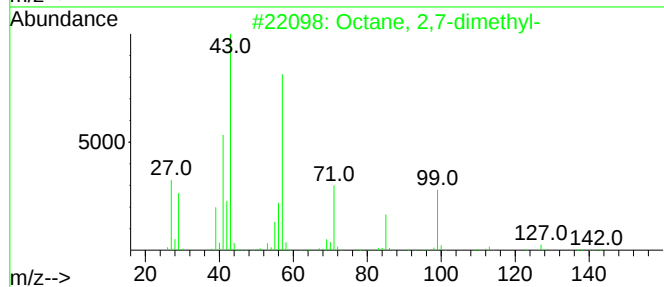
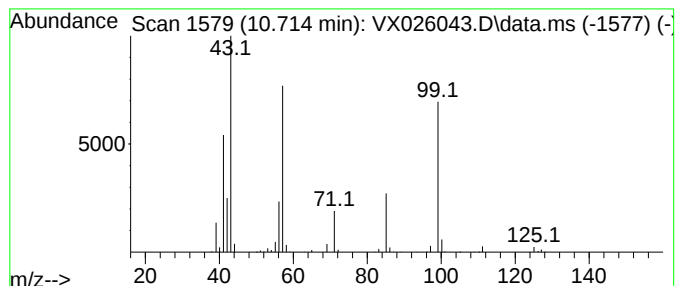
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 40

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.714 | 24.39 ug/L | 369966 | Chlorobenzene-d5 | 10.055 |

| Hit# | of                                  | 5 | Tentative ID | MW      | MolForm | CAS#        | Qual |
|------|-------------------------------------|---|--------------|---------|---------|-------------|------|
| 1    | Octane, 2,7-dimethyl-               |   | 142          | C10H22  |         | 001072-16-8 | 56   |
| 2    | 4H-Imidazol-4-one, 2-amino-1,5-d... |   | 99           | C3H5N3O |         | 000503-86-6 | 53   |
| 3    | Decane, 2,5,6-trimethyl-            |   | 184          | C13H28  |         | 062108-23-0 | 50   |
| 4    | Nonane                              |   | 128          | C9H20   |         | 000111-84-2 | 50   |
| 5    | 3,6-Heptanedione                    |   | 128          | C7H12O2 |         | 001703-51-1 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

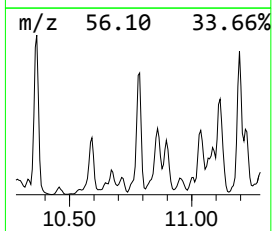
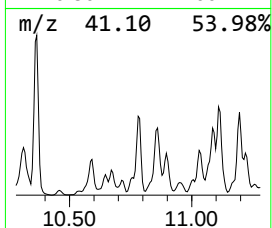
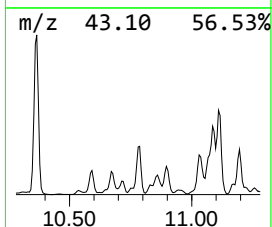
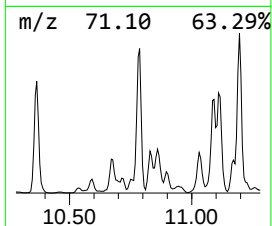
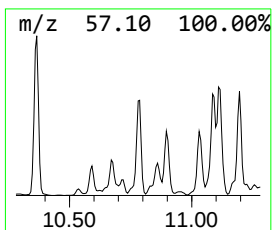
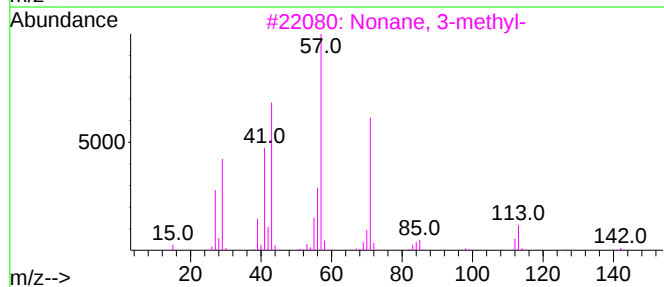
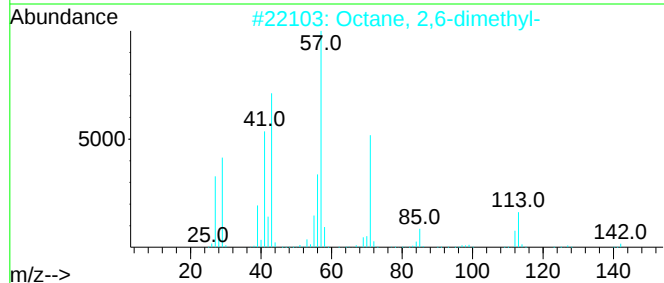
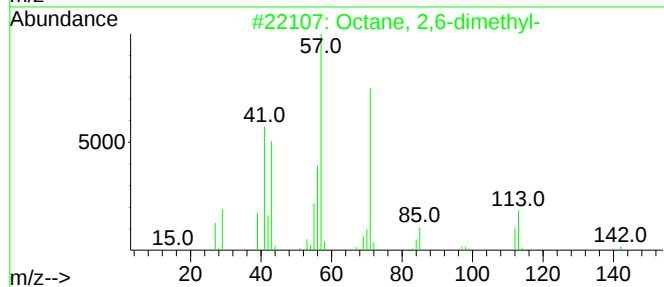
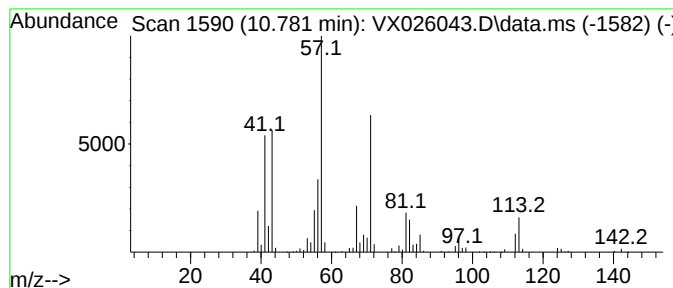
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 10.781 | 244.23 ug/L | 3705170 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 94   |
| 2    |    |   | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 81   |
| 3    |    |   | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 81   |
| 4    |    |   | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 76   |
| 5    |    |   | Octane, 3,6-dimethyl- | 142 | C10H22  | 015869-94-0 | 74   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

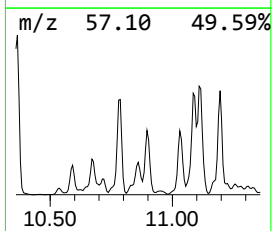
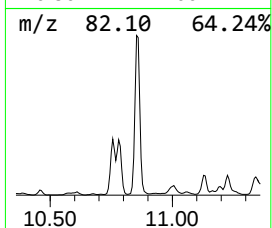
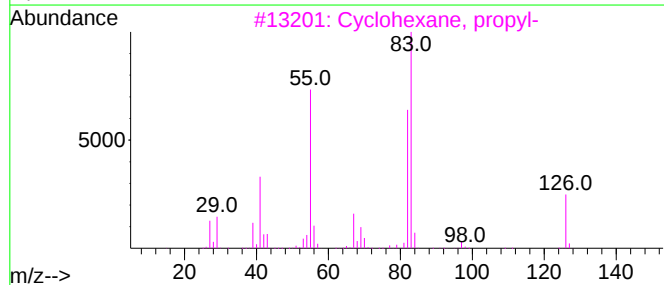
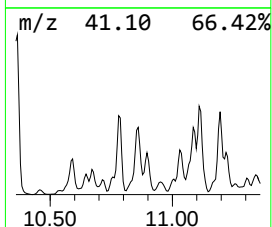
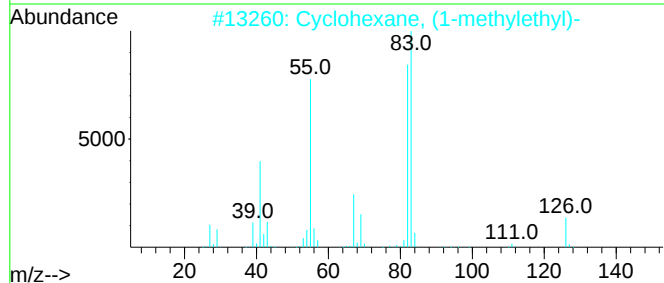
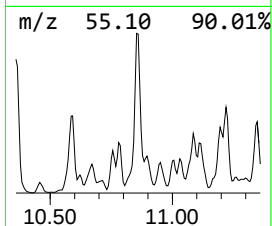
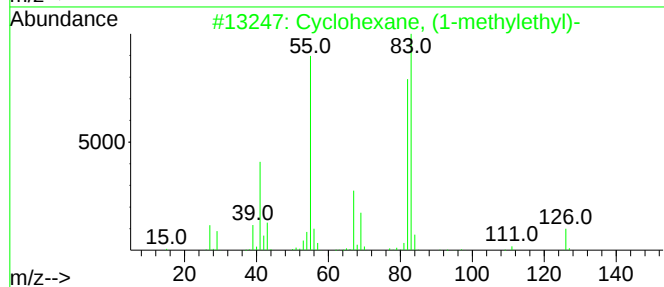
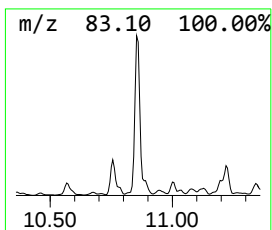
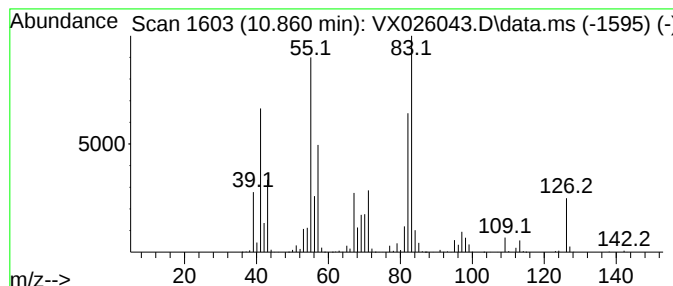
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 (DEL) Alkane: Cyclic10.860 Concentration Rank 9

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 10.860 | 251.41 ug/L | 3814070 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 81   |
| 2    |    |   | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 81   |
| 3    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 64   |
| 4    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 64   |
| 5    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

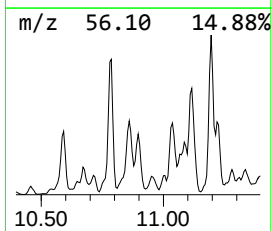
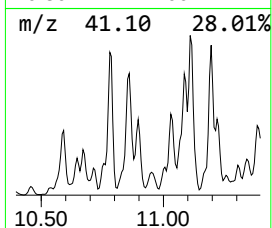
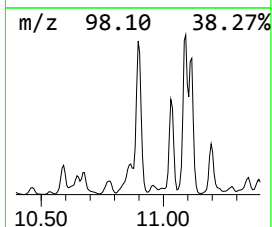
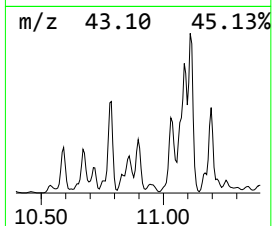
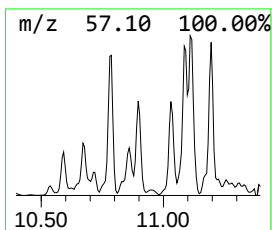
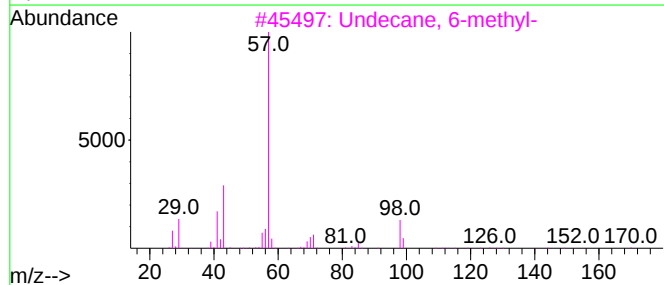
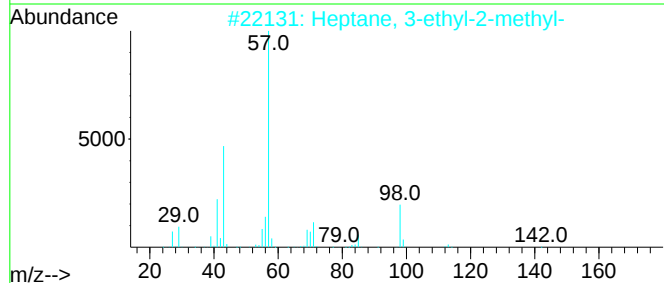
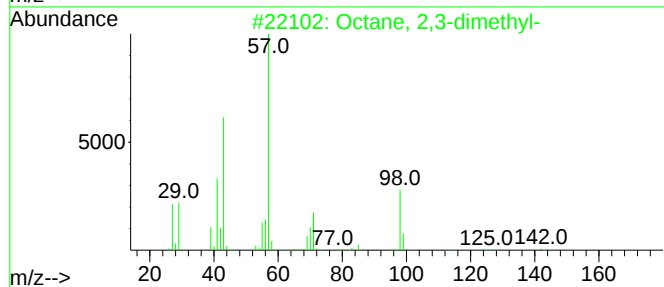
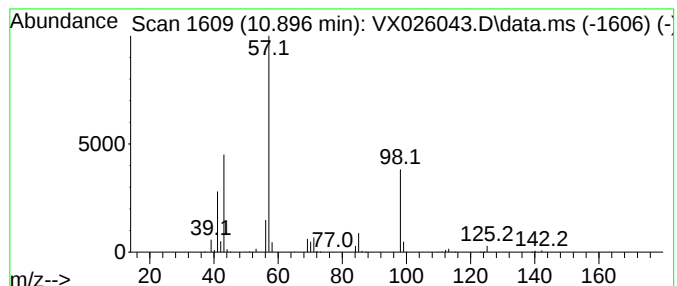
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 20

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 10.896 | 103.28 ug/L | 1566760 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2,3-dimethyl-      | 142 | C10H22  | 007146-60-3 | 64   |
| 2    |    |   | Heptane, 3-ethyl-2-methyl- | 142 | C10H22  | 014676-29-0 | 64   |
| 3    |    |   | Undecane, 6-methyl-        | 170 | C12H26  | 017302-33-9 | 59   |
| 4    |    |   | Heptane, 4-propyl-         | 142 | C10H22  | 003178-29-8 | 56   |
| 5    |    |   | Heptane, 4-propyl-         | 142 | C10H22  | 003178-29-8 | 56   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

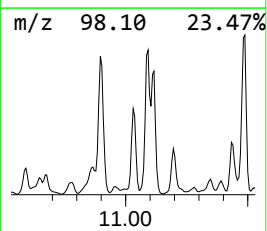
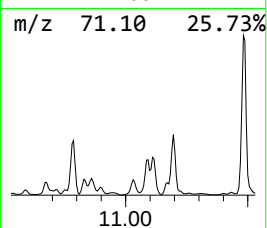
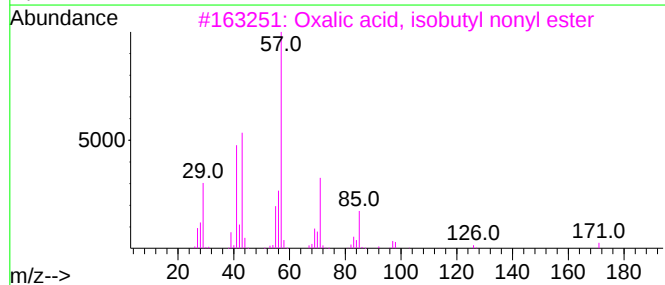
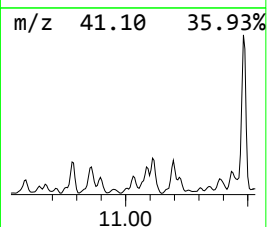
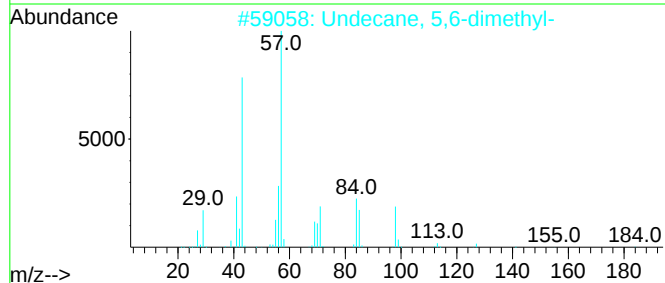
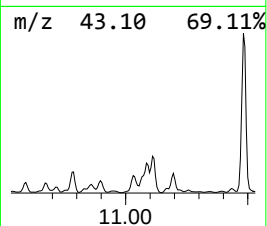
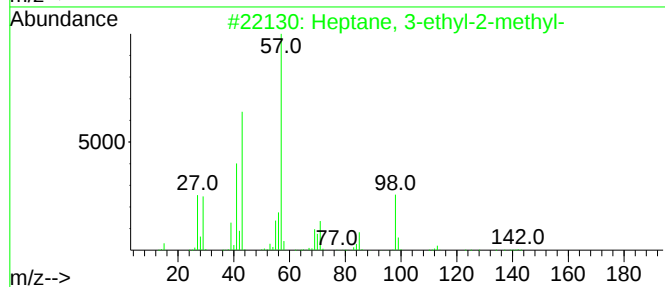
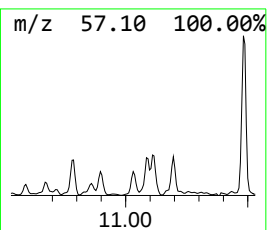
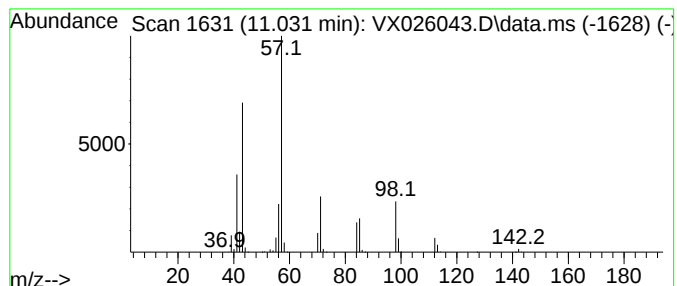
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 26

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 11.031 | 67.59 ug/L | 1025400 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Heptane, 3-ethyl-2-methyl-          | 142 | C10H22     | 014676-29-0  | 64   |
| 2    |    |   | Undecane, 5,6-dimethyl-             | 184 | C13H28     | 017615-91-7  | 50   |
| 3    |    |   | Oxalic acid, isobutyl nonyl ester   | 272 | C15H28O4   | 1010309-37-4 | 50   |
| 4    |    |   | Octane, 2,3-dimethyl-               | 142 | C10H22     | 007146-60-3  | 49   |
| 5    |    |   | 2-Ethyl-1-hexanol, trifluoroacetate | 226 | C10H17F3O2 | 053800-08-1  | 47   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

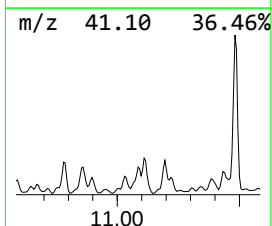
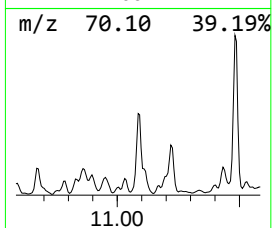
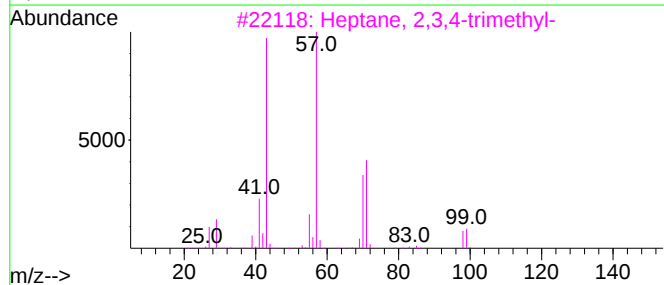
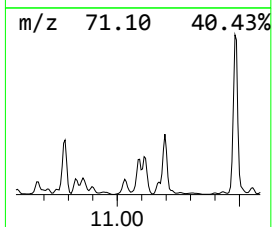
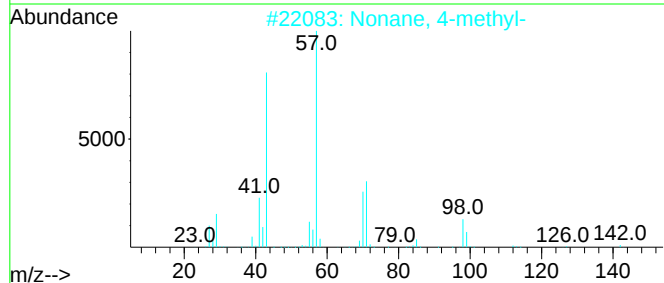
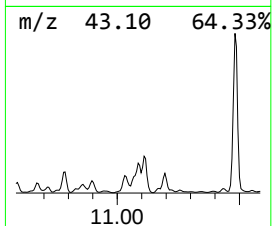
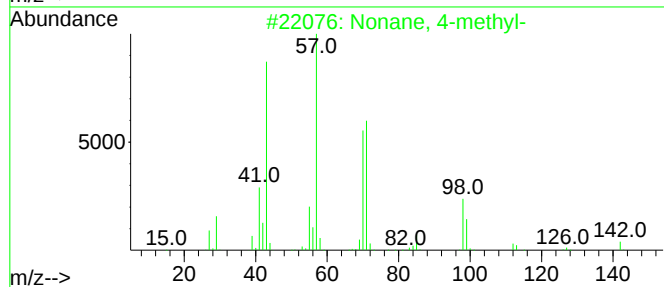
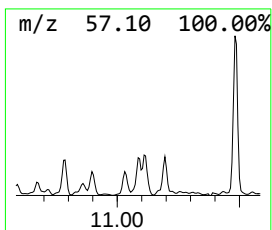
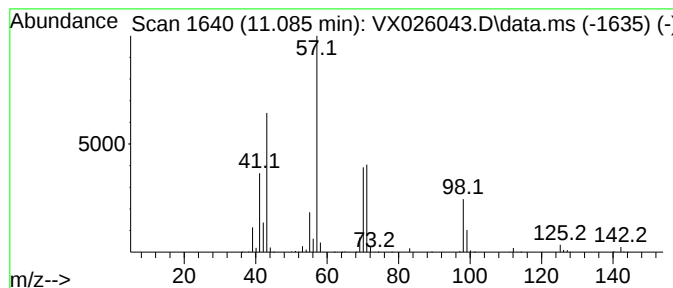
Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.      | EstConc                       | Area    | Relative to ISTD       |         |             | R.T.   |
|-----------|-------------------------------|---------|------------------------|---------|-------------|--------|
| 11.085    | 104.83 ug/L                   | 2523560 | 1,4-Dichlorobenzene-d4 |         |             | 12.024 |
| Hit# of 5 | Tentative ID                  |         | MW                     | MolForm | CAS#        | Qual   |
| 1         | Nonane, 4-methyl-             |         | 142                    | C10H22  | 017301-94-9 | 72     |
| 2         | Nonane, 4-methyl-             |         | 142                    | C10H22  | 017301-94-9 | 72     |
| 3         | Heptane, 2,3,4-trimethyl-     |         | 142                    | C10H22  | 052896-95-4 | 64     |
| 4         | Pentane, 2,2,3,3-tetramethyl- |         | 128                    | C9H20   | 007154-79-2 | 59     |
| 5         | Octane, 3,5-dimethyl-         |         | 142                    | C10H22  | 015869-93-9 | 58     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

