

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
 Data File : VW021852.D  
 Acq On : 11 Jan 2022 23:20  
 Operator : SY/VA  
 Sample : N1084-10  
 Misc : 6.57g/10mL/MSVOA\_W/SOIL  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_W  
 ClientSampleId :  
 BGLN3

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\_W\Method\SFAMWLM011122SMA.M  
 Title : SFAM01.0

Signal : TIC: VW021852.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         | 3.026       | 175           | 184         | 196          | rVB2     | 68448          | 159459        | 0.18%           | 0.015%        |
| 2         | 3.385       | 235           | 243         | 259          | rVB      | 73291          | 179967        | 0.20%           | 0.017%        |
| 3         | 4.019       | 336           | 347         | 356          | rBV      | 78111          | 226336        | 0.25%           | 0.022%        |
| 4         | 4.964       | 479           | 502         | 519          | rBV      | 162016         | 715506        | 0.80%           | 0.068%        |
| 5         | 5.440       | 567           | 580         | 598          | rBV2     | 158073         | 544829        | 0.61%           | 0.052%        |
| 6         | 5.940       | 650           | 662         | 678          | rBV      | 355096         | 1072414       | 1.20%           | 0.102%        |
| 7         | 6.988       | 823           | 834         | 845          | rVV      | 290024         | 863296        | 0.97%           | 0.082%        |
| 8         | 7.653       | 931           | 943         | 949          | rBV      | 110822         | 297551        | 0.33%           | 0.028%        |
| 9         | 7.927       | 979           | 988         | 1000         | rVV4     | 149775         | 545288        | 0.61%           | 0.052%        |
| 10        | 8.159       | 1016          | 1026        | 1034         | rBV      | 261409         | 601585        | 0.67%           | 0.057%        |
| 11        | 8.329       | 1038          | 1054        | 1064         | rVV      | 12763904       | 26874708      | 30.11%          | 2.559%        |
| 12        | 8.433       | 1064          | 1071        | 1075         | rVV3     | 66006          | 184204        | 0.21%           | 0.018%        |
| 13        | 8.488       | 1075          | 1080        | 1089         | rVV3     | 68149          | 239438        | 0.27%           | 0.023%        |
| 14        | 8.695       | 1104          | 1114        | 1123         | rBV      | 444618         | 863839        | 0.97%           | 0.082%        |
| 15        | 8.842       | 1128          | 1138        | 1150         | rBV      | 226339         | 483139        | 0.54%           | 0.046%        |
| 16        | 9.281       | 1202          | 1210        | 1213         | rBV2     | 135000         | 311332        | 0.35%           | 0.030%        |
| 17        | 9.335       | 1213          | 1219        | 1226         | rVB      | 541948         | 1100510       | 1.23%           | 0.105%        |
| 18        | 9.957       | 1316          | 1321        | 1325         | rBV      | 210586         | 379886        | 0.43%           | 0.036%        |
| 19        | 10.097      | 1338          | 1344        | 1357         | rVB      | 263103         | 591973        | 0.66%           | 0.056%        |
| 20        | 10.323      | 1374          | 1381        | 1385         | rBV      | 762511         | 1430135       | 1.60%           | 0.136%        |
| 21        | 10.390      | 1385          | 1392        | 1400         | rVB      | 19508054       | 31534253      | 35.33%          | 3.003%        |
| 22        | 10.506      | 1404          | 1411        | 1418         | rVB2     | 570983         | 1034151       | 1.16%           | 0.098%        |
| 23        | 10.670      | 1432          | 1438        | 1445         | rVB      | 336080         | 638517        | 0.72%           | 0.061%        |
| 24        | 10.762      | 1445          | 1453        | 1458         | rBV      | 355090         | 752599        | 0.84%           | 0.072%        |
| 25        | 10.847      | 1463          | 1467        | 1473         | rVB      | 195285         | 337028        | 0.38%           | 0.032%        |
| 26        | 10.939      | 1473          | 1482        | 1488         | rBV2     | 449610         | 912776        | 1.02%           | 0.087%        |
| 27        | 11.042      | 1488          | 1499        | 1506         | rBV      | 417369         | 993184        | 1.11%           | 0.095%        |
| 28        | 11.116      | 1506          | 1511        | 1517         | rBV      | 578786         | 1354569       | 1.52%           | 0.129%        |
| 29        | 11.183      | 1517          | 1522        | 1526         | rVB      | 654408         | 1081184       | 1.21%           | 0.103%        |
| 30        | 11.237      | 1526          | 1531        | 1536         | rVB      | 617163         | 1059770       | 1.19%           | 0.101%        |
| 31        | 11.286      | 1536          | 1539        | 1543         | rVB2     | 151388         | 214810        | 0.24%           | 0.020%        |
| 32        | 11.347      | 1543          | 1549        | 1552         | rBV      | 569858         | 1046640       | 1.17%           | 0.100%        |
| 33        | 11.420      | 1552          | 1561        | 1572         | rVB5     | 2232663        | 6818498       | 7.64%           | 0.649%        |
| 34        | 11.524      | 1572          | 1578        | 1590         | rVB      | 1673828        | 3716141       | 4.16%           | 0.354%        |
| 35        | 11.628      | 1590          | 1595        | 1605         | rBV3     | 299158         | 826306        | 0.93%           | 0.079%        |
| 36        | 11.731      | 1605          | 1612        | 1614         | rBV3     | 37910834       | 61193349      | 68.56%          | 5.827%        |