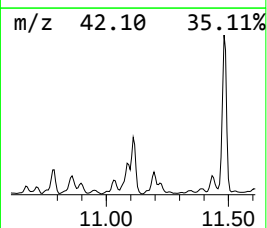
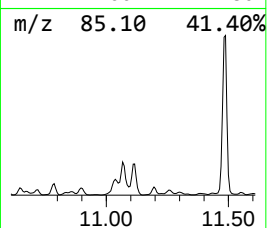
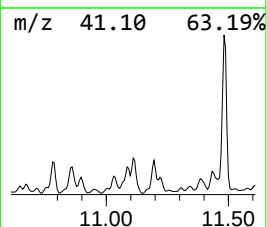
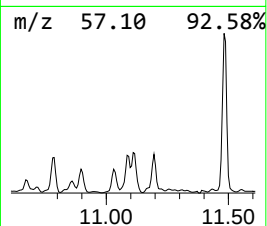
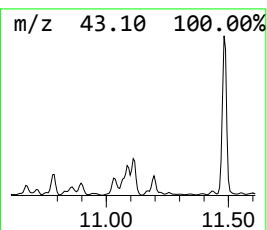
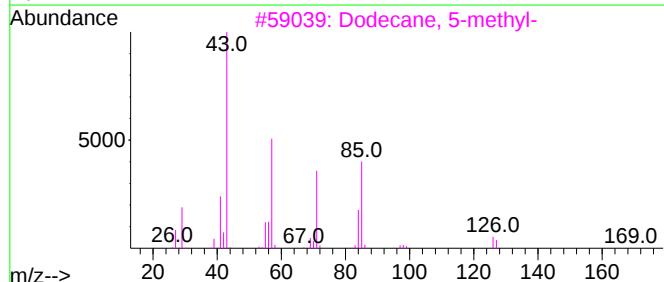
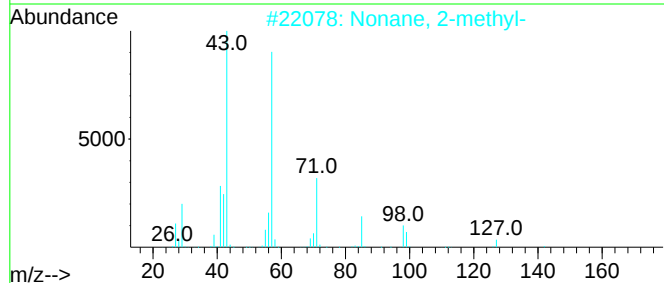
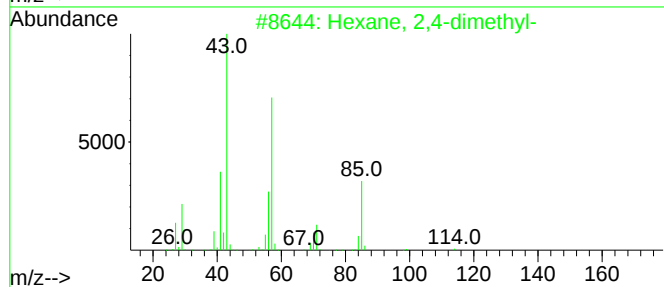
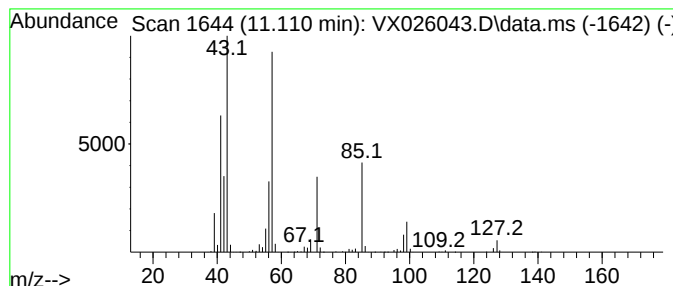
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.   | EstConc     | Area    | Relative to ISTD       | R.T.   |
|--------|-------------|---------|------------------------|--------|
| 11.110 | 140.89 ug/L | 3391540 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexane, 2,4-dimethyl-   | 114 | C8H18   | 000589-43-5 | 53   |
| 2    |    |   | Nonane, 2-methyl-       | 142 | C10H22  | 000871-83-0 | 53   |
| 3    |    |   | Dodecane, 5-methyl-     | 184 | C13H28  | 017453-93-9 | 53   |
| 4    |    |   | Undecane, 2,7-dimethyl- | 184 | C13H28  | 017301-24-5 | 53   |
| 5    |    |   | Hexane, 2,4-dimethyl-   | 114 | C8H18   | 000589-43-5 | 53   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

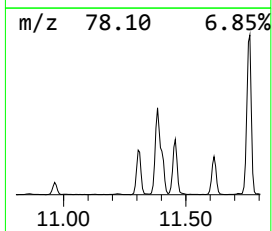
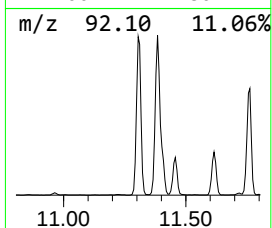
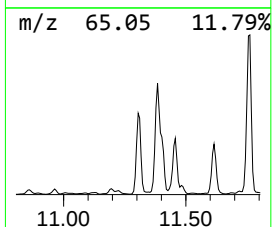
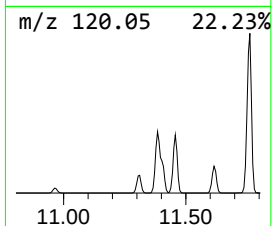
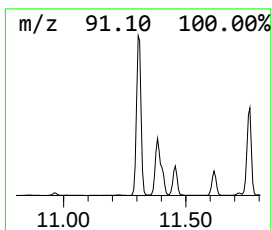
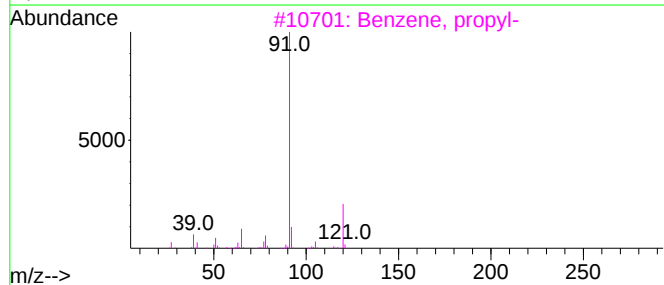
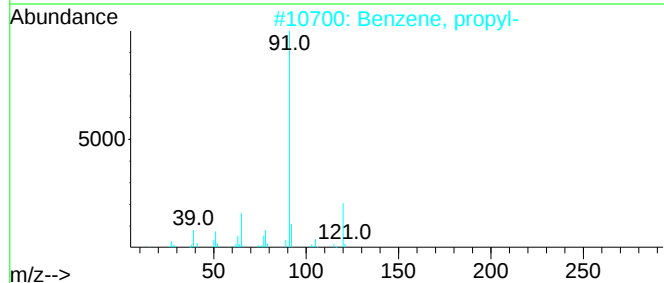
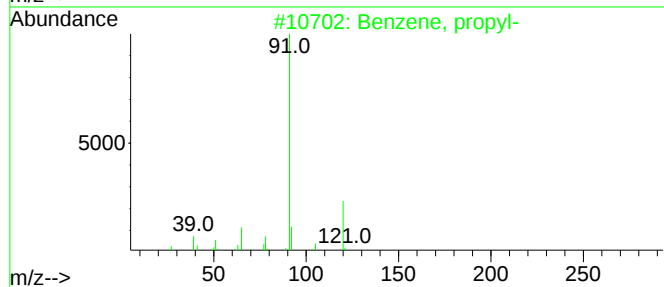
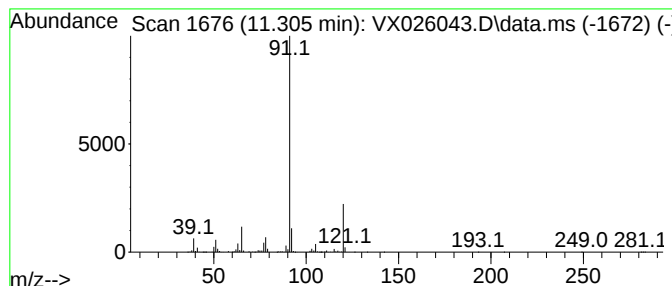
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 Benzene, propyl- Concentration Rank 12

| R.T.   | EstConc     | Area    | Relative to ISTD       | R.T.   |
|--------|-------------|---------|------------------------|--------|
| 11.305 | 151.81 ug/L | 3654330 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, propyl-          | 120 | C9H12   | 000103-65-1 | 94   |
| 2    |    |   | Benzene, propyl-          | 120 | C9H12   | 000103-65-1 | 94   |
| 3    |    |   | Benzene, propyl-          | 120 | C9H12   | 000103-65-1 | 90   |
| 4    |    |   | Benzene, propyl-          | 120 | C9H12   | 000103-65-1 | 87   |
| 5    |    |   | N-Benzyl-2-phenethylamine | 211 | C15H17N | 003647-71-0 | 83   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

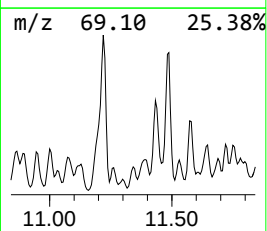
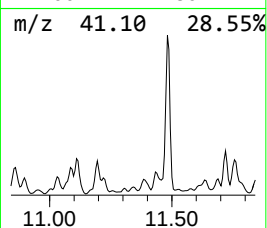
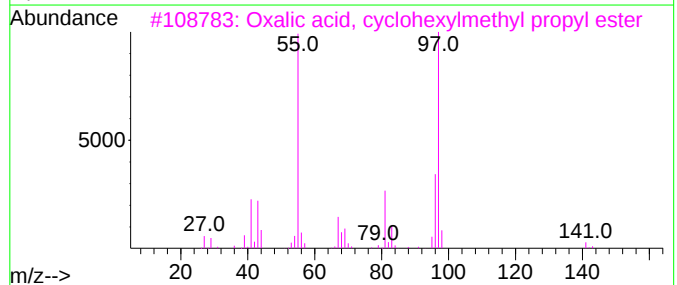
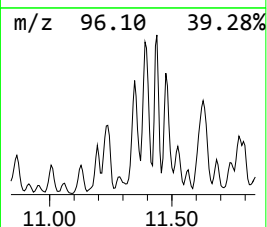
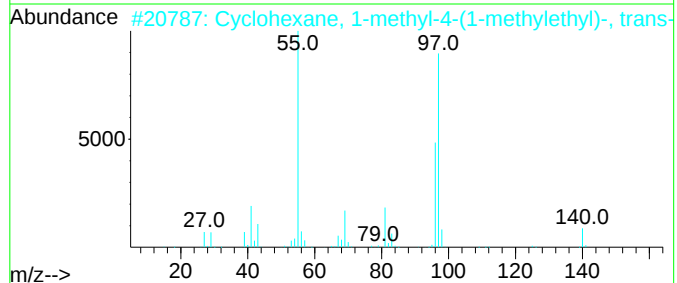
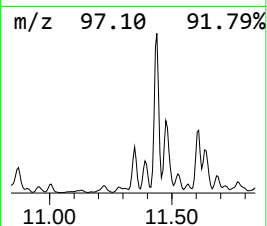
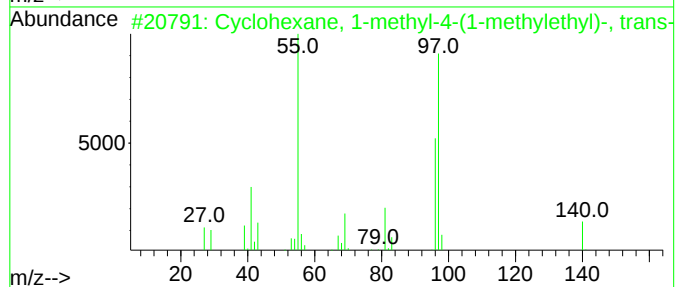
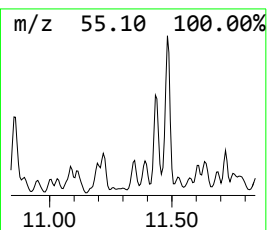
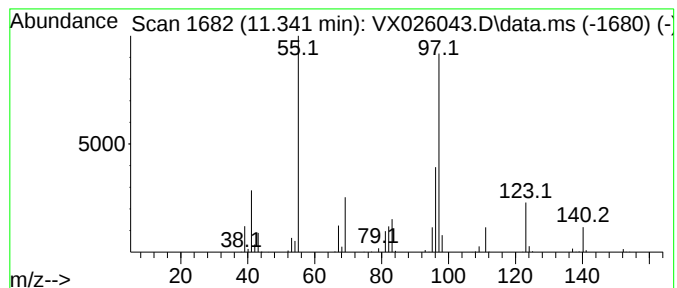
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 (DEL) Alkane: Cyclic11.342 Concentration Rank 37

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 11.342 | 32.15 ug/L | 773879 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Cyclohexane, 1-methyl-4-(1-methy... | 140 | C10H20   | 001678-82-6  | 64   |
| 2    |    |   | Cyclohexane, 1-methyl-4-(1-methy... | 140 | C10H20   | 001678-82-6  | 59   |
| 3    |    |   | Oxalic acid, cyclohexylmethyl pr... | 228 | C12H20O4 | 1010309-68-1 | 59   |
| 4    |    |   | m-Menthane, (1S,3S)-(+)-            | 140 | C10H20   | 013837-67-7  | 50   |
| 5    |    |   | 3-Hexene, 3-ethyl-2,5-dimethyl-     | 140 | C10H20   | 062338-08-3  | 47   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

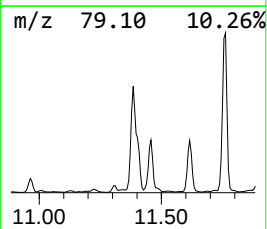
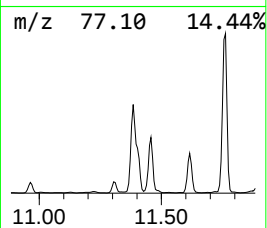
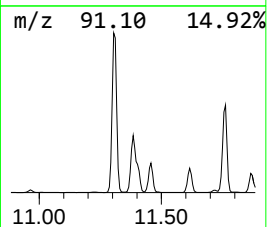
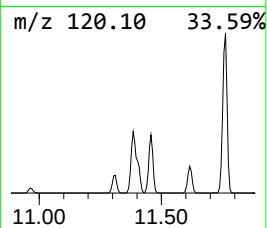
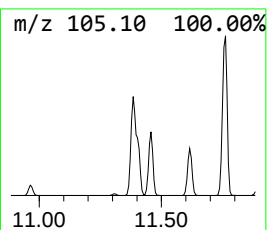
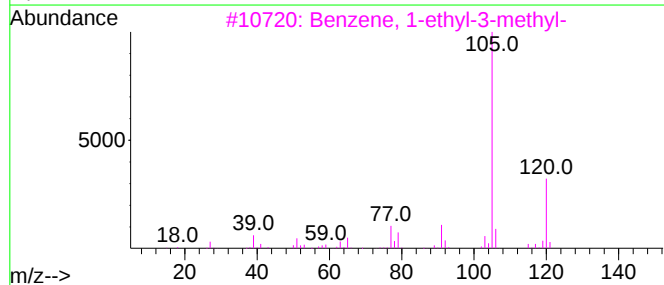
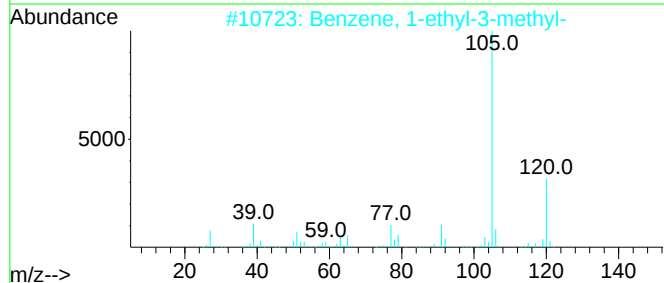
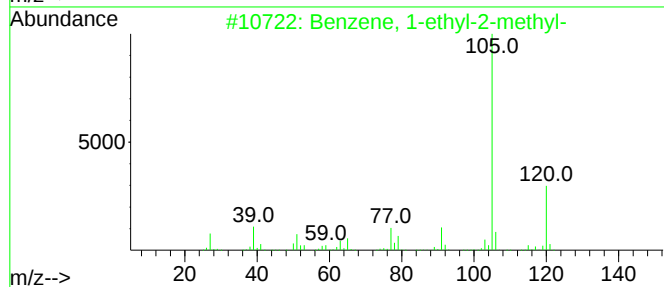
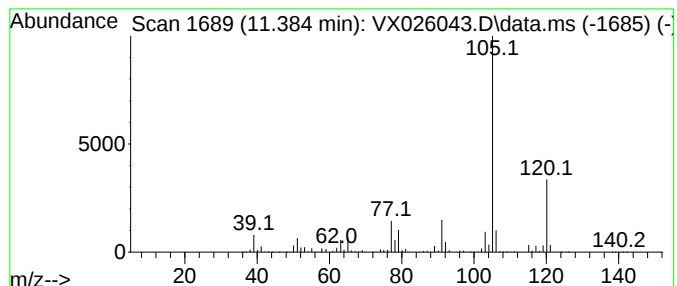
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

| R.T.   | EstConc     | Area     | Relative to ISTD       | R.T.   |
|--------|-------------|----------|------------------------|--------|
| 11.384 | 713.15 ug/L | 17166800 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 2    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 94   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 94   |
| 4    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 91   |
| 5    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

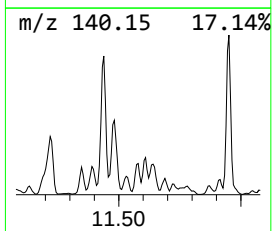
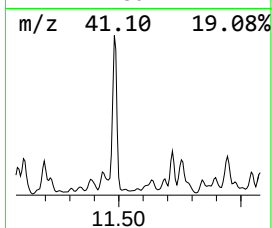
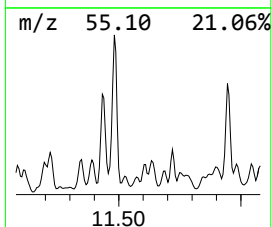
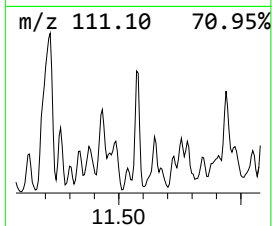
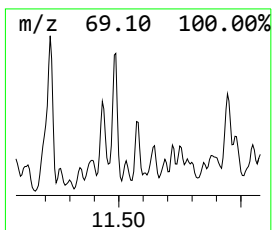
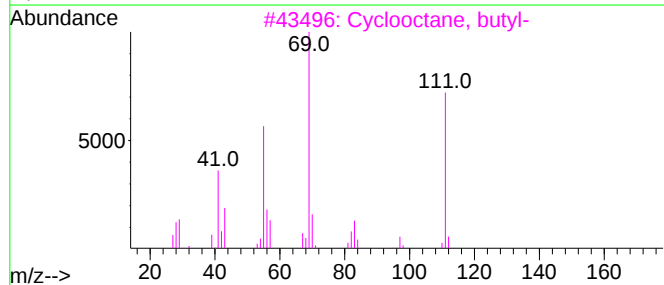
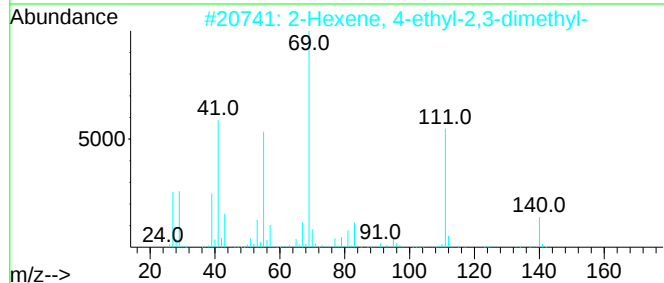
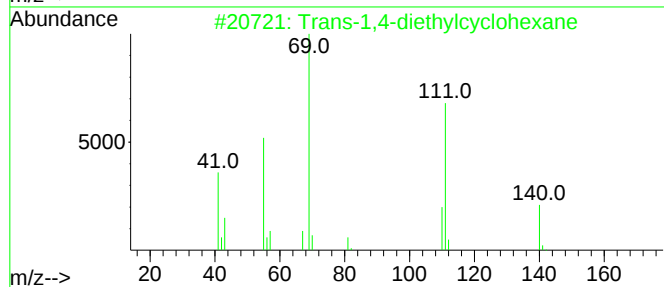
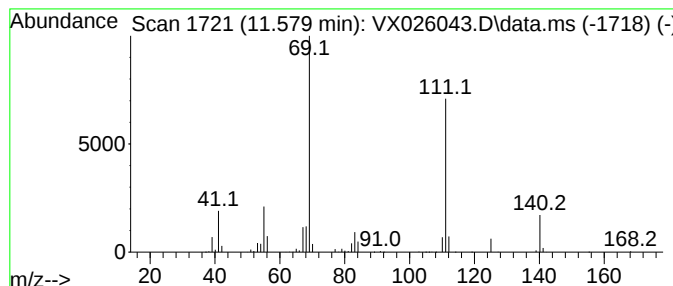
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 (DEL) Alkane: Cyclic11.579 Concentration Rank 54

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 11.579 | 8.07 ug/L | 194309 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Trans-1,4-diethylcyclohexane        | 140 | C10H20  | 013990-93-7  | 64   |
| 2    |    |   | 2-Hexene, 4-ethyl-2,3-dimethyl-     | 140 | C10H20  | 1000149-19-7 | 64   |
| 3    |    |   | Cyclooctane, butyl-                 | 168 | C12H24  | 016538-93-5  | 50   |
| 4    |    |   | 2-Octene, 2,7-dimethyl-             | 140 | C10H20  | 002050-81-9  | 50   |
| 5    |    |   | Cyclohexane, 1-ethyl-1,3-dimethy... | 140 | C10H20  | 062238-29-3  | 45   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

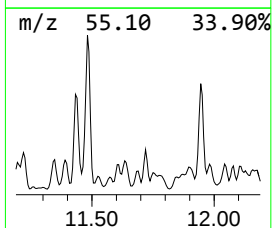
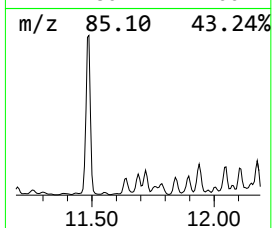
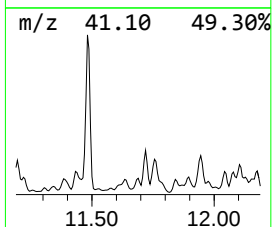
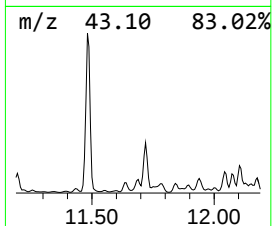
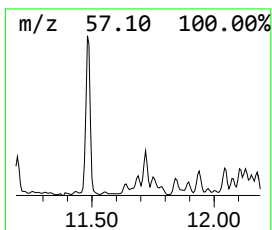
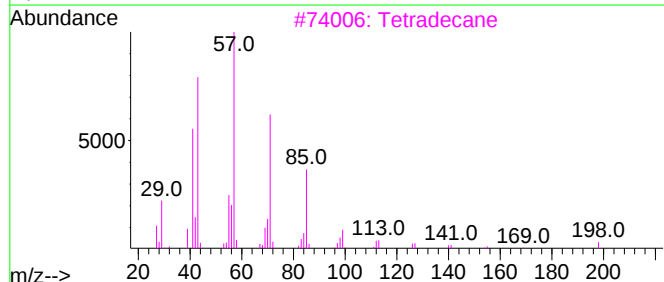
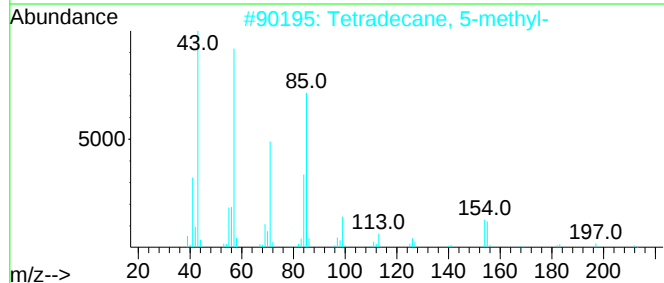
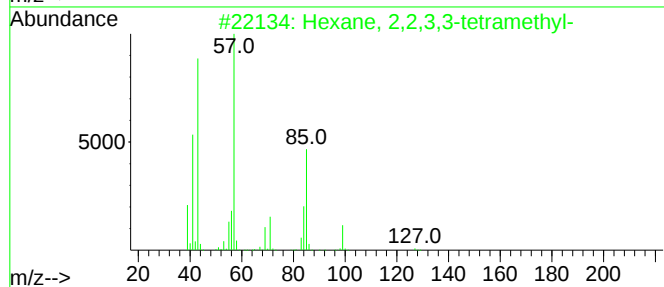
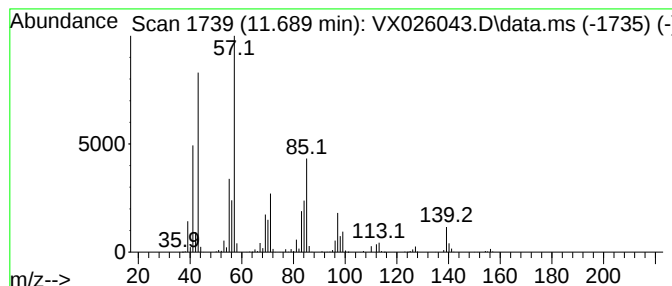
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 11.689 | 48.61 ug/L | 1170190 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexane, 2,2,3,3-tetramethyl- | 142 | C10H22  | 013475-81-5 | 53   |
| 2    |    |   | Tetradecane, 5-methyl-       | 212 | C15H32  | 025117-32-2 | 53   |
| 3    |    |   | Tetradecane                  | 198 | C14H30  | 000629-59-4 | 47   |
| 4    |    |   | Hexane, 2,3,3-trimethyl-     | 128 | C9H20   | 016747-28-7 | 47   |
| 5    |    |   | Hexane, 2,3,5-trimethyl-     | 128 | C9H20   | 001069-53-0 | 43   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

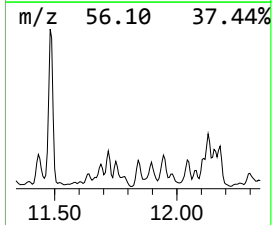
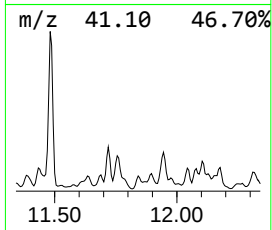
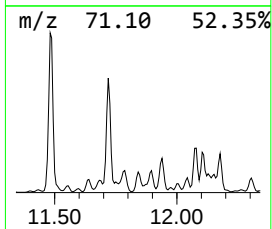
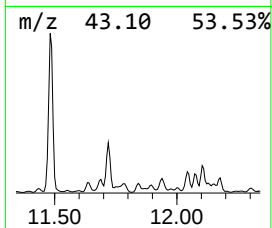
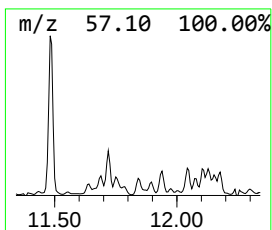
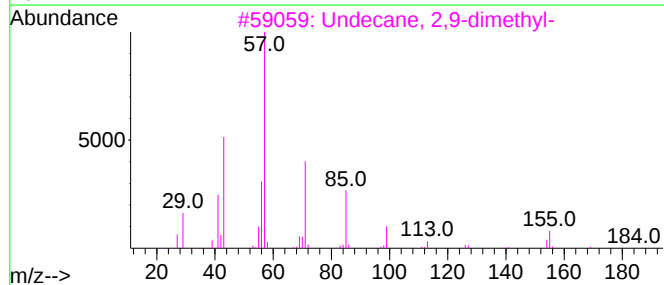
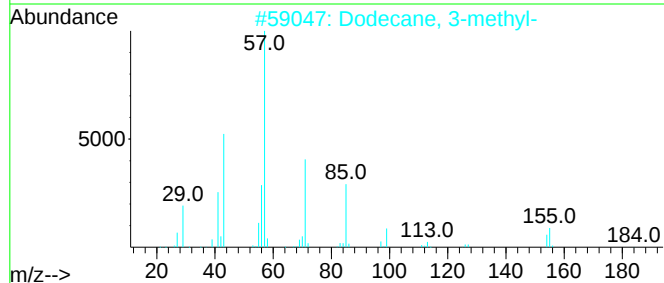
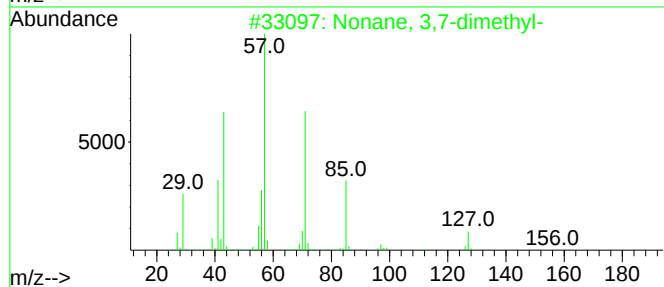
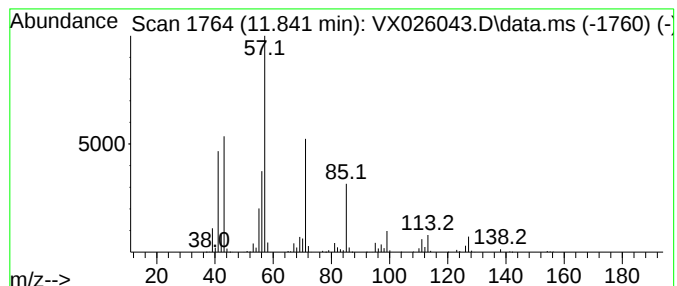
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 31

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 11.841 | 53.37 ug/L | 1284660 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 3,7-dimethyl-   | 156 | C11H24  | 017302-32-8 | 81   |
| 2    |    |   | Dodecane, 3-methyl-     | 184 | C13H28  | 017312-57-1 | 72   |
| 3    |    |   | Undecane, 2,9-dimethyl- | 184 | C13H28  | 017301-26-7 | 72   |
| 4    |    |   | Undecane, 2-methyl-     | 170 | C12H26  | 007045-71-8 | 64   |
| 5    |    |   | Undecane, 5-ethyl-      | 184 | C13H28  | 017453-94-0 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

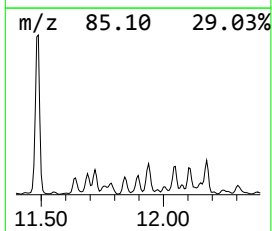
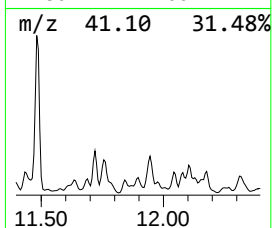
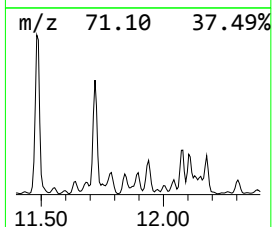
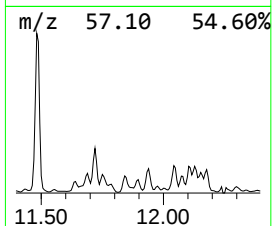
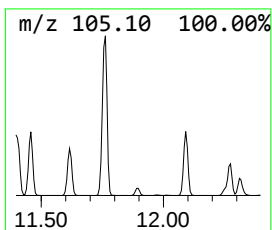
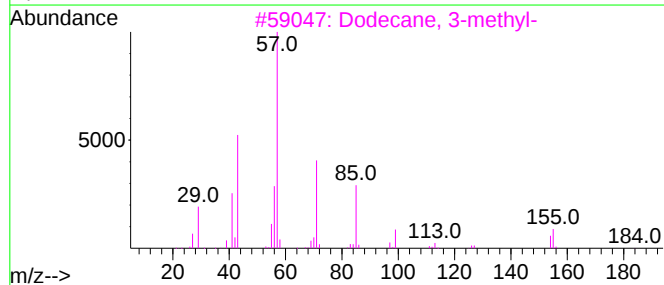
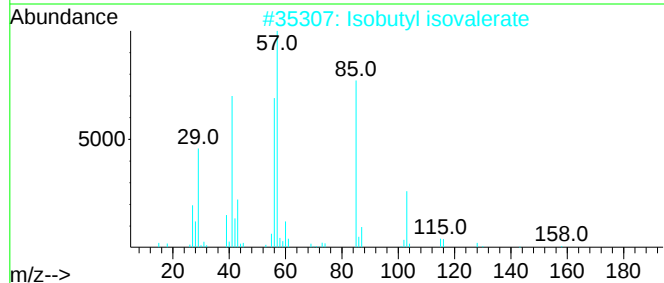
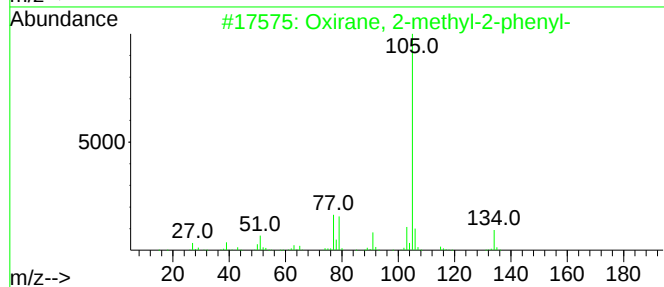
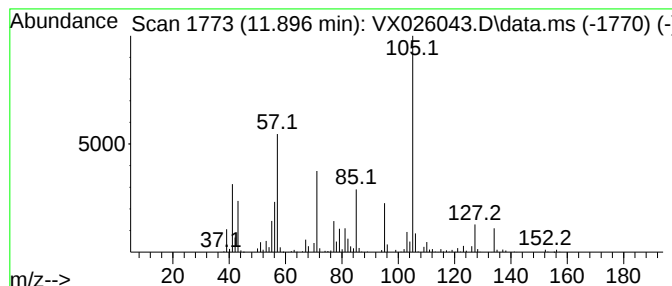
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 Oxirane, 2-methyl-2-phenyl- Concentration Rank 25

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 11.896 | 67.97 ug/L | 1636040 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Oxirane, 2-methyl-2-phenyl-         | 134 | C9H10O  | 002085-88-3 | 35   |
| 2    |    |   | Isobutyl isovalerate                | 158 | C9H18O2 | 000589-59-3 | 22   |
| 3    |    |   | Dodecane, 3-methyl-                 | 184 | C13H28  | 017312-57-1 | 22   |
| 4    |    |   | Benzene, 1-methyl-4-propyl-         | 134 | C10H14  | 001074-55-1 | 18   |
| 5    |    |   | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O  | 000093-53-8 | 14   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

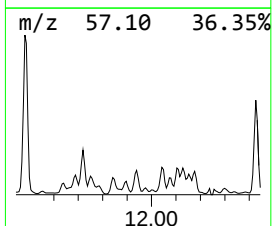
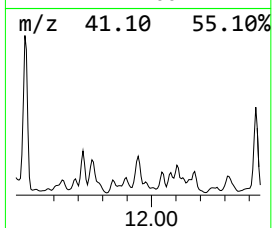
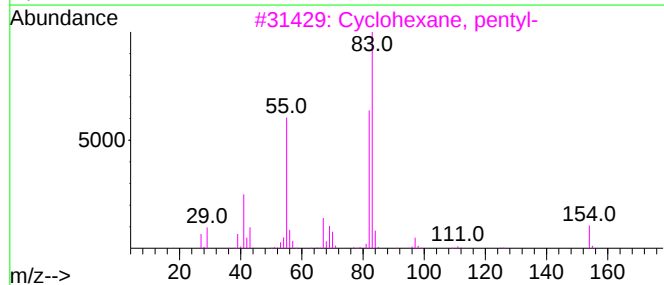
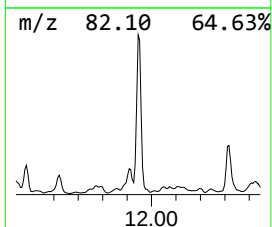
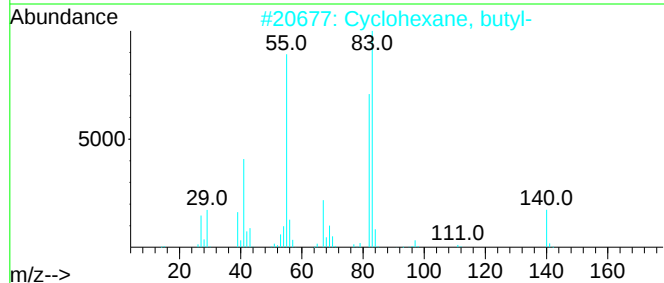
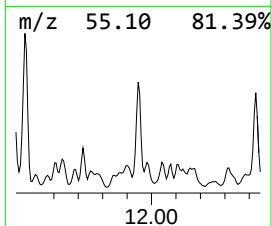
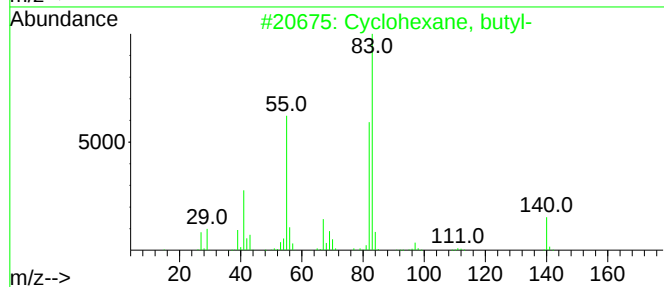
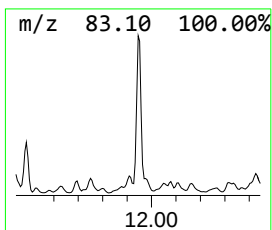
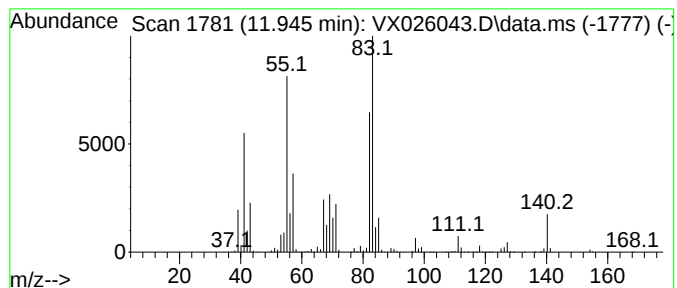
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 (DEL) Alkane: Cyclic11.945 Concentration Rank 11

| R.T.   | EstConc     | Area    | Relative to ISTD       | R.T.   |
|--------|-------------|---------|------------------------|--------|
| 11.945 | 175.08 ug/L | 4214570 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, butyl-   | 140 | C10H20  | 001678-93-9 | 76   |
| 2    |    |   | Cyclohexane, butyl-   | 140 | C10H20  | 001678-93-9 | 72   |
| 3    |    |   | Cyclohexane, pentyl-  | 154 | C11H22  | 004292-92-6 | 64   |
| 4    |    |   | Cyclohexane, undecyl- | 238 | C17H34  | 054105-66-7 | 64   |
| 5    |    |   | Cyclohexane, butyl-   | 140 | C10H20  | 001678-93-9 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

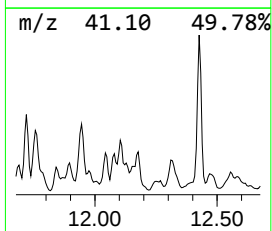
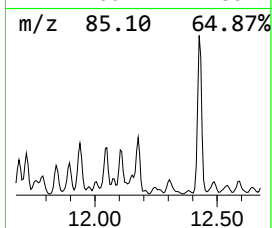
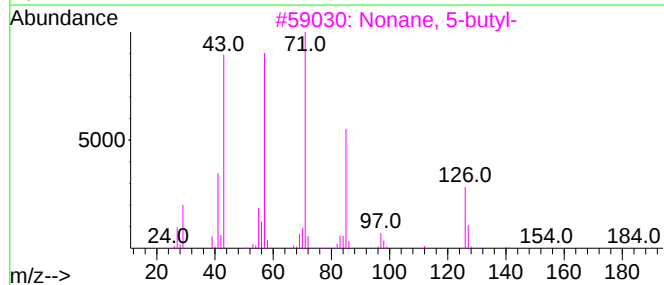
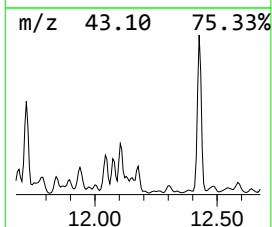
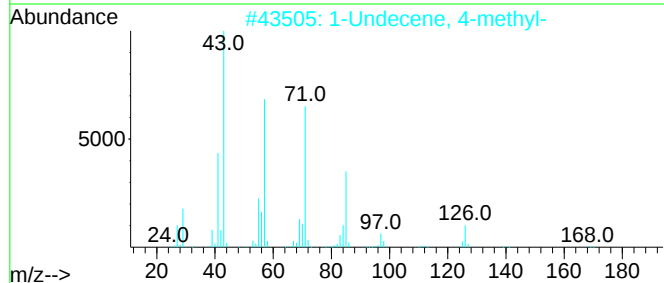
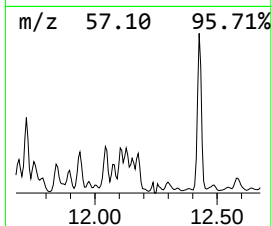
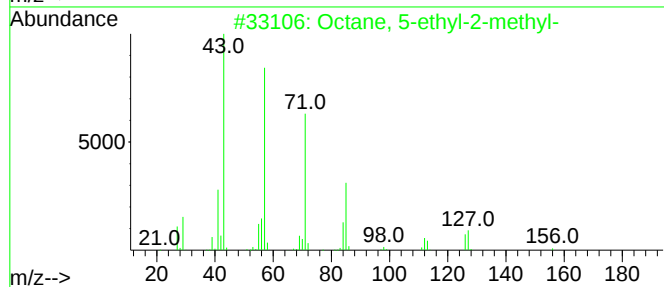
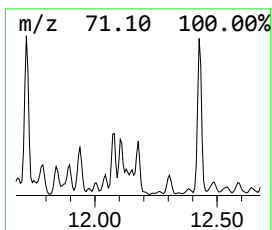
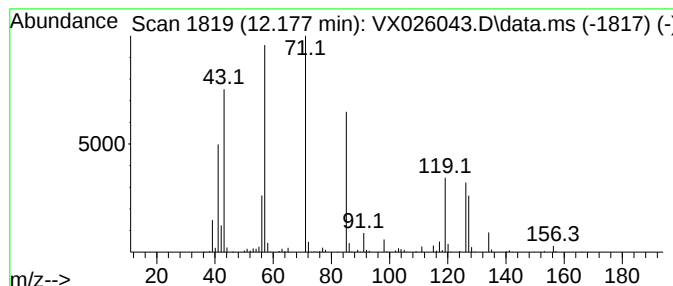
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 12.177 | 88.27 ug/L | 2124920 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of                         | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----------------------------|---|--------------|-----|---------|-------------|------|
| 1    | Octane, 5-ethyl-2-methyl-  |   |              | 156 | C11H24  | 062016-18-6 | 59   |
| 2    | 1-Undecene, 4-methyl-      |   |              | 168 | C12H24  | 074630-39-0 | 59   |
| 3    | Nonane, 5-butyl-           |   |              | 184 | C13H28  | 017312-63-9 | 53   |
| 4    | Nonane, 5-methyl-5-propyl- |   |              | 184 | C13H28  | 017312-75-3 | 53   |
| 5    | Undecane, 2-methyl-        |   |              | 170 | C12H26  | 007045-71-8 | 53   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

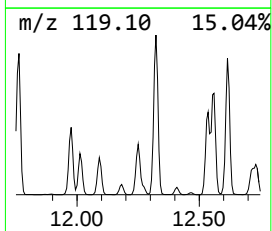
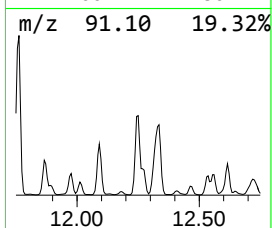
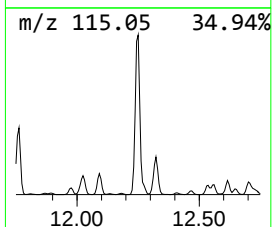
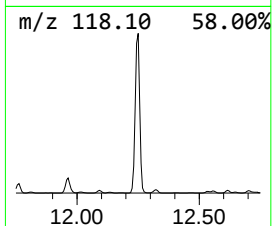
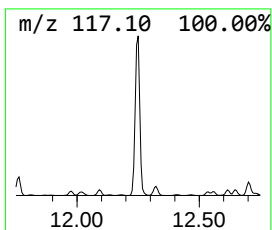
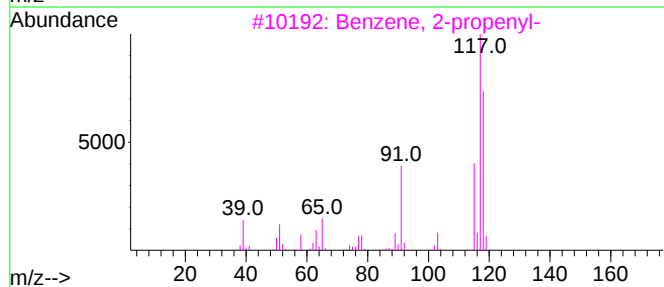
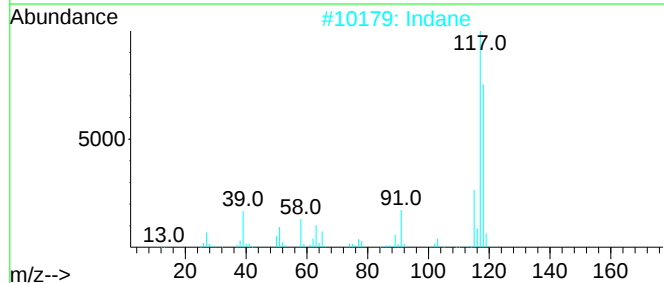
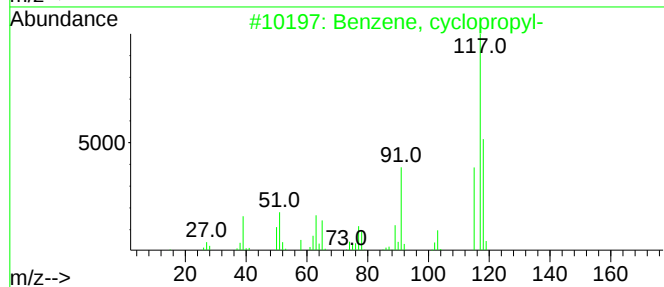
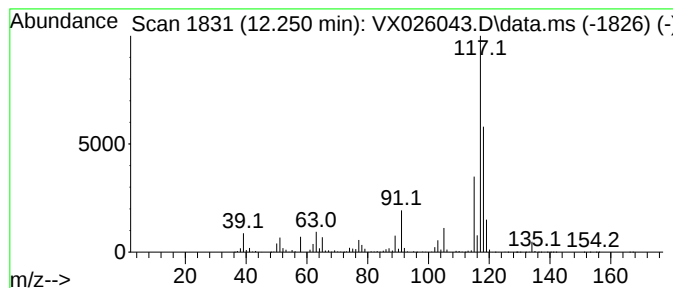
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 37 Benzene, cyclopropyl- Concentration Rank 3

| R.T.   | EstConc     | Area    | Relative to ISTD       | R.T.   |
|--------|-------------|---------|------------------------|--------|
| 12.250 | 401.47 ug/L | 9663960 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, cyclopropyl- | 118 | C9H10   | 000873-49-4 | 76   |
| 2    |    |   | Indane                | 118 | C9H10   | 000496-11-7 | 76   |
| 3    |    |   | Benzene, 2-propenyl-  | 118 | C9H10   | 000300-57-2 | 76   |
| 4    |    |   | Indane                | 118 | C9H10   | 000496-11-7 | 76   |
| 5    |    |   | Benzene, cyclopropyl- | 118 | C9H10   | 000873-49-4 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

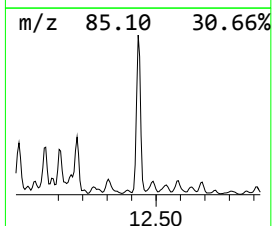
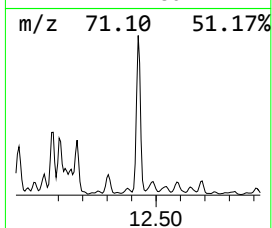
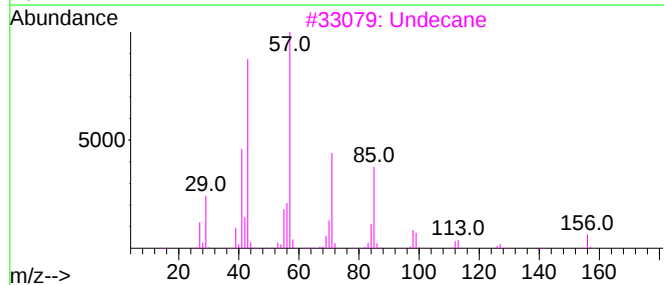
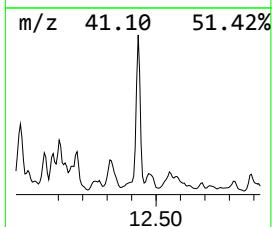
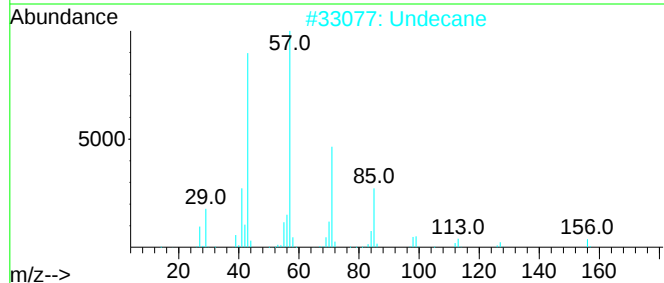
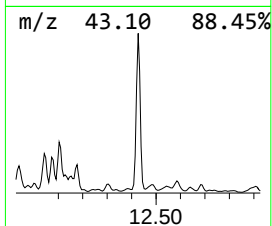
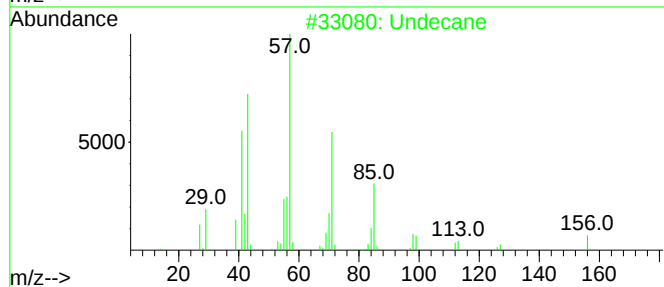
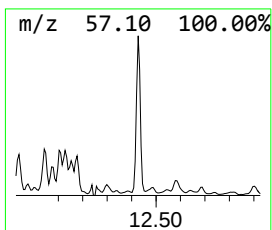
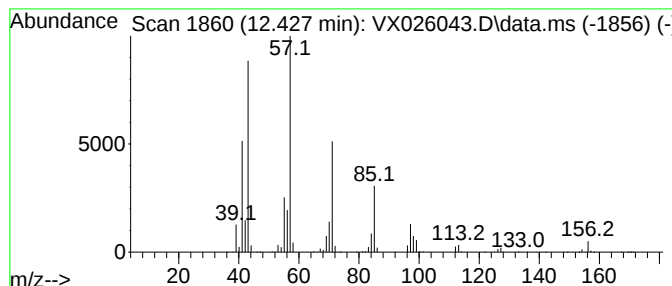
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.   | EstConc     | Area    | Relative to ISTD       | R.T.   |
|--------|-------------|---------|------------------------|--------|
| 12.427 | 272.57 ug/L | 6561160 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of       | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----------|---|--------------|-----|---------|-------------|------|
| 1    | Undecane |   |              | 156 | C11H24  | 001120-21-4 | 96   |
| 2    | Undecane |   |              | 156 | C11H24  | 001120-21-4 | 94   |
| 3    | Undecane |   |              | 156 | C11H24  | 001120-21-4 | 94   |
| 4    | Undecane |   |              | 156 | C11H24  | 001120-21-4 | 90   |
| 5    | Undecane |   |              | 156 | C11H24  | 001120-21-4 | 90   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

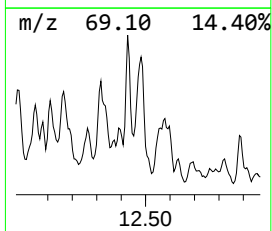
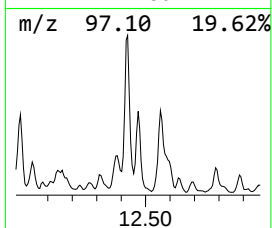
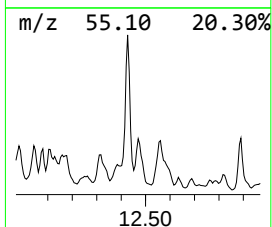
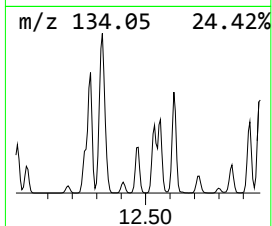
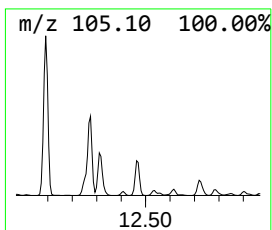
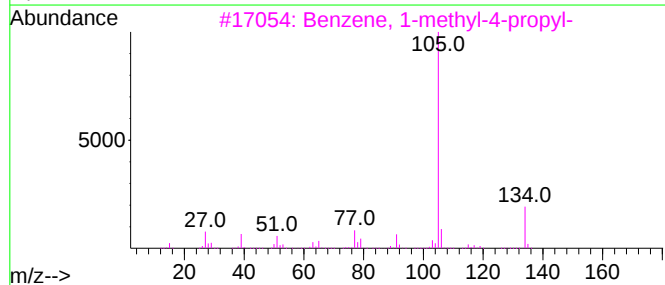
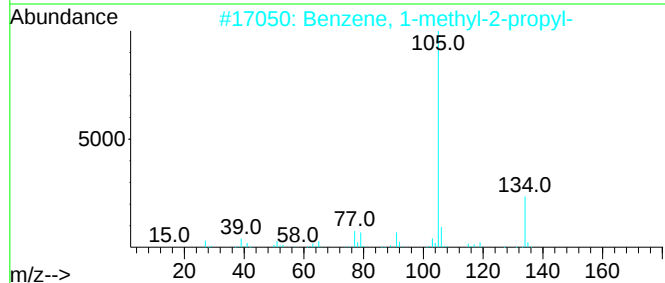
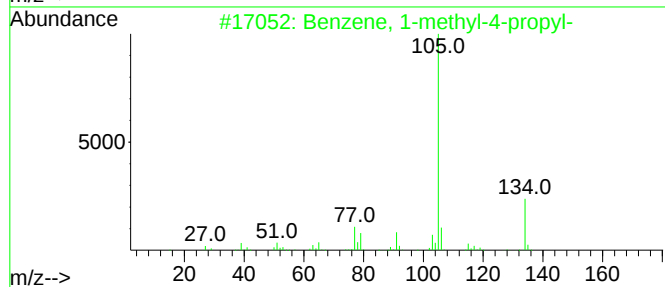
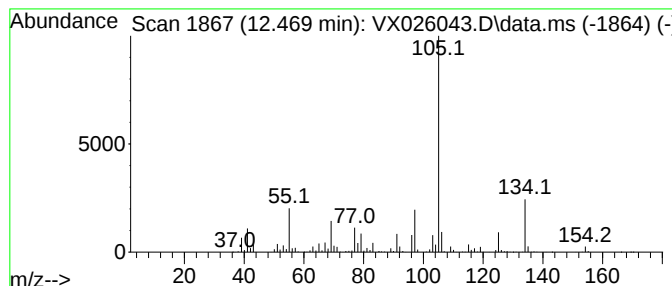
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 Benzene, 1-methyl-4-propyl- Concentration Rank 17

| R.T.   | EstConc     | Area    | Relative to ISTD       | R.T.   |
|--------|-------------|---------|------------------------|--------|
| 12.469 | 109.30 ug/L | 2631060 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-4-propyl-         | 134 | C10H14  | 001074-55-1 | 90   |
| 2    |    |   | Benzene, 1-methyl-2-propyl-         | 134 | C10H14  | 001074-17-5 | 90   |
| 3    |    |   | Benzene, 1-methyl-4-propyl-         | 134 | C10H14  | 001074-55-1 | 60   |
| 4    |    |   | Benzene, (1-methylpropyl)-          | 134 | C10H14  | 000135-98-8 | 60   |
| 5    |    |   | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O  | 000093-53-8 | 60   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

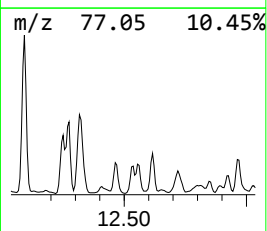
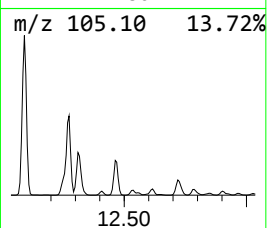
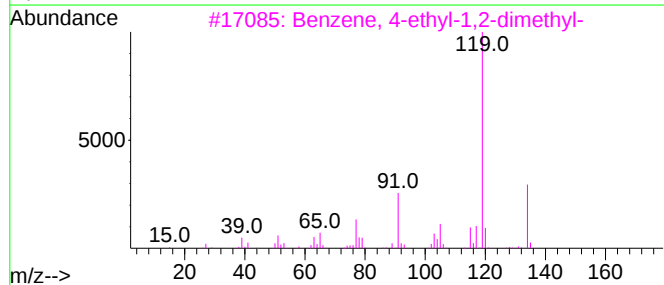
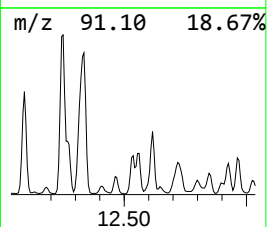
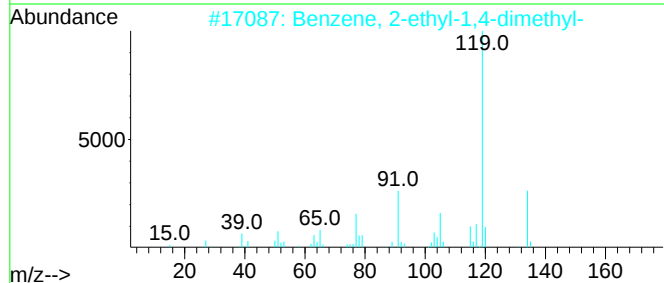
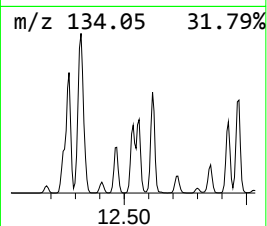
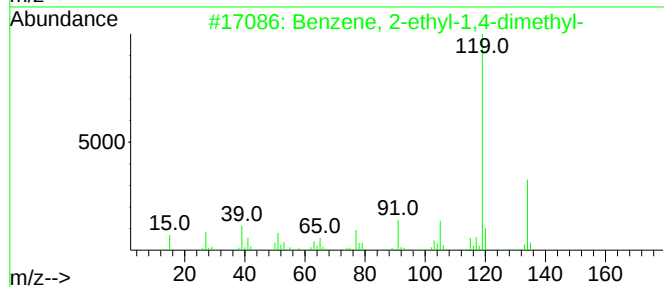
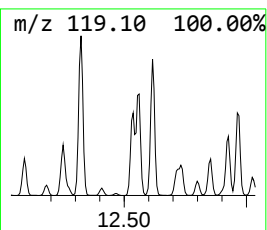
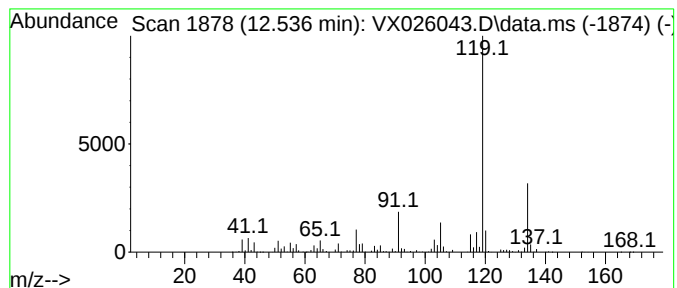
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 40 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 24

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 12.536 | 72.99 ug/L | 1757100 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 97   |
| 2    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 96   |
| 3    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 96   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 95   |
| 5    |    |   | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 95   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

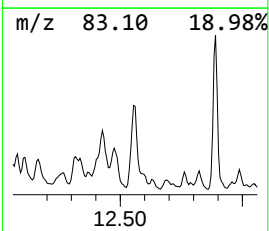
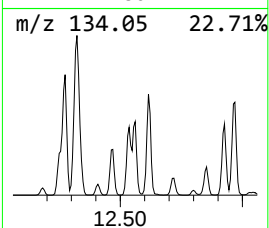
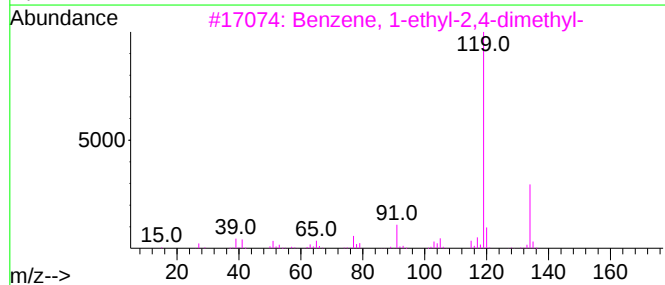
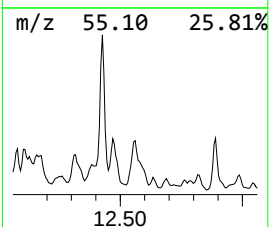
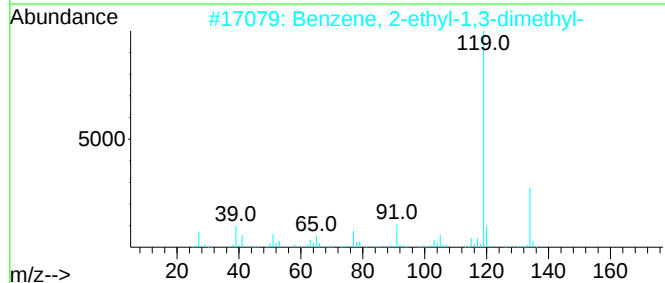
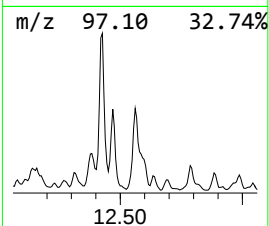
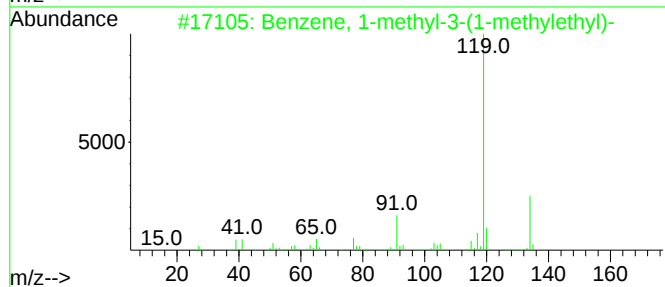
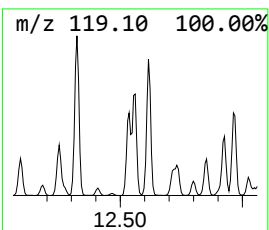
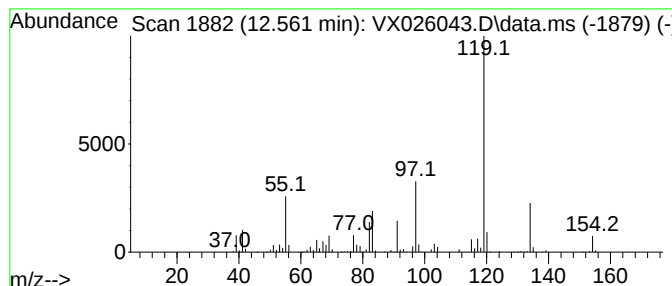
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 41 Benzene, 1-methyl-3-(1-meth... Concentration Rank 19

| R.T.   | EstConc     | Area    | Relative to ISTD       | R.T.   |
|--------|-------------|---------|------------------------|--------|
| 12.561 | 104.37 ug/L | 2512340 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 89   |
| 2    |    |   | Benzene, 2-ethyl-1,3-dimethyl-      | 134 | C10H14  | 002870-04-4 | 76   |
| 3    |    |   | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9 | 76   |
| 4    |    |   | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 76   |
| 5    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 76   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

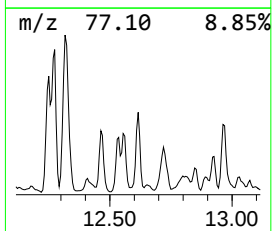
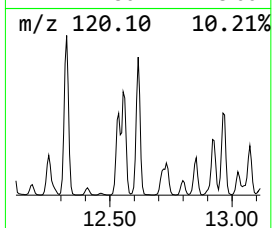
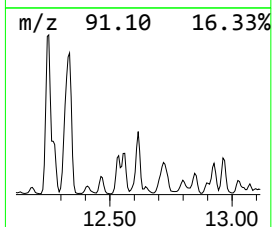
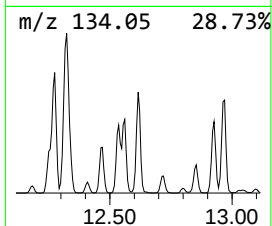
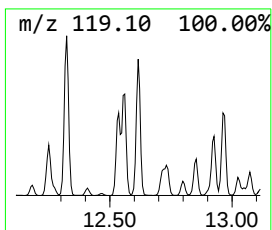
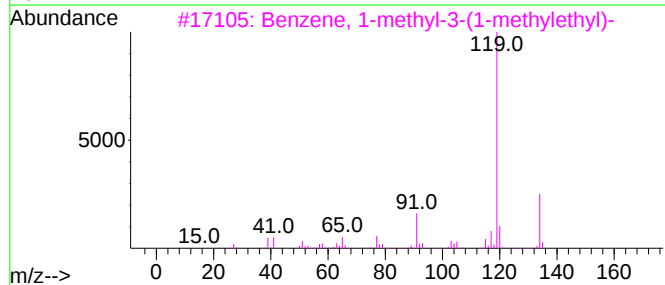
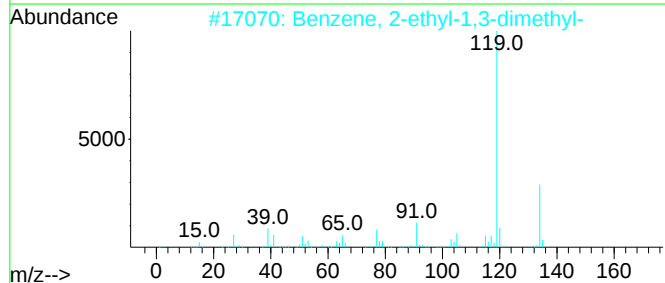
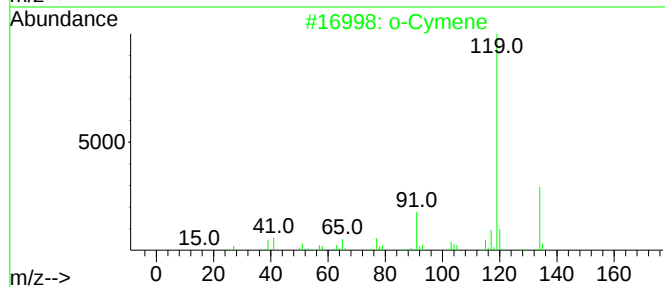
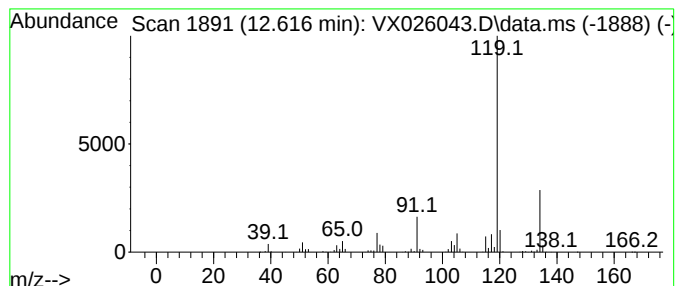
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 42 o-Cymene Concentration Rank 23

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 12.616 | 75.77 ug/L | 1824020 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|---------|-------------|------|
| 1    |    |   | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 97   |
| 2    |    |   | Benzene, 2-ethyl-1,3-dimethyl-       | 134 | C10H14  | 002870-04-4 | 95   |
| 3    |    |   | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 95   |
| 4    |    |   | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 95   |
| 5    |    |   | p-Cymene                             | 134 | C10H14  | 000099-87-6 | 95   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

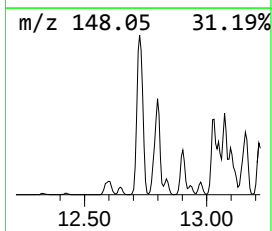
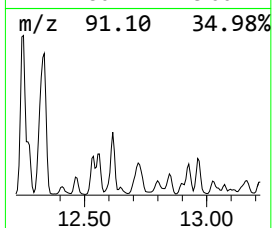
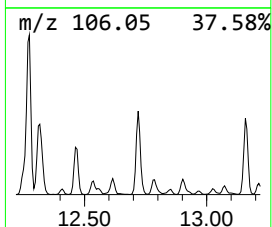
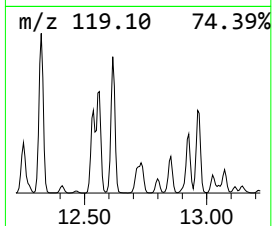
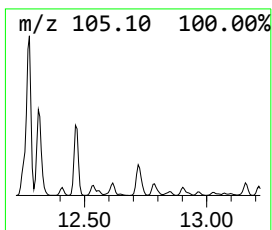
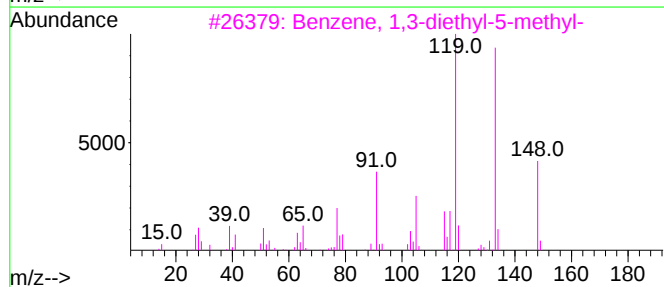
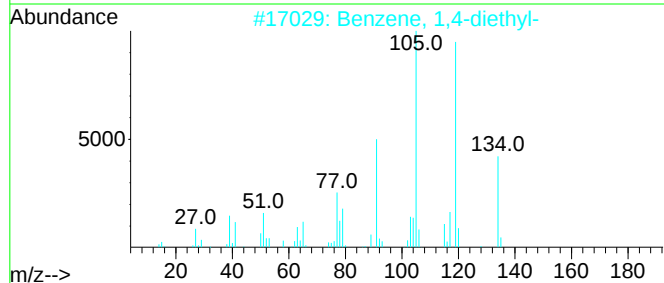
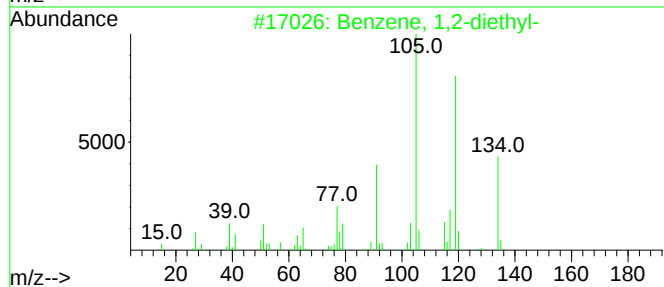
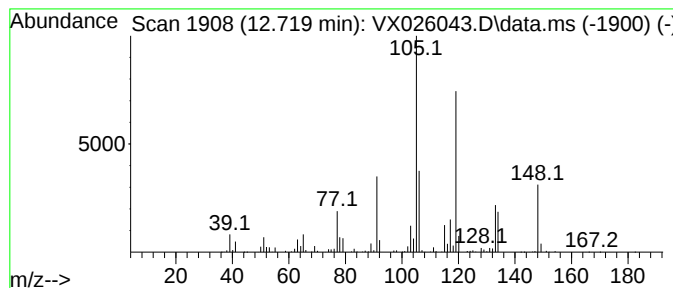
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 43 Benzene, 1,2-diethyl- Concentration Rank 21

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 12.719 | 89.42 ug/L | 2152530 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 70   |
| 2    |    |   | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 70   |
| 3    |    |   | Benzene, 1,3-diethyl-5-methyl- | 148 | C11H16  | 002050-24-0 | 62   |
| 4    |    |   | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 58   |
| 5    |    |   | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 58   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

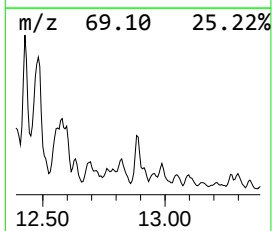
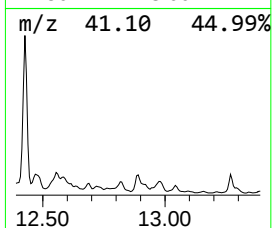
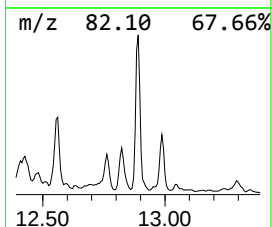
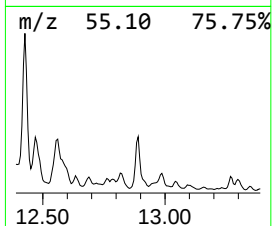
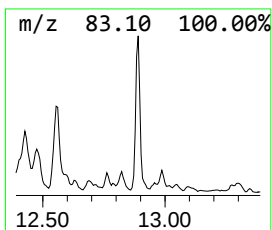
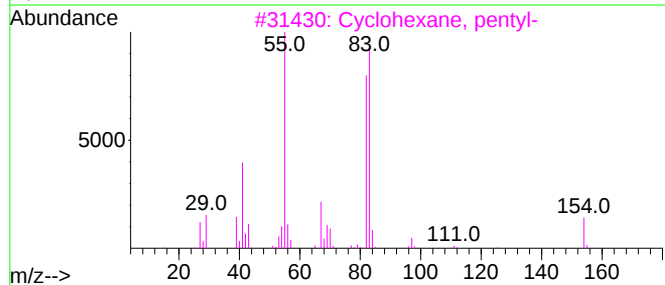
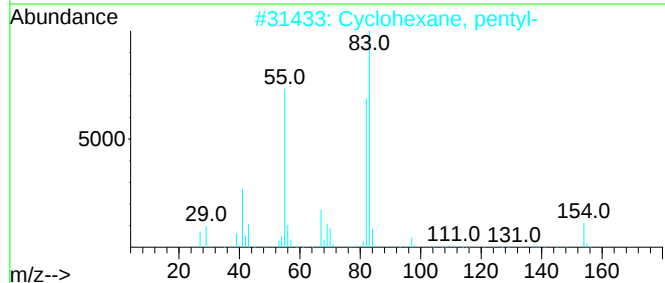
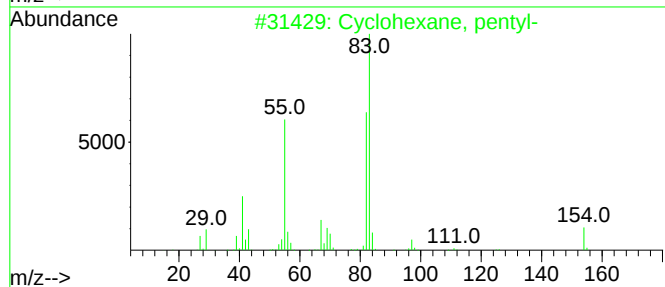
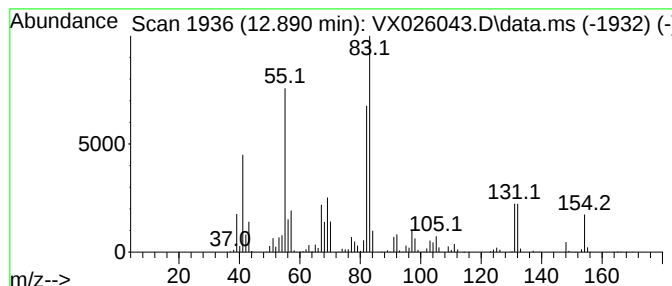
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 44 (DEL) Alkane: Cyclic12.890 Concentration Rank 29

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 12.890 | 59.71 ug/L | 1437290 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 81   |
| 2    |    |   | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 76   |
| 3    |    |   | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 68   |
| 4    |    |   | Cyclohexane, (4-methylpentyl)- | 168 | C12H24  | 061142-20-9 | 68   |
| 5    |    |   | Cyclohexane, octyl-            | 196 | C14H28  | 001795-15-9 | 59   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

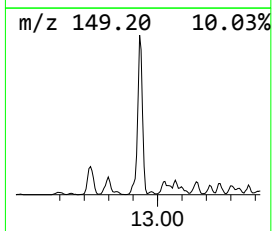
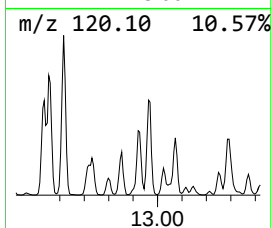
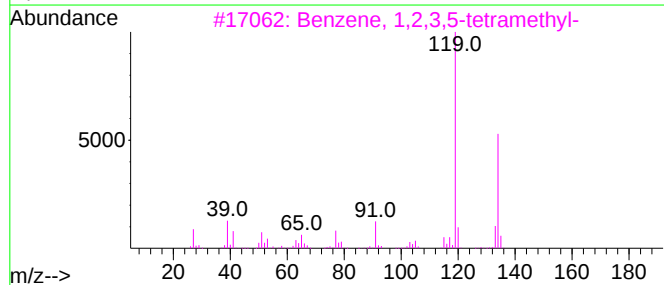
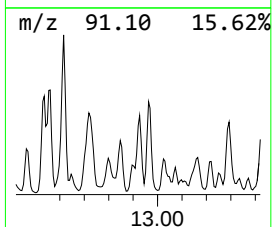
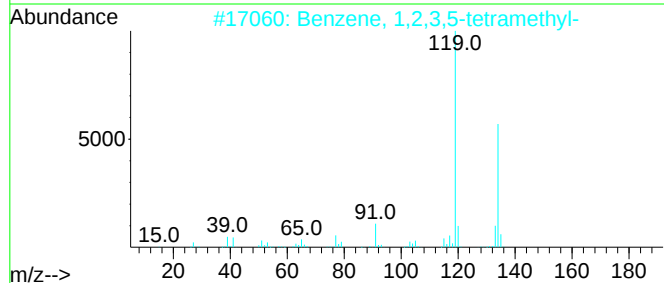
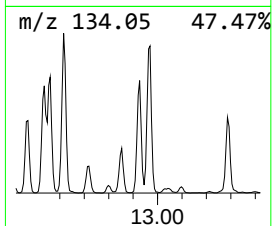
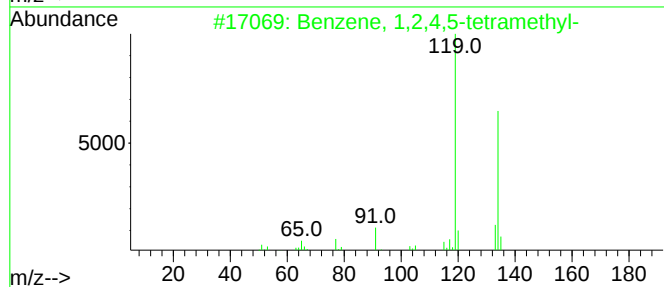
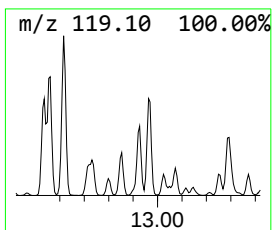
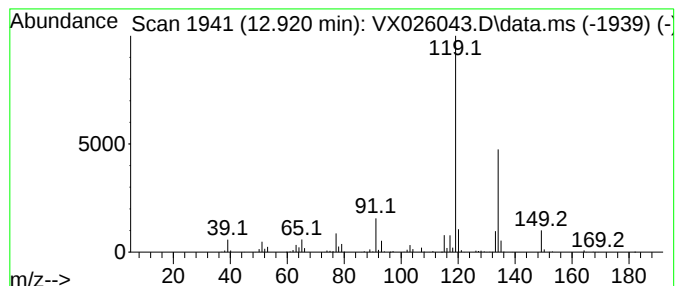
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 45 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 30

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 12.921 | 54.84 ug/L | 1319970 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 94   |
| 2    |    |   | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 94   |
| 3    |    |   | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 94   |
| 4    |    |   | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 93   |
| 5    |    |   | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 92   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

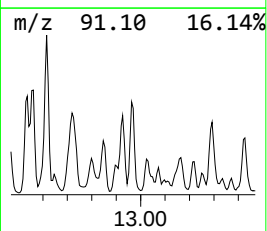
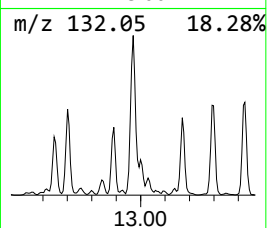
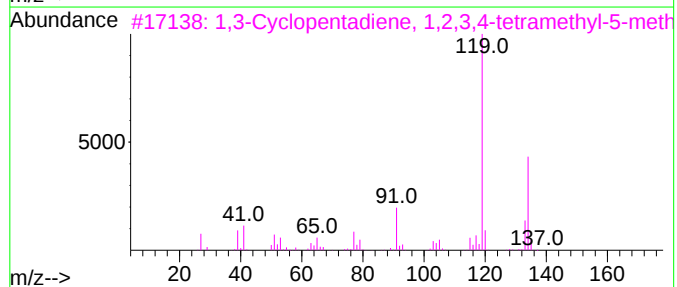
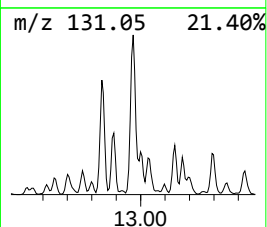
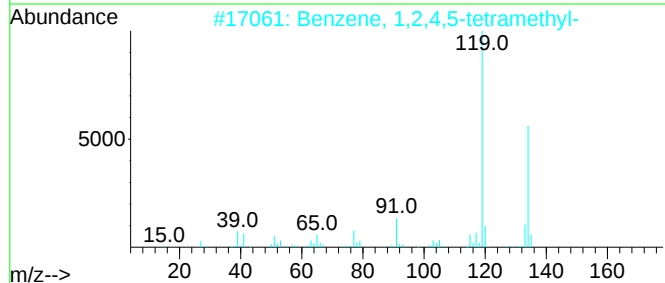
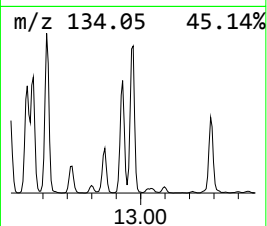
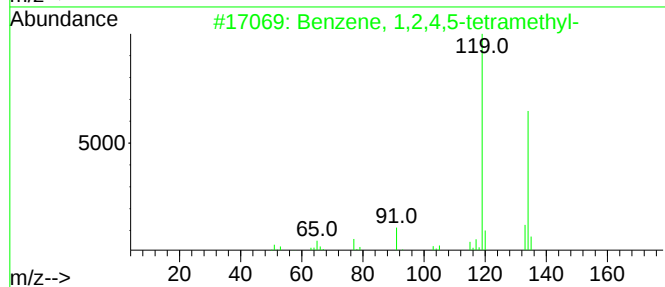
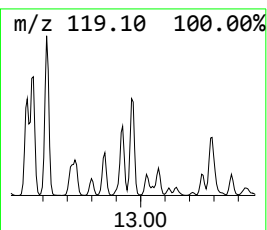
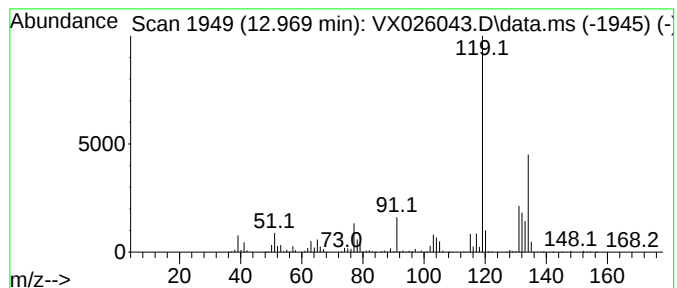
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 46 1,3-Cyclopentadiene, 1,2,3,... Concentration Rank 16

| R.T.   | EstConc     | Area    | Relative to ISTD       | R.T.   |
|--------|-------------|---------|------------------------|--------|
| 12.969 | 113.79 ug/L | 2739050 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 81   |
| 2    |    |   | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 81   |
| 3    |    |   | 1,3-Cyclopentadiene, 1,2,3,4-tet... | 134 | C10H14  | 076089-59-3 | 81   |
| 4    |    |   | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 81   |
| 5    |    |   | Benzene, 1,2,3,5-tetramethyl-       | 134 | C10H14  | 000527-53-7 | 81   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

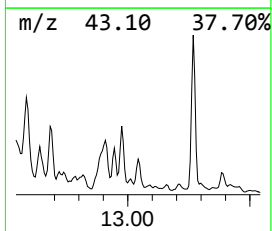
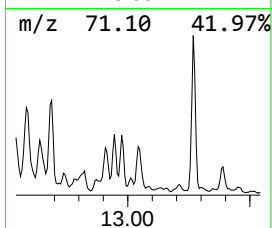
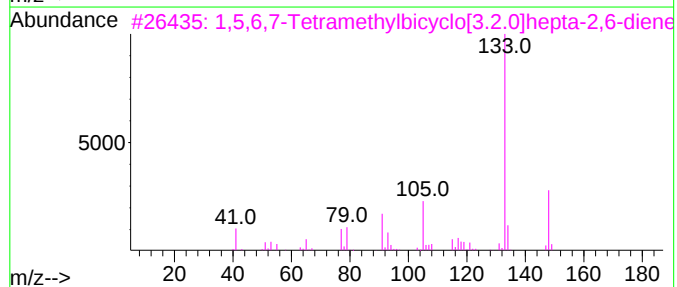
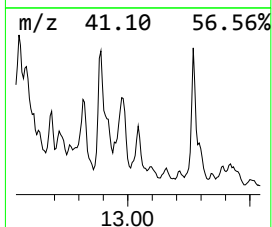
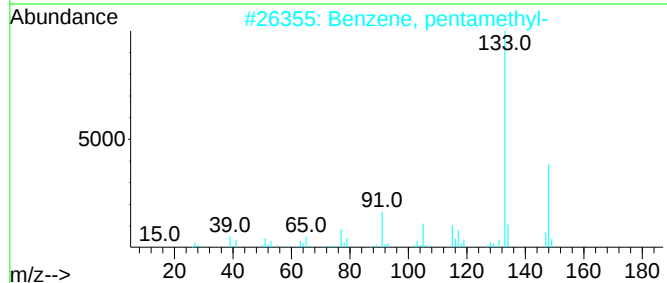
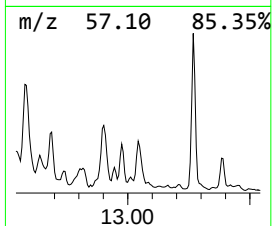
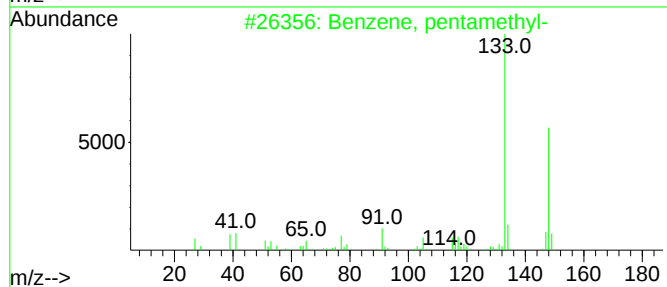
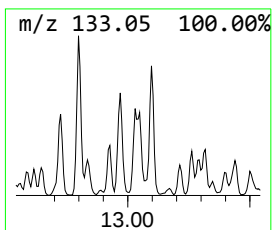
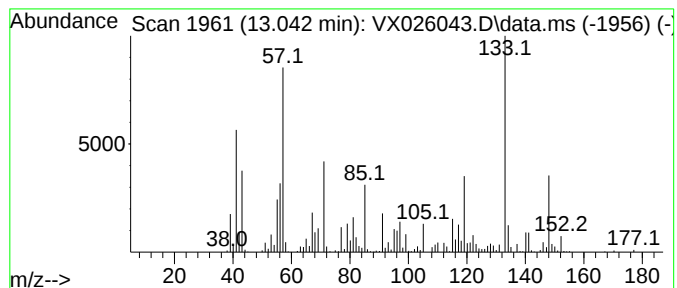
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 47 Benzene, pentamethyl- Concentration Rank 38

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 13.042 | 30.29 ug/L | 729037 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                                     | MW  | MolForm | CAS#         | Qual |
|------|----|---|--------------------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Benzene, pentamethyl-                            | 148 | C11H16  | 000700-12-9  | 55   |
| 2    |    |   | Benzene, pentamethyl-                            | 148 | C11H16  | 000700-12-9  | 55   |
| 3    |    |   | 1,5,6,7-Tetramethylbicyclo[3.2.0]hepta-2,6-diene | 148 | C11H16  | 134329-46-7  | 50   |
| 4    |    |   | 3,4-Dimethylcumene                               | 148 | C11H16  | 1000370-34-1 | 50   |
| 5    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)-              | 148 | C11H16  | 004218-48-8  | 49   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

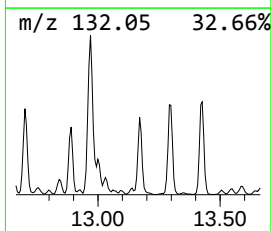
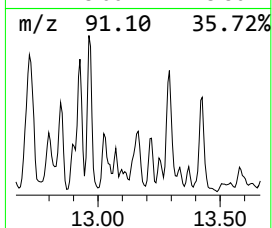
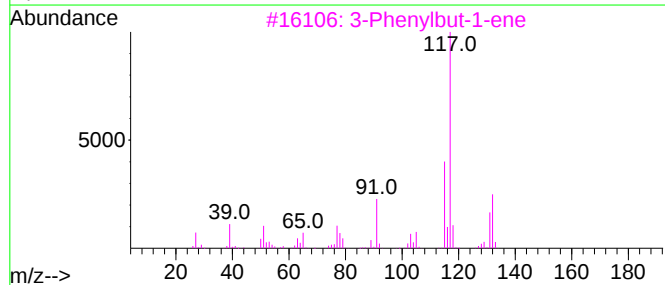
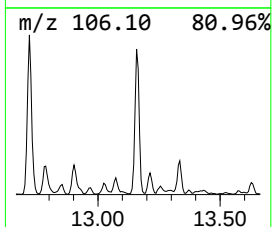
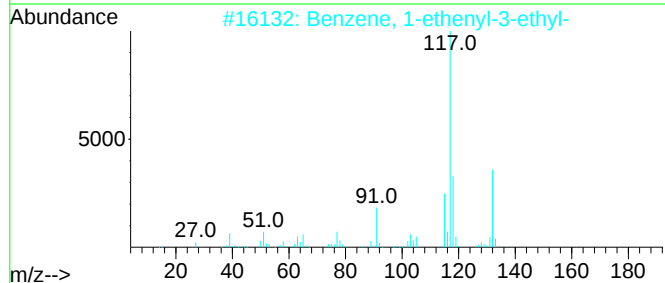
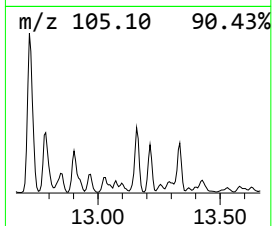
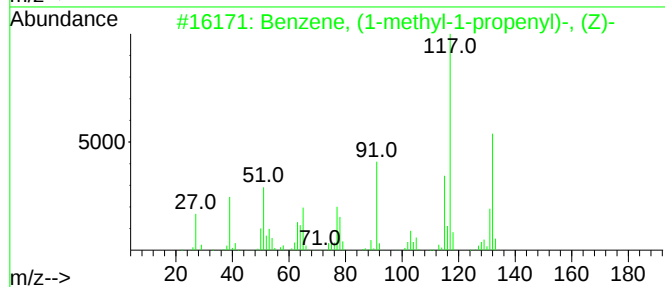
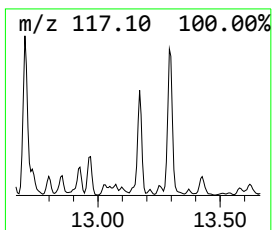
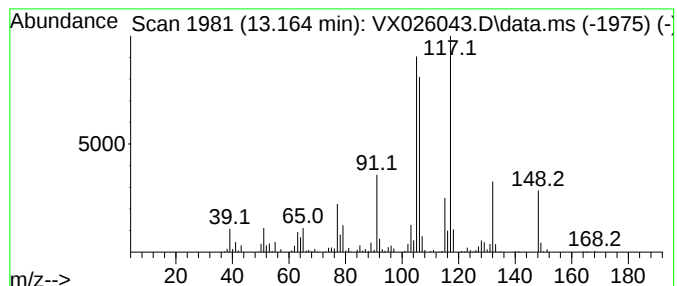
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 48 Benzene, (1-methyl-1-propen... Concentration Rank 36

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 13.164 | 35.13 ug/L | 845600 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000767-99-7 | 55   |
| 2    |    |   | Benzene, 1-ethenyl-3-ethyl-         | 132 | C10H12  | 007525-62-4 | 50   |
| 3    |    |   | 3-Phenylbut-1-ene                   | 132 | C10H12  | 000934-10-1 | 50   |
| 4    |    |   | Benzene, 2-butenyl-                 | 132 | C10H12  | 001560-06-1 | 50   |
| 5    |    |   | 1-Methyl-2-phenylcyclopropane       | 132 | C10H12  | 003145-76-4 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

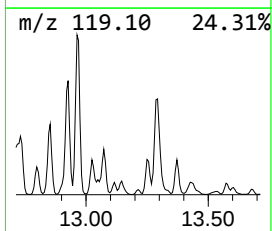
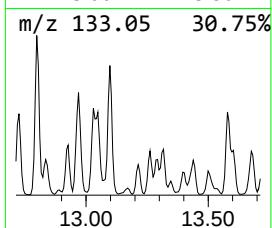
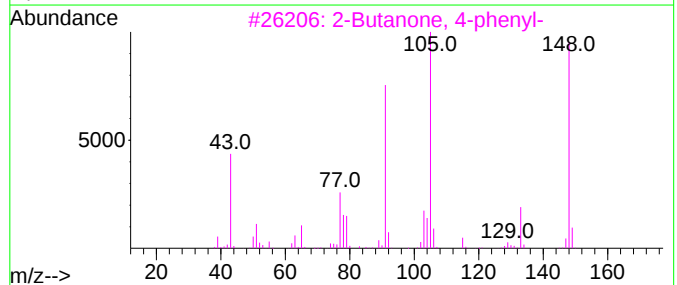
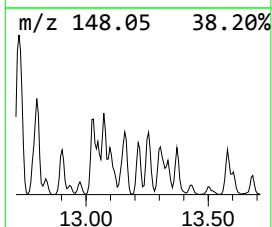
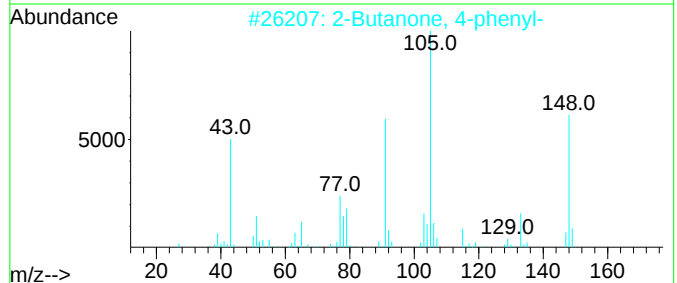
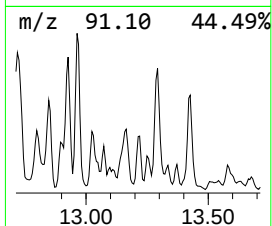
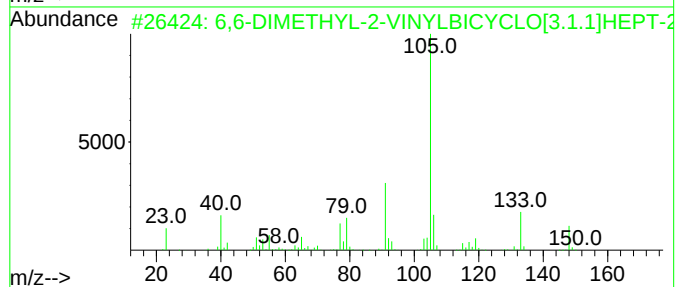
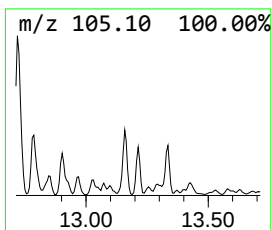
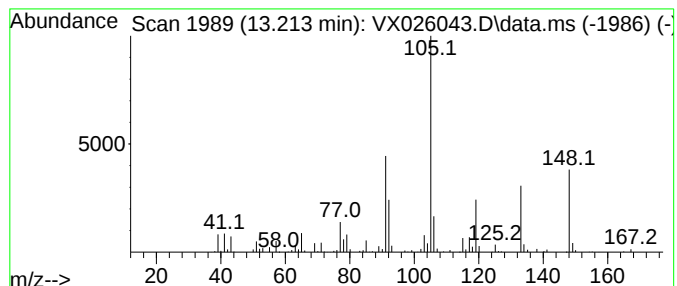
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 49 6,6-DIMETHYL-2-VINYLBICYCLO... Concentration Rank 50

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 13.213 | 10.04 ug/L | 241757 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 6,6-DIMETHYL-2-VINYLBICYCLO[3.1.... | 148 | C11H16       | 000473-00-7 | 59      |      |      |
| 2    | 2-Butanone, 4-phenyl-               | 148 | C10H12O      | 002550-26-7 | 43      |      |      |
| 3    | 2-Butanone, 4-phenyl-               | 148 | C10H12O      | 002550-26-7 | 43      |      |      |
| 4    | 2-Methylindan-2-ol                  | 148 | C10H12O      | 033223-84-6 | 38      |      |      |
| 5    | Benzene, (1,2-dimethylpropyl)-      | 148 | C11H16       | 004481-30-5 | 38      |      |      |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

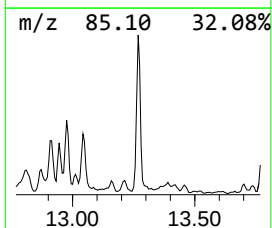
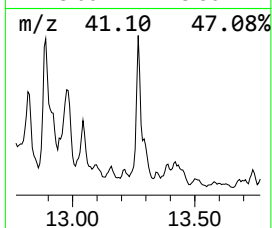
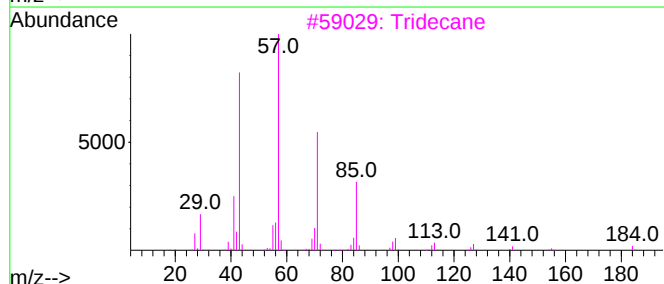
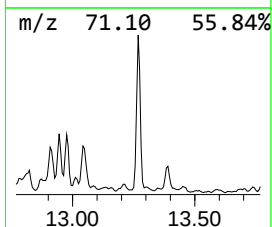
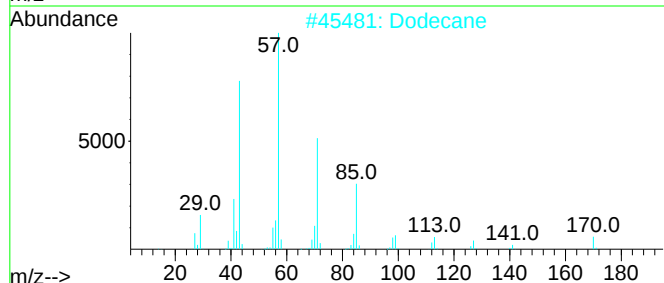
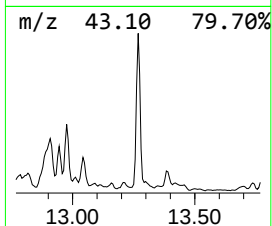
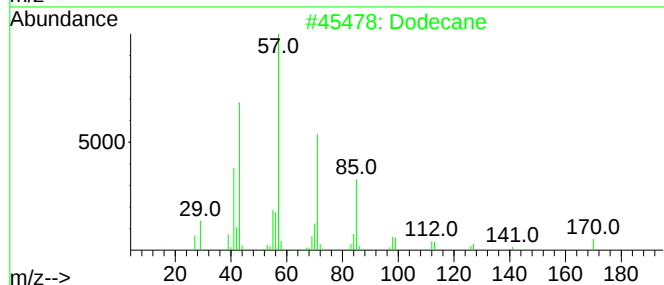
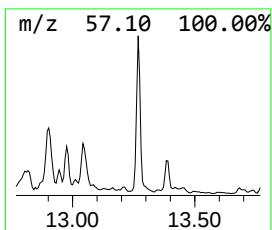
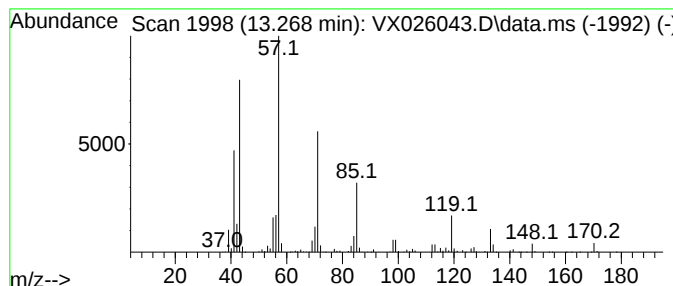
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 34

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 13.268 | 45.45 ug/L | 1094000 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane     | 170 | C12H26  | 000112-40-3 | 96   |
| 2    |    |   | Dodecane     | 170 | C12H26  | 000112-40-3 | 62   |
| 3    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 58   |
| 4    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 58   |
| 5    |    |   | Undecane     | 156 | C11H24  | 001120-21-4 | 58   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

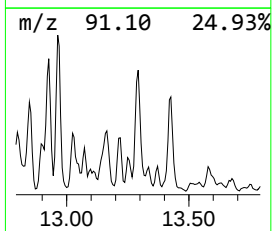
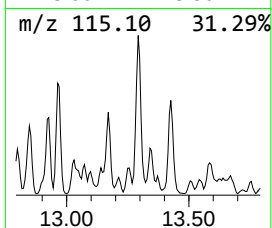
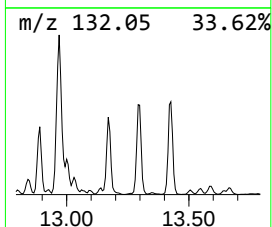
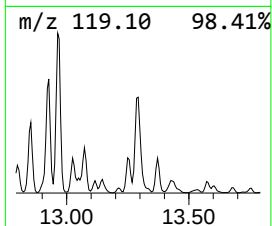
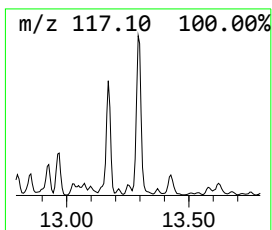
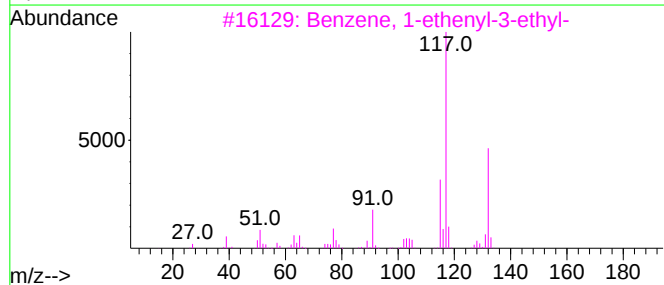
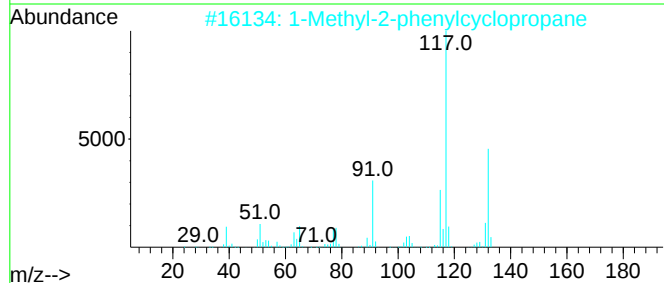
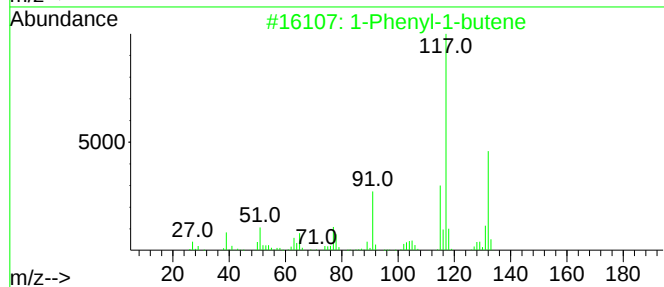
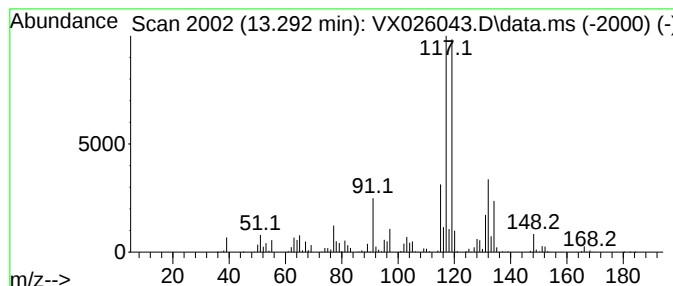
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 51 1-Phenyl-1-butene Concentration Rank 28

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 13.292 | 59.93 ug/L | 1442630 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Phenyl-1-butene                | 132 | C10H12  | 000824-90-8 | 91   |
| 2    |    |   | 1-Methyl-2-phenylcyclopropane    | 132 | C10H12  | 003145-76-4 | 87   |
| 3    |    |   | Benzene, 1-ethenyl-3-ethyl-      | 132 | C10H12  | 007525-62-4 | 86   |
| 4    |    |   | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12  | 002039-89-6 | 86   |
| 5    |    |   | Benzene, 2-butenyl-              | 132 | C10H12  | 001560-06-1 | 84   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

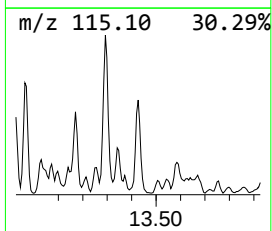
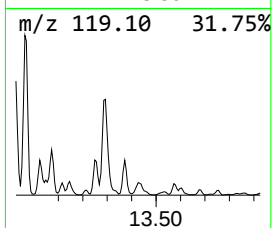
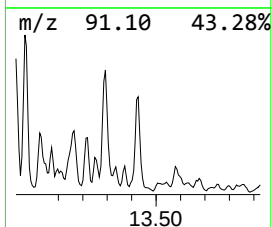
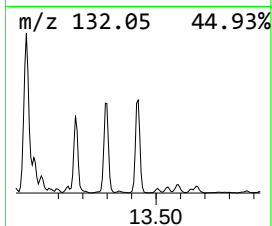
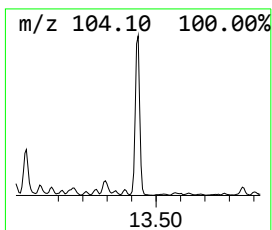
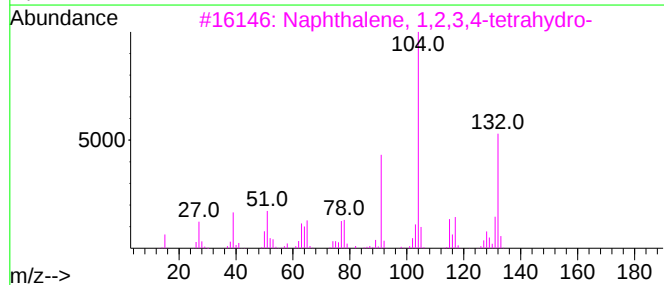
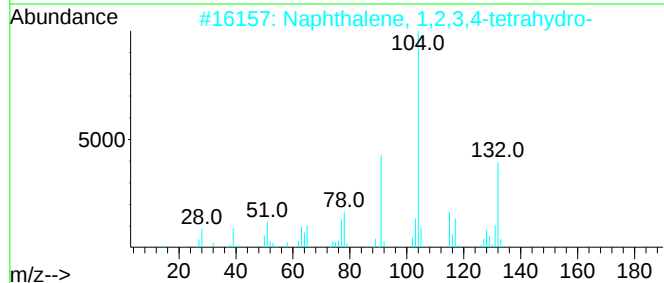
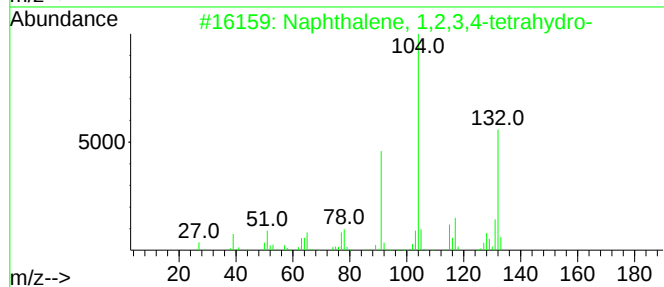
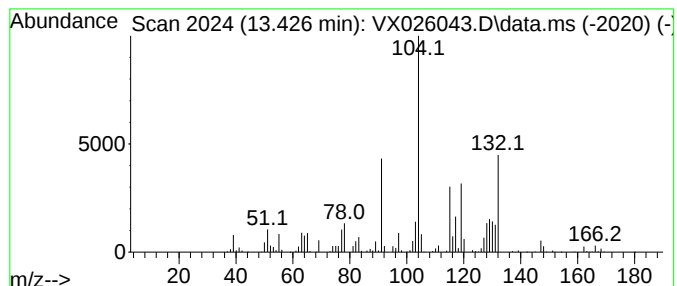
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 53 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 32

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 13.427 | 53.31 ug/L | 1283270 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 94   |
| 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 | 76   |
| 4    |    |   | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 76   |
| 5    |    |   | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 49   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

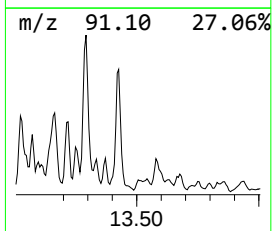
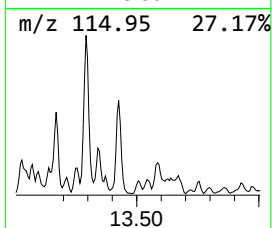
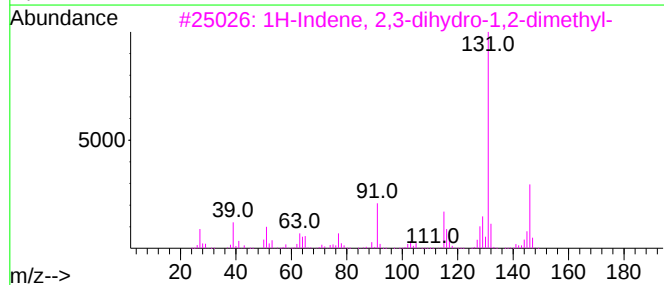
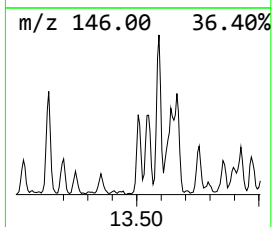
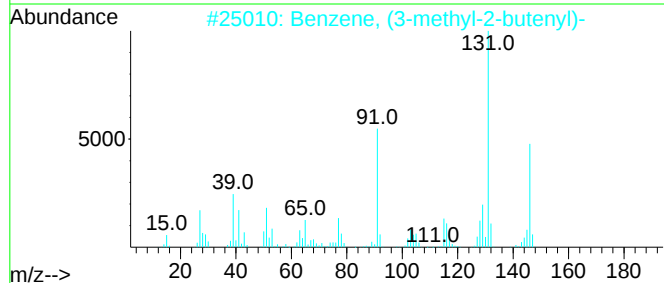
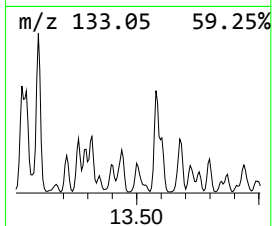
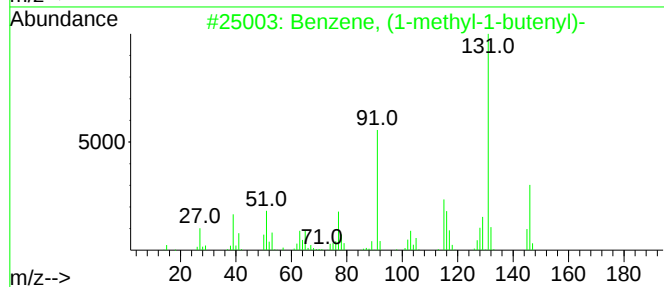
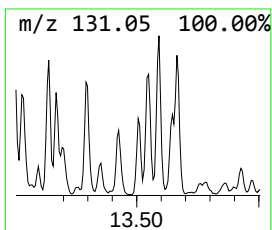
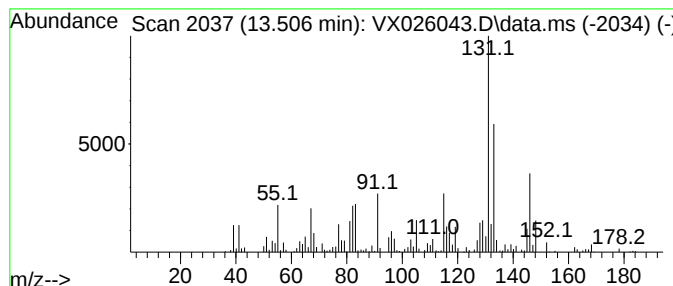
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 54 Benzene, (1-methyl-1-butenyl)- Concentration Rank 55

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.506 | 7.71 ug/L | 185473 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 86   |
| 2    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 86   |
| 3    |    |   | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 70   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146 | C11H14  | 004175-53-5 | 70   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 70   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

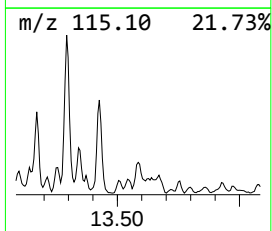
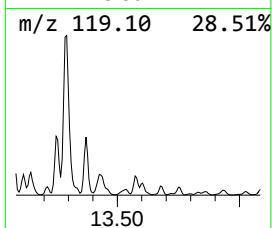
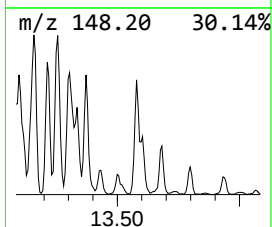
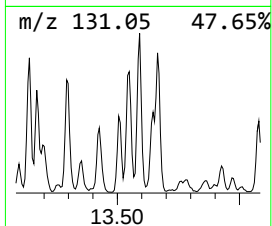
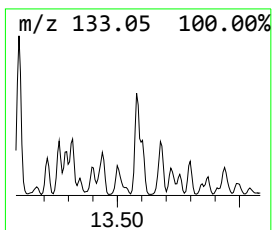
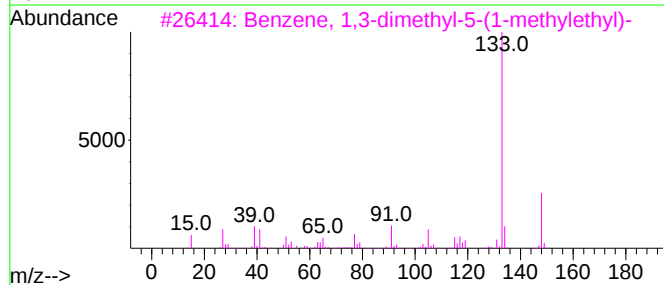
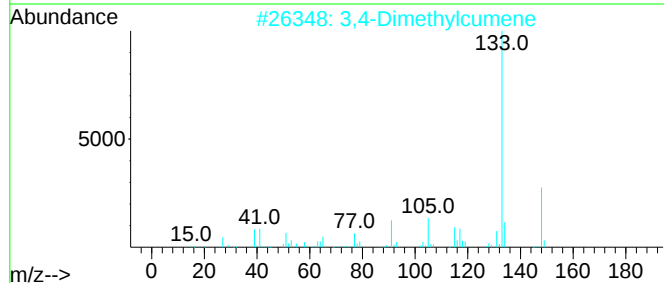
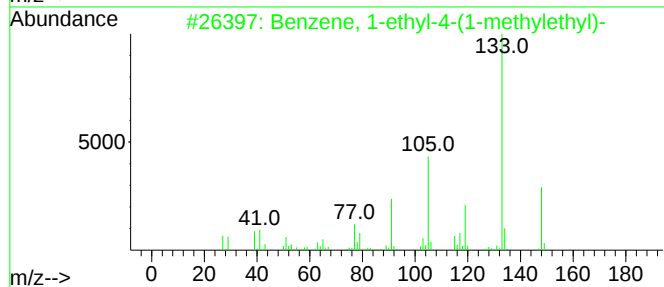
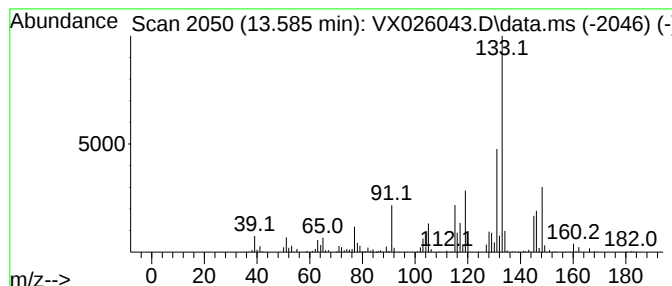
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 55 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 47

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 13.585 | 15.60 ug/L | 375398 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 | C11H16  | 004218-48-8  | 60   |
| 2    |    |   | 3,4-Dimethylcumene                  | 148 | C11H16  | 1000370-34-1 | 46   |
| 3    |    |   | Benzene, 1,3-dimethyl-5-(1-methy... | 148 | C11H16  | 004706-90-5  | 46   |
| 4    |    |   | Ethanone, 1-(4-ethylphenyl)-        | 148 | C10H12O | 000937-30-4  | 46   |
| 5    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 | C11H16  | 004218-48-8  | 46   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

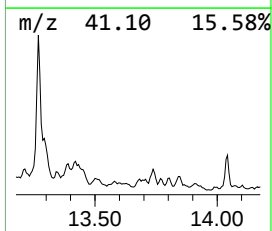
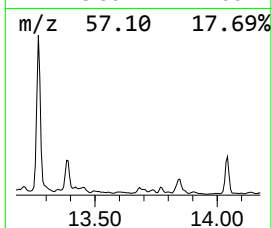
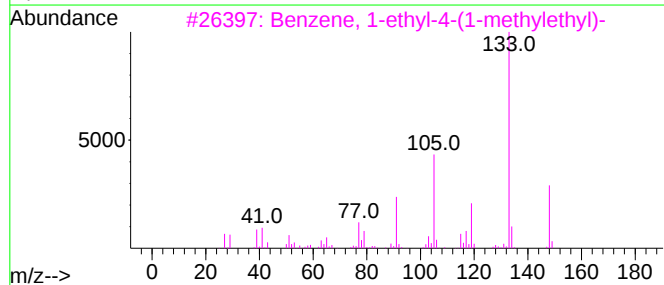
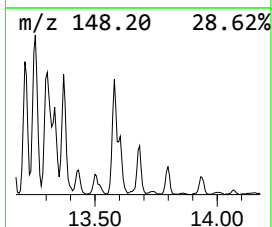
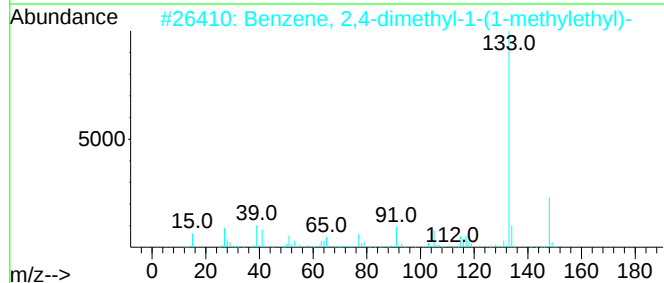
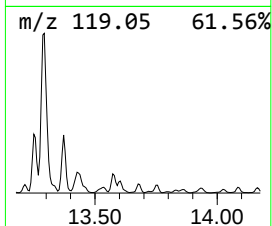
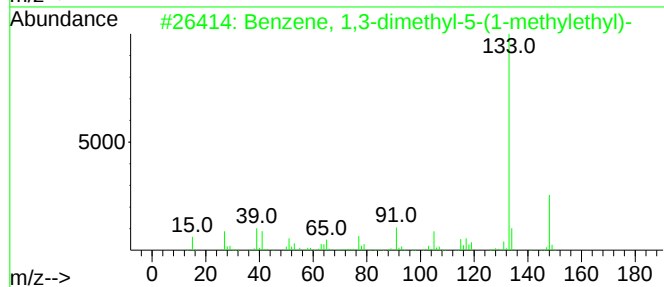
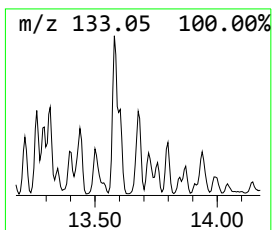
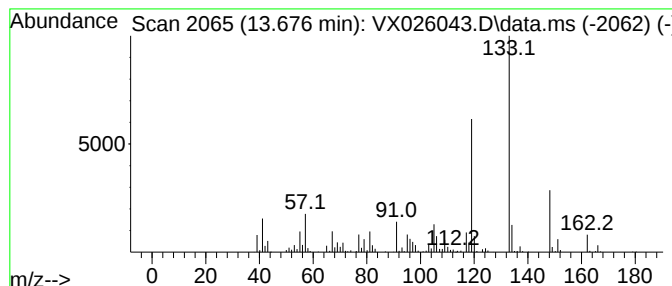
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 56 Benzene, 1,3-dimethyl-5-(1-... Concentration Rank 62

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.676 | 5.74 ug/L | 138083 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,3-dimethyl-5-(1-methy... | 148 | C11H16  | 004706-90-5 | 58   |
| 2    |    |   | Benzene, 2,4-dimethyl-1-(1-methy... | 148 | C11H16  | 004706-89-2 | 58   |
| 3    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 | C11H16  | 004218-48-8 | 58   |
| 4    |    |   | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9 | 53   |
| 5    |    |   | Benzene, 1,3-diethyl-5-methyl-      | 148 | C11H16  | 002050-24-0 | 52   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

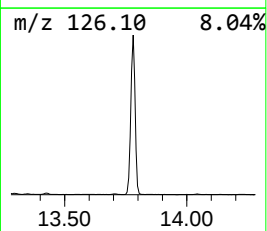
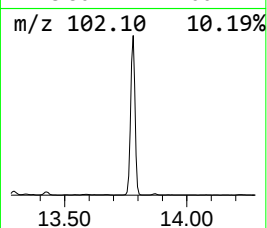
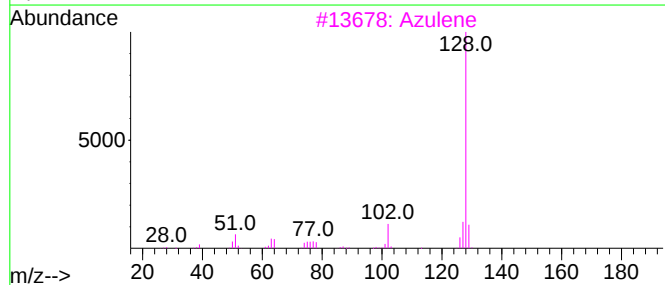
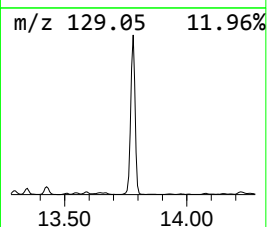
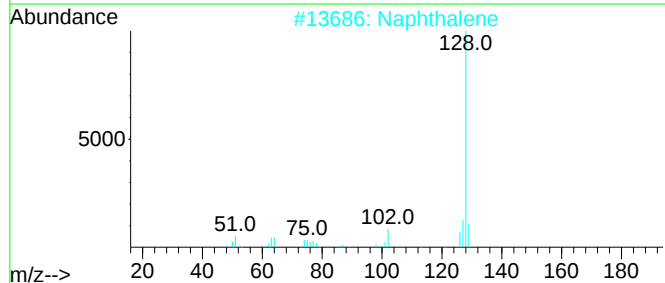
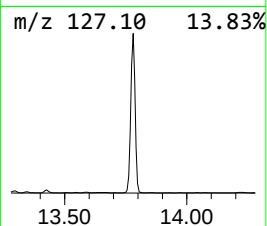
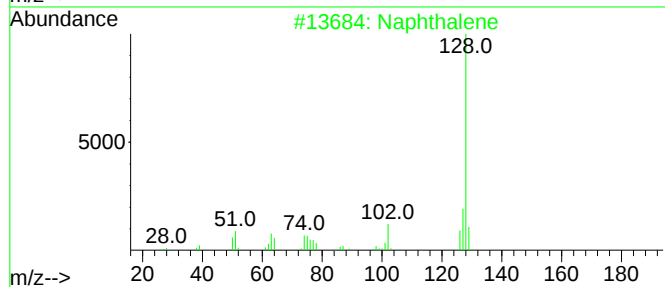
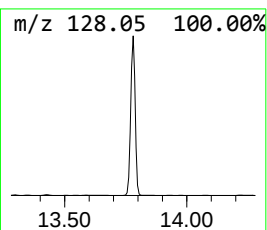
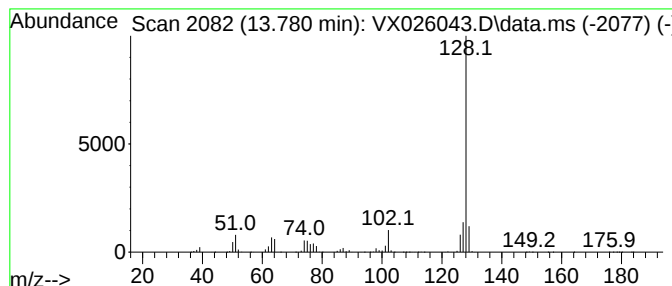
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 57 Azulene Concentration Rank 6

| R.T.   | EstConc     | Area    | Relative to ISTD       | R.T.   |
|--------|-------------|---------|------------------------|--------|
| 13.780 | 327.94 ug/L | 7894070 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene  | 128 | C10H8   | 000091-20-3 | 97   |
| 2    |    |   | Naphthalene  | 128 | C10H8   | 000091-20-3 | 95   |
| 3    |    |   | Azulene      | 128 | C10H8   | 000275-51-4 | 94   |
| 4    |    |   | Azulene      | 128 | C10H8   | 000275-51-4 | 91   |
| 5    |    |   | Azulene      | 128 | C10H8   | 000275-51-4 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

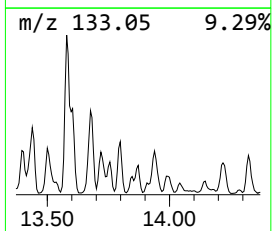
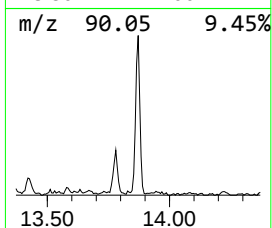
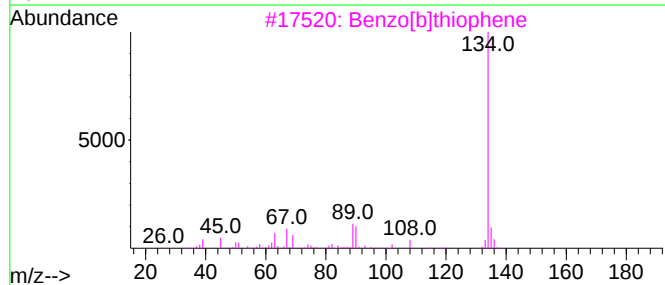
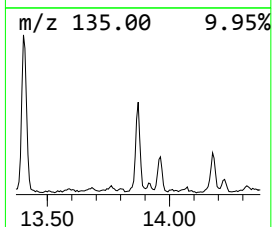
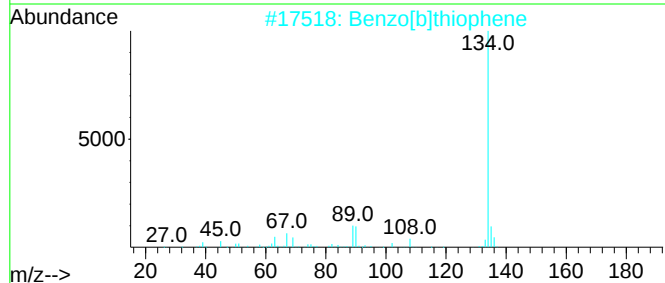
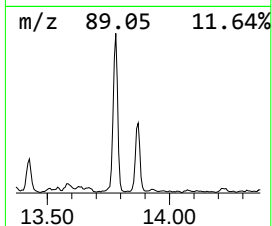
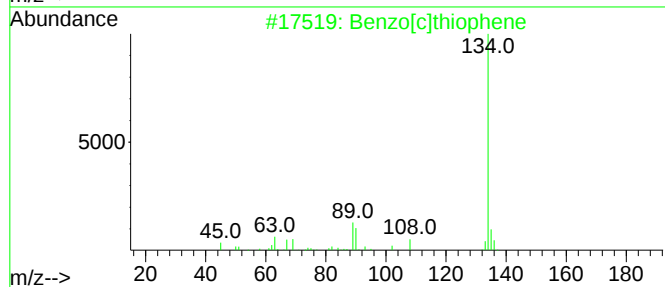
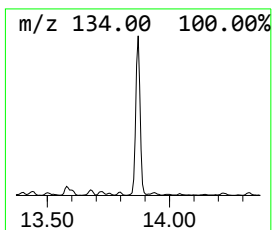
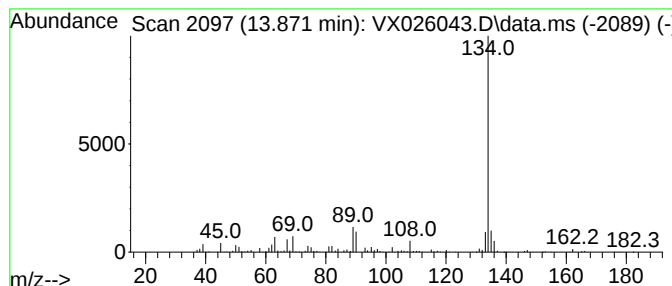
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 58 Benzo[c]thiophene Concentration Rank 46

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 13.871 | 16.79 ug/L | 404243 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzo[c]thiophene          | 134 | C8H6S   | 000270-82-6 | 94   |
| 2    |    |   | Benzo[b]thiophene          | 134 | C8H6S   | 000095-15-8 | 93   |
| 3    |    |   | Benzo[b]thiophene          | 134 | C8H6S   | 000095-15-8 | 90   |
| 4    |    |   | Cyclopenta[c]thiapyran     | 134 | C8H6S   | 000270-63-3 | 90   |
| 5    |    |   | 5-Phenyl-1,2,3-thiadiazole | 162 | C8H6N2S | 018212-29-8 | 56   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

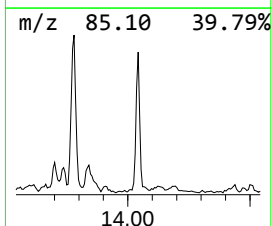
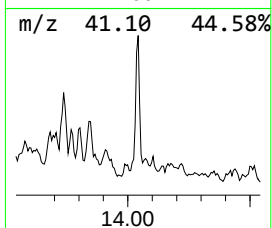
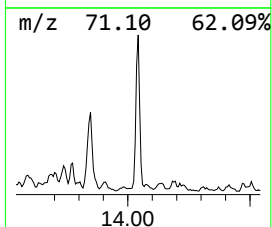
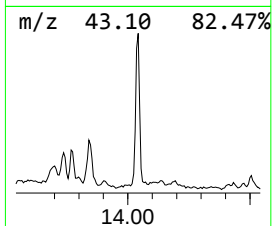
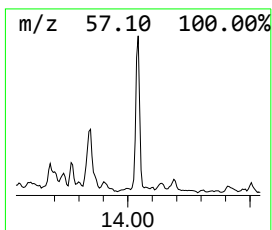
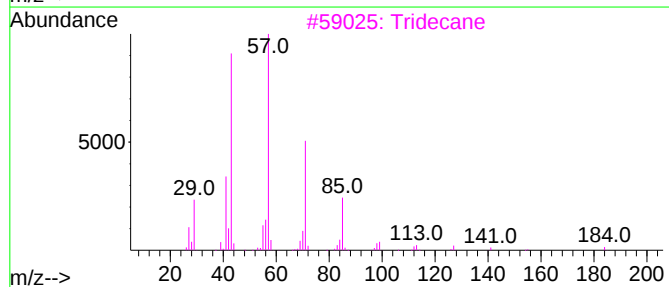
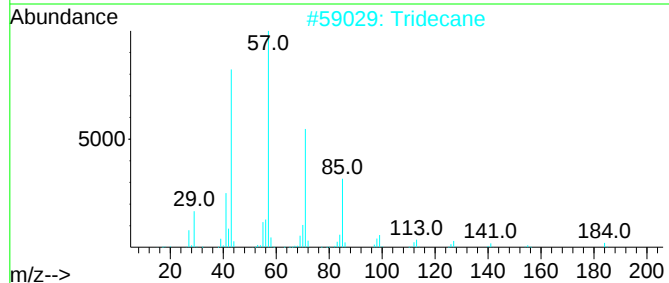
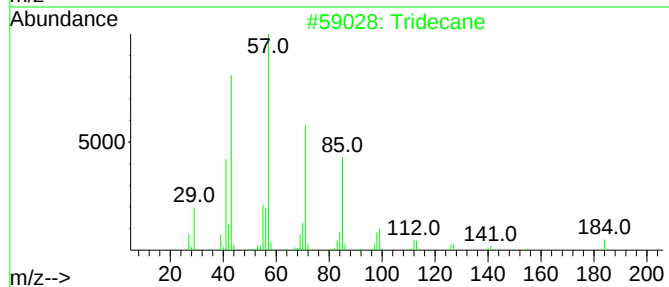
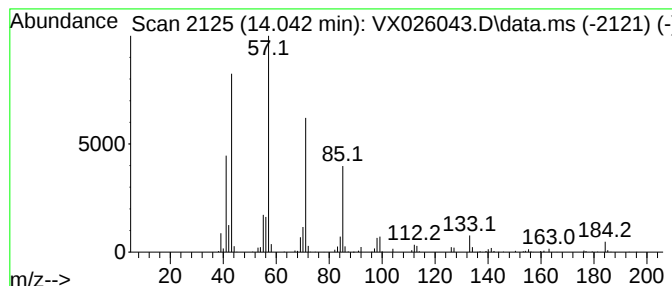
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 56

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 14.042 | 7.27 ug/L | 174946 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|------|-------------------------------------|-----|----------|--------------|------|
| 1    |      | Tridecane                           | 184 | C13H28   | 000629-50-5  | 97   |
| 2    |      | Tridecane                           | 184 | C13H28   | 000629-50-5  | 91   |
| 3    |      | Tridecane                           | 184 | C13H28   | 000629-50-5  | 90   |
| 4    |      | Hexadecane                          | 226 | C16H34   | 000544-76-3  | 87   |
| 5    |      | Carbonic acid, hexadecyl prop-1-... | 326 | C20H38O3 | 1000382-90-3 | 86   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

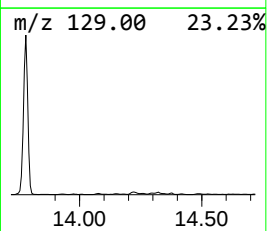
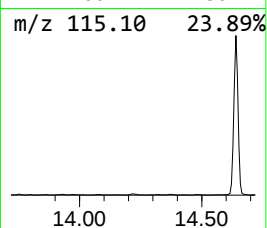
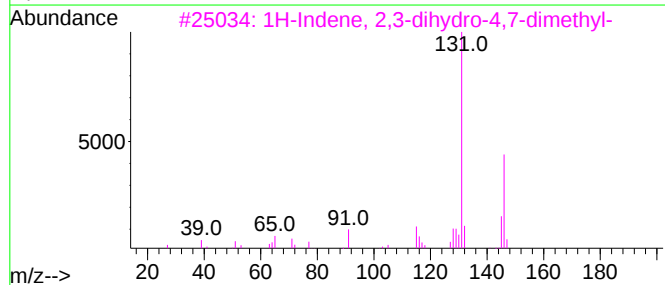
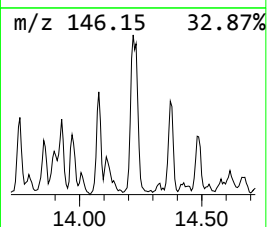
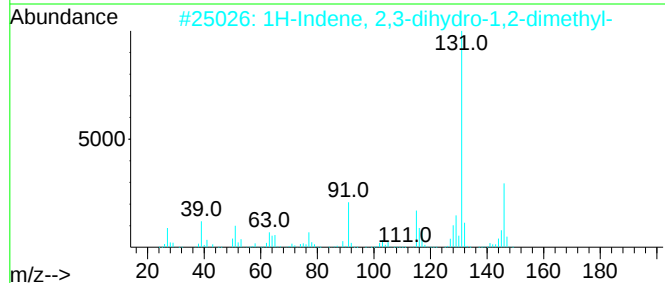
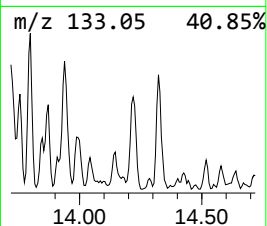
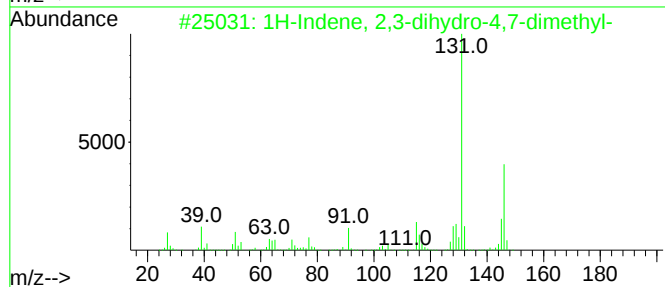
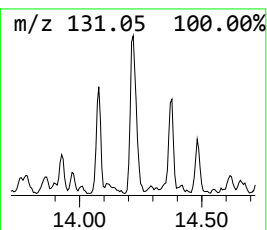
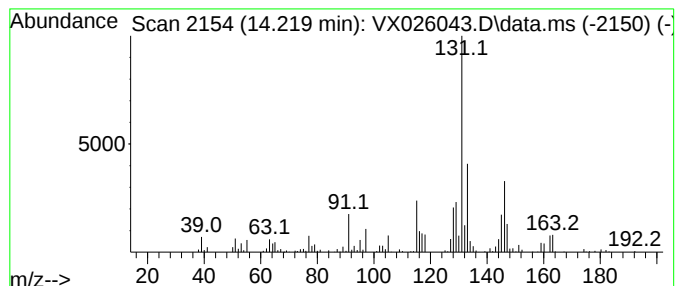
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 60 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 53

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 14.219 | 8.62 ug/L | 207570 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 64   |
| 2    |    |   | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 62   |
| 3    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 58   |
| 4    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 58   |
| 5    |    |   | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 | C11H14  | 000769-25-5 | 58   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

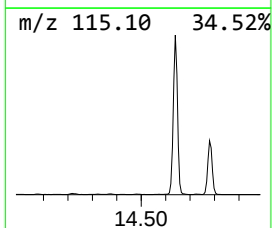
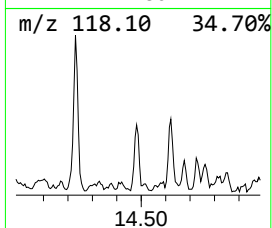
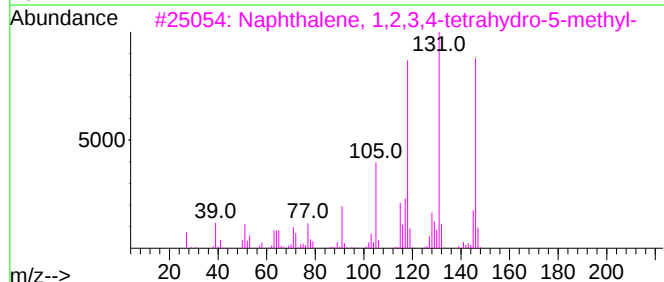
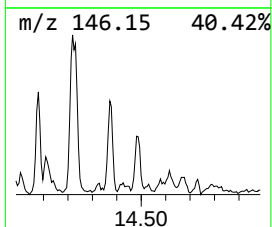
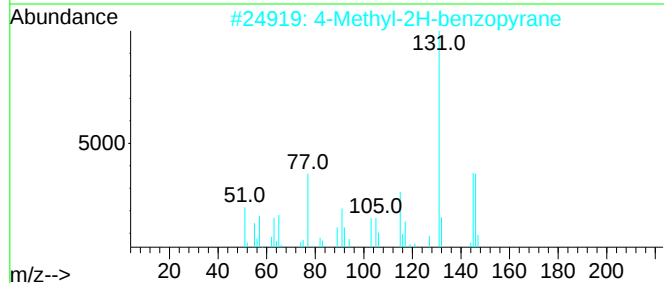
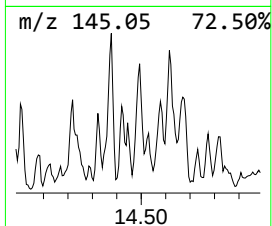
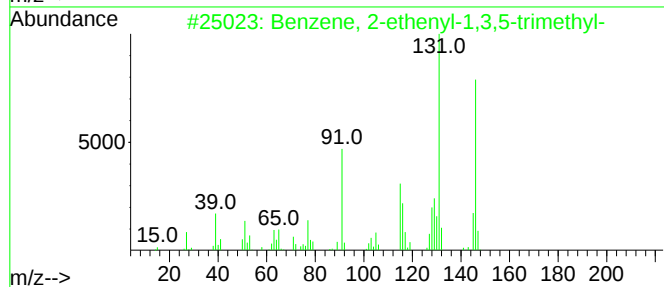
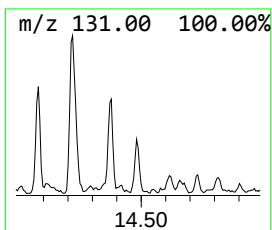
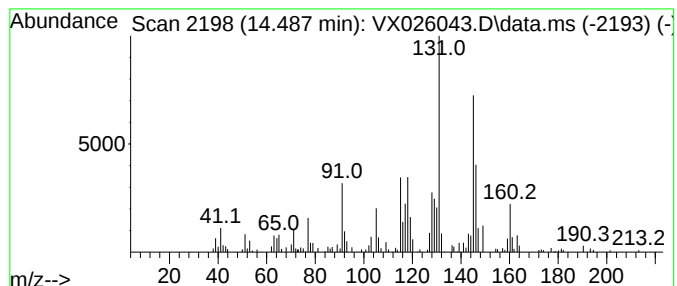
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 61 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 66

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 14.487 | 4.47 ug/L | 107710 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 | C11H14  | 000769-25-5 | 52   |
| 2    |    |   | 4-Methyl-2H-benzopyrane             | 146 | C10H100 | 021776-94-3 | 46   |
| 3    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5 | 46   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-5,6-dimet... | 146 | C11H14  | 001075-22-5 | 43   |
| 5    |    |   | Benzene, 4-(2-butenyl)-1,2-dimet... | 160 | C12H16  | 054340-86-2 | 43   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

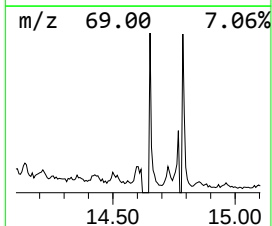
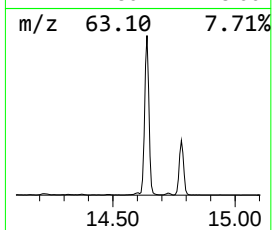
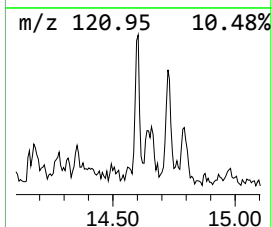
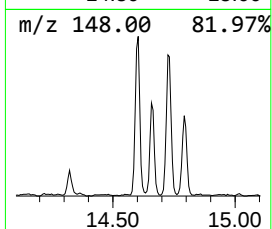
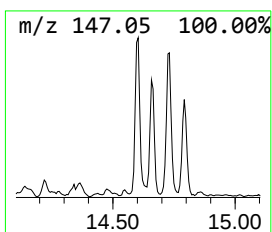
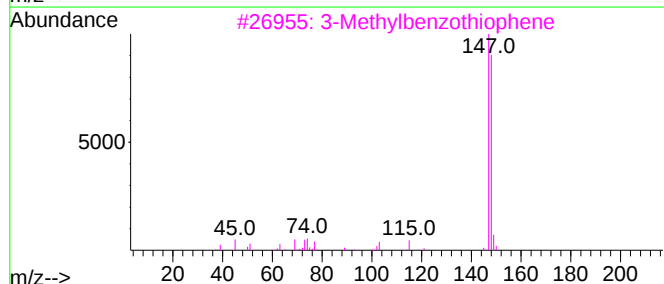
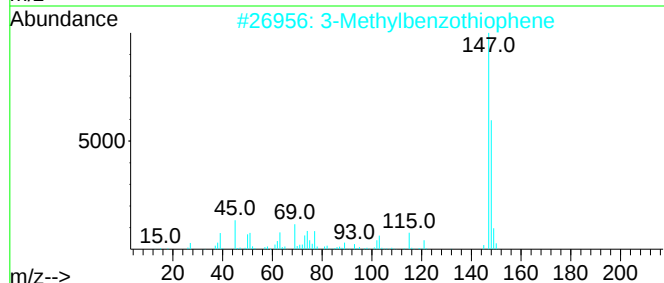
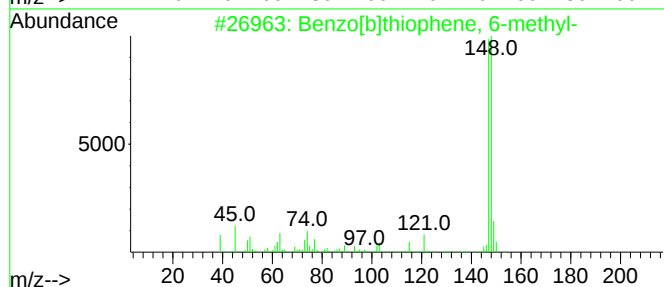
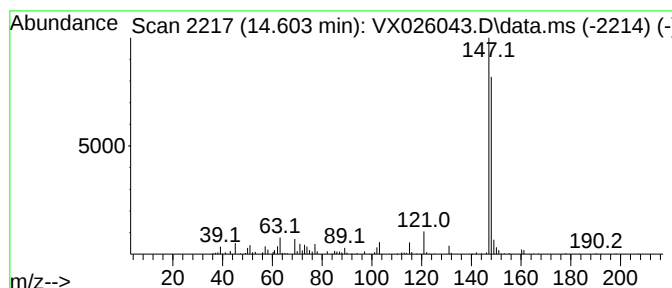
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 62 Benzo[b]thiophene, 6-methyl- Concentration Rank 69

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 14.603 | 3.14 ug/L | 75489 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzo[b]thiophene, 6-methyl-   | 148 | C9H8S   | 016587-47-6 | 86   |
| 2    |    |   | 3-Methylbenzothiophene         | 148 | C9H8S   | 001455-18-1 | 83   |
| 3    |    |   | 3-Methylbenzothiophene         | 148 | C9H8S   | 001455-18-1 | 83   |
| 4    |    |   | Benzaldehyde, 2,4,6-trimethyl- | 148 | C10H12O | 000487-68-3 | 72   |
| 5    |    |   | 2-Propenoic acid, 3-phenyl-    | 148 | C9H8O2  | 000621-82-9 | 72   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

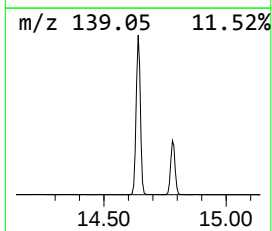
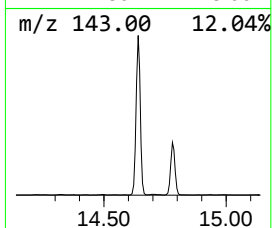
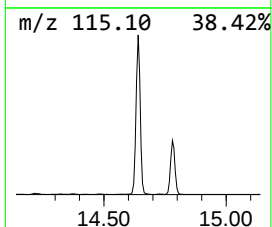
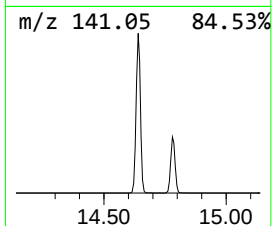
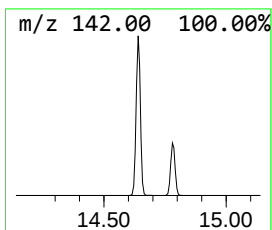
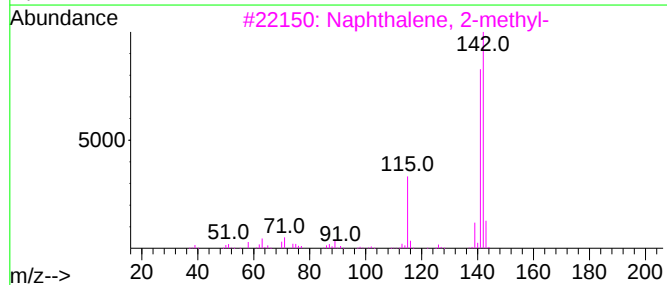
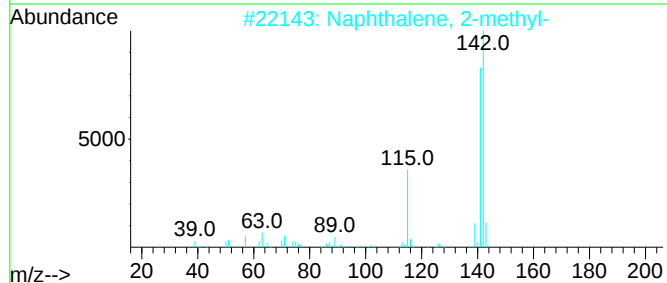
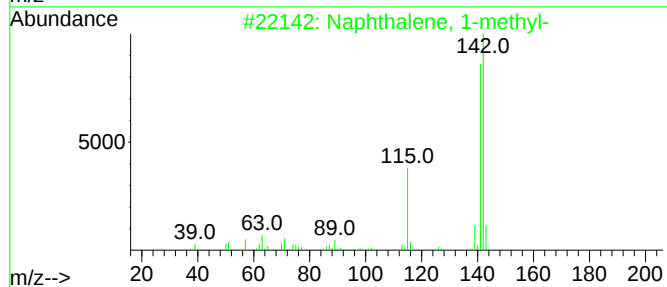
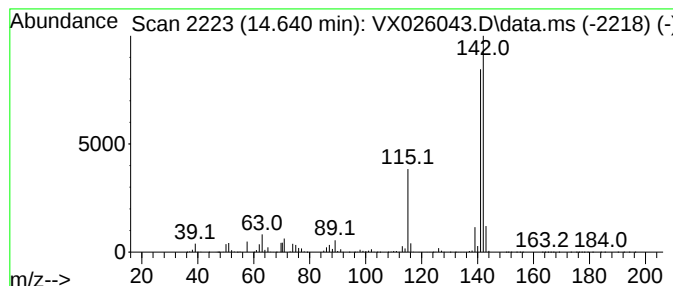
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 63 Naphthalene, 1-methyl- Concentration Rank 5

| R.T.   | EstConc     | Area    | Relative to ISTD       | R.T.   |
|--------|-------------|---------|------------------------|--------|
| 14.640 | 381.45 ug/L | 9182140 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 96   |
| 2    |    |   | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 96   |
| 3    |    |   | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 96   |
| 4    |    |   | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 94   |
| 5    |    |   | 1,4-Methanonaphthalene, 1,4-dihy... | 142 | C11H10  | 004453-90-1 | 94   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

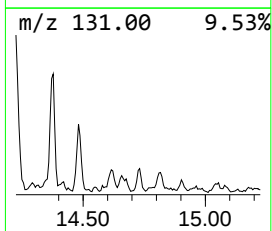
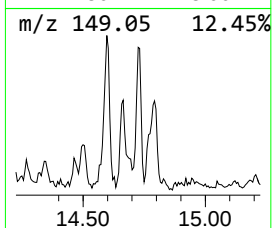
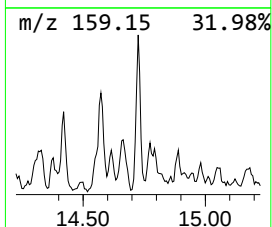
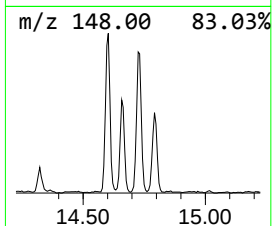
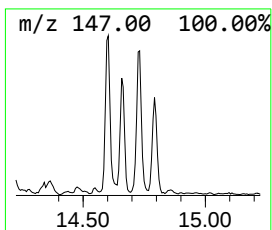
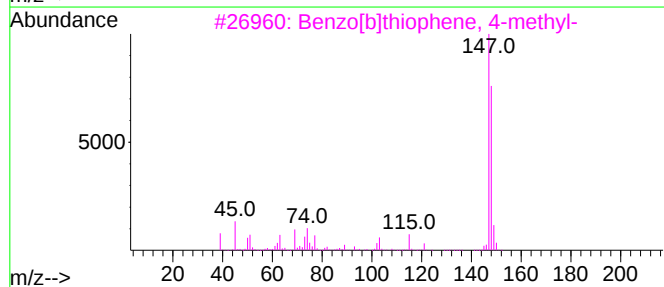
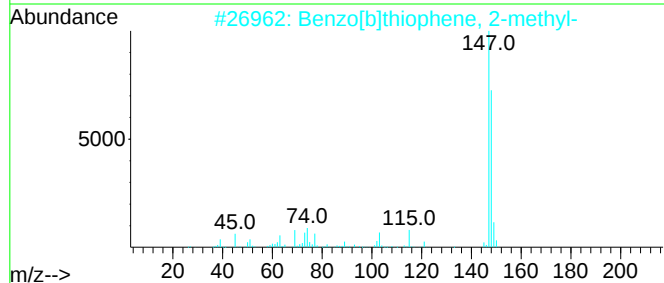
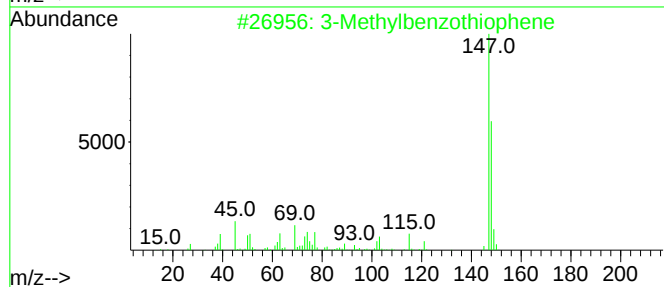
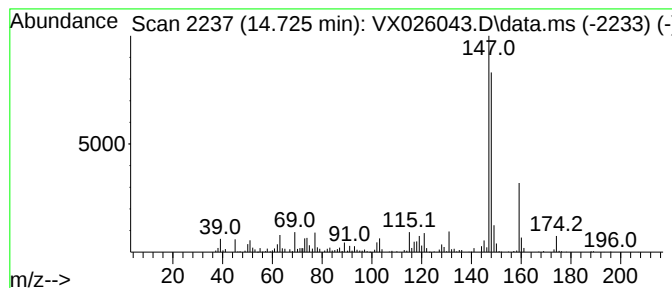
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXML122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 64 3-Methylbenzothiophene Concentration Rank 59

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 14.725 | 6.12 ug/L | 147389 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | 3-Methylbenzothiophene       | 148 | C9H8S   | 001455-18-1 | 76   |
| 2    |    |   | Benzo[b]thiophene, 2-methyl- | 148 | C9H8S   | 001195-14-8 | 76   |
| 3    |    |   | Benzo[b]thiophene, 4-methyl- | 148 | C9H8S   | 014315-11-8 | 76   |
| 4    |    |   | Benzo[b]thiophene, 5-methyl- | 148 | C9H8S   | 014315-14-1 | 76   |
| 5    |    |   | Benzo[b]thiophene, 2-methyl- | 148 | C9H8S   | 001195-14-8 | 76   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
Data File : VX026043.D  
Acq On : 25 Dec 2021 18:51  
Operator : JC/MD  
Sample : M5213-05 10X  
Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
BGLD7

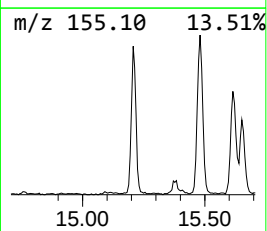
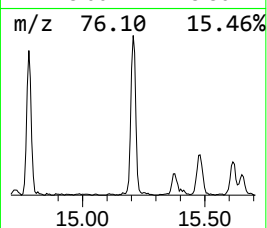
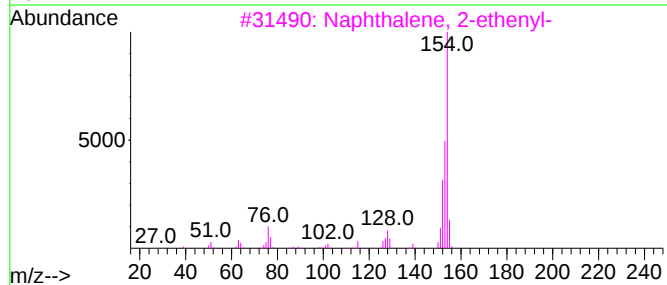
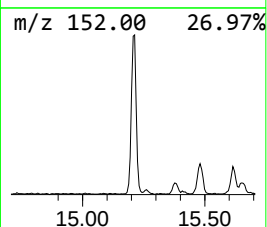
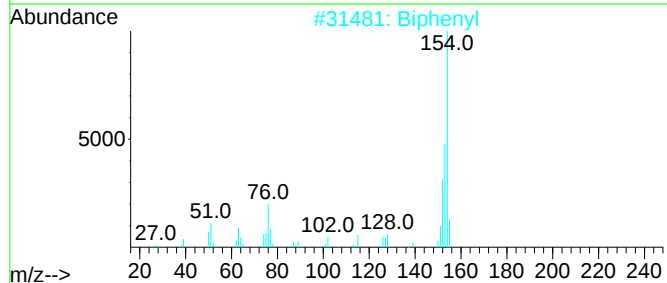
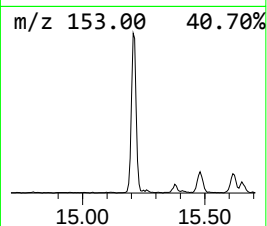
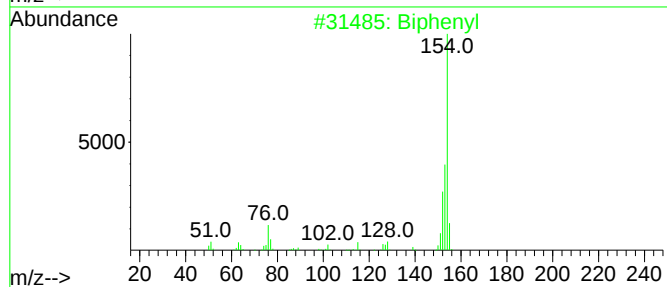
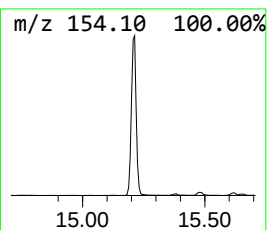
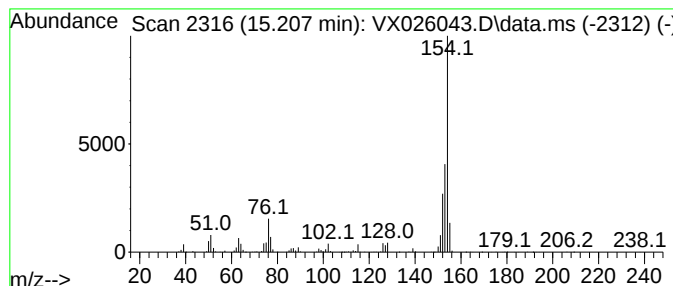
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXLM122321WMA.M  
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 66 Biphenyl Concentration Rank 45

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 15.207 | 17.70 ug/L | 426102 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Biphenyl                | 154 | C12H10  | 000092-52-4 | 95   |
| 2    |    |   | Biphenyl                | 154 | C12H10  | 000092-52-4 | 95   |
| 3    |    |   | Naphthalene, 2-ethenyl- | 154 | C12H10  | 000827-54-3 | 87   |
| 4    |    |   | Biphenyl                | 154 | C12H10  | 000092-52-4 | 87   |
| 5    |    |   | Naphthalene, 2-ethenyl- | 154 | C12H10  | 000827-54-3 | 81   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX122521\  
 Data File : VX026043.D  
 Acq On : 25 Dec 2021 18:51  
 Operator : JC/MD  
 Sample : M5213-05 10X  
 Misc : 6.57G/5.0mL/100uL/5.0mL/MSVOA\_X/MEOH  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 BGLD7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\SFAMXML122321WMA.M  
 Quant Title : VOC Analysis

TIC Library : C:\Database\NIST0.L  
 TIC Integration Parameters: LSCINT.P

|                    |        |         |       |          | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
| TIC Top Hit name   | RT     | EstConc | Units | Response | #                     | RT     | Resp    | Conc |
| (DEL) Alkane: S... | 6.611  | 3.1     | ug/L  | 36845    | 1                     | 6.763  | 601471  | 50.0 |
| (DEL) Alkane: S... | 8.275  | 5.9     | ug/L  | 70874    | 1                     | 6.763  | 601471  | 50.0 |
| (DEL) Alkane: S... | 8.440  | 4.2     | ug/L  | 64255    | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: C... | 8.598  | 6.8     | ug/L  | 103993   | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: C... | 9.104  | 3.4     | ug/L  | 52059    | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 9.293  | 4.8     | ug/L  | 73488    | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 9.500  | 19.8    | ug/L  | 300486   | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: C... | 9.561  | 18.5    | ug/L  | 281292   | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: C... | 9.622  | 26.7    | ug/L  | 405676   | 2                     | 10.055 | 758529  | 50.0 |
| 1-Undecanol        | 9.762  | 4.9     | ug/L  | 74360    | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 9.823  | 40.5    | ug/L  | 614082   | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 9.909  | 65.2    | ug/L  | 989047   | 2                     | 10.055 | 758529  | 50.0 |
| Pentalene, octa... | 10.122 | 9.9     | ug/L  | 150494   | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 10.366 | 384.5   | ug/L  | 5833720  | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: C... | 10.458 | 21.7    | ug/L  | 329780   | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 10.537 | 11.5    | ug/L  | 174409   | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: C... | 10.592 | 125.3   | ug/L  | 1900370  | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 10.714 | 24.4    | ug/L  | 369966   | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 10.781 | 244.2   | ug/L  | 3705170  | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: C... | 10.860 | 251.4   | ug/L  | 3814070  | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 10.896 | 103.3   | ug/L  | 1566760  | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 11.031 | 67.6    | ug/L  | 1025400  | 2                     | 10.055 | 758529  | 50.0 |
| (DEL) Alkane: S... | 11.085 | 104.8   | ug/L  | 2523560  | 3                     | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: S... | 11.110 | 140.9   | ug/L  | 3391540  | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, propyl-   | 11.305 | 151.8   | ug/L  | 3654330  | 3                     | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: C... | 11.342 | 32.1    | ug/L  | 773879   | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, 1-ethy... | 11.384 | 713.1   | ug/L  | 17166800 | 3                     | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: C... | 11.579 | 8.1     | ug/L  | 194309   | 3                     | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: S... | 11.689 | 48.6    | ug/L  | 1170190  | 3                     | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: S... | 11.841 | 53.4    | ug/L  | 1284660  | 3                     | 12.024 | 1203580 | 50.0 |
| Oxirane, 2-meth... | 11.896 | 68.0    | ug/L  | 1636040  | 3                     | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: C... | 11.945 | 175.1   | ug/L  | 4214570  | 3                     | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: S... | 12.177 | 88.3    | ug/L  | 2124920  | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, cyclop... | 12.250 | 401.5   | ug/L  | 9663960  | 3                     | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: S... | 12.427 | 272.6   | ug/L  | 6561160  | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, 1-meth... | 12.469 | 109.3   | ug/L  | 2631060  | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, 2-ethy... | 12.536 | 73.0    | ug/L  | 1757100  | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, 1-meth... | 12.561 | 104.4   | ug/L  | 2512340  | 3                     | 12.024 | 1203580 | 50.0 |
| o-Cymene           | 12.616 | 75.8    | ug/L  | 1824020  | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, 1,2-di... | 12.719 | 89.4    | ug/L  | 2152530  | 3                     | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: C... | 12.890 | 59.7    | ug/L  | 1437290  | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, 1,2,4,... | 12.921 | 54.8    | ug/L  | 1319970  | 3                     | 12.024 | 1203580 | 50.0 |
| 1,3-Cyclopentad... | 12.969 | 113.8   | ug/L  | 2739050  | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, pentam... | 13.042 | 30.3    | ug/L  | 729037   | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, (1-met... | 13.164 | 35.1    | ug/L  | 845600   | 3                     | 12.024 | 1203580 | 50.0 |
| 6,6-DIMETHYL-2-... | 13.213 | 10.0    | ug/L  | 241757   | 3                     | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: S... | 13.268 | 45.5    | ug/L  | 1094000  | 3                     | 12.024 | 1203580 | 50.0 |
| 1-Phenyl-1-butene  | 13.292 | 59.9    | ug/L  | 1442630  | 3                     | 12.024 | 1203580 | 50.0 |
| Naphthalene, 1,... | 13.427 | 53.3    | ug/L  | 1283270  | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, (1-met... | 13.506 | 7.7     | ug/L  | 185473   | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, 1-ethy... | 13.585 | 15.6    | ug/L  | 375398   | 3                     | 12.024 | 1203580 | 50.0 |
| Benzene, 1,3-di... | 13.676 | 5.7     | ug/L  | 138083   | 3                     | 12.024 | 1203580 | 50.0 |

|                    |        |       |      |         |   |        |         |      |
|--------------------|--------|-------|------|---------|---|--------|---------|------|
| Azulene            | 13.780 | 327.9 | ug/L | 7894070 | 3 | 12.024 | 1203580 | 50.0 |
| Benzo[c]thiophene  | 13.871 | 16.8  | ug/L | 404243  | 3 | 12.024 | 1203580 | 50.0 |
| (DEL) Alkane: S... | 14.042 | 7.3   | ug/L | 174946  | 3 | 12.024 | 1203580 | 50.0 |
| 1H-Indene, 2,3-... | 14.219 | 8.6   | ug/L | 207570  | 3 | 12.024 | 1203580 | 50.0 |
| Benzene, 2-ethe... | 14.487 | 4.5   | ug/L | 107710  | 3 | 12.024 | 1203580 | 50.0 |
| Benzo[b]thiophe... | 14.603 | 3.1   | ug/L | 75489   | 3 | 12.024 | 1203580 | 50.0 |
| Naphthalene, 1-... | 14.640 | 381.4 | ug/L | 9182140 | 3 | 12.024 | 1203580 | 50.0 |
| 3-Methylbenzoth... | 14.725 | 6.1   | ug/L | 147389  | 3 | 12.024 | 1203580 | 50.0 |
| Biphenyl           | 15.207 | 17.7  | ug/L | 426102  | 3 | 12.024 | 1203580 | 50.0 |

Instrument :

MSVOA\_X

ClientSampleId :

BGLD7