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 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_W  
 ClientSampleId :  
 BGLN3

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\_W\Method\SFAMWLM011122SMA.M  
 Title : SFAM01.0

|    |        |      |      |      |      |          |          |         |        |
|----|--------|------|------|------|------|----------|----------|---------|--------|
| 37 | 11.835 | 1624 | 1629 | 1631 | rBV2 | 39551368 | 59137517 | 66.25%  | 5.631% |
| 38 | 11.939 | 1644 | 1646 | 1650 | rVB  | 1650169  | 2004814  | 2.25%   | 0.191% |
| 39 | 11.981 | 1650 | 1653 | 1655 | rBV2 | 198775   | 271132   | 0.30%   | 0.026% |
| 40 | 12.012 | 1655 | 1658 | 1661 | rVB  | 208620   | 246736   | 0.28%   | 0.023% |
| 41 | 12.067 | 1661 | 1667 | 1672 | rBV3 | 2043815  | 3713443  | 4.16%   | 0.354% |
| 42 | 12.170 | 1672 | 1684 | 1691 | rBV  | 32690098 | 68050111 | 76.24%  | 6.480% |
| 43 | 12.256 | 1692 | 1698 | 1703 | rVB  | 5592208  | 9187027  | 10.29%  | 0.875% |
| 44 | 12.341 | 1703 | 1712 | 1714 | rBV5 | 3533613  | 9090944  | 10.18%  | 0.866% |
| 45 | 12.384 | 1714 | 1719 | 1727 | rVV4 | 4397639  | 13062170 | 14.63%  | 1.244% |
| 46 | 12.469 | 1727 | 1733 | 1736 | rVV2 | 5032082  | 10116978 | 11.33%  | 0.963% |
| 47 | 12.518 | 1736 | 1741 | 1742 | rVV2 | 4204957  | 8141975  | 9.12%   | 0.775% |
| 48 | 12.548 | 1743 | 1746 | 1748 | rVV  | 7274313  | 11448102 | 12.83%  | 1.090% |
| 49 | 12.573 | 1748 | 1750 | 1757 | rVV2 | 7821538  | 12936641 | 14.49%  | 1.232% |
| 50 | 12.658 | 1760 | 1764 | 1768 | rVV  | 6689930  | 9909280  | 11.10%  | 0.944% |
| 51 | 12.707 | 1768 | 1772 | 1776 | rVV4 | 1708866  | 3332759  | 3.73%   | 0.317% |
| 52 | 12.804 | 1777 | 1788 | 1793 | rVB  | 12984337 | 27538716 | 30.85%  | 2.622% |
| 53 | 12.871 | 1793 | 1799 | 1803 | rBV2 | 37126214 | 76783337 | 86.02%  | 7.311% |
| 54 | 12.908 | 1803 | 1805 | 1807 | rVV  | 22743354 | 28394862 | 31.81%  | 2.704% |
| 55 | 12.951 | 1807 | 1812 | 1817 | rVV3 | 39541591 | 89259642 | 100.00% | 8.499% |
| 56 | 12.999 | 1817 | 1820 | 1825 | rVB3 | 4438272  | 6747846  | 7.56%   | 0.643% |
| 57 | 13.109 | 1830 | 1838 | 1845 | rBV2 | 19454618 | 47195813 | 52.87%  | 4.494% |
| 58 | 13.170 | 1845 | 1848 | 1856 | rVB  | 16399512 | 23688558 | 26.54%  | 2.256% |
| 59 | 13.255 | 1856 | 1862 | 1863 | rBV3 | 35199672 | 53175288 | 59.57%  | 5.063% |
| 60 | 13.353 | 1874 | 1878 | 1881 | rBV2 | 3797028  | 6360012  | 7.13%   | 0.606% |
| 61 | 13.383 | 1881 | 1883 | 1885 | rVV2 | 1707630  | 1570114  | 1.76%   | 0.150% |
| 62 | 13.420 | 1885 | 1889 | 1890 | rBV  | 4485947  | 5547821  | 6.22%   | 0.528% |
| 63 | 13.450 | 1890 | 1894 | 1897 | rVB3 | 10626819 | 15243332 | 17.08%  | 1.451% |
| 64 | 13.493 | 1897 | 1901 | 1904 | rBV2 | 8397766  | 10064166 | 11.28%  | 0.958% |
| 65 | 13.524 | 1904 | 1906 | 1909 | rVB  | 5137551  | 4726663  | 5.30%   | 0.450% |
| 66 | 13.560 | 1909 | 1912 | 1913 | rBV  | 6385406  | 7115276  | 7.97%   | 0.678% |
| 67 | 13.591 | 1913 | 1917 | 1921 | rBV  | 25723431 | 40787517 | 45.70%  | 3.884% |
| 68 | 13.719 | 1930 | 1938 | 1940 | rBV5 | 3963257  | 8226671  | 9.22%   | 0.783% |
| 69 | 13.755 | 1940 | 1944 | 1947 | rVV2 | 19314833 | 32600575 | 36.52%  | 3.104% |
| 70 | 13.792 | 1947 | 1950 | 1959 | rVB2 | 14442908 | 26939305 | 30.18%  | 2.565% |
| 71 | 13.883 | 1959 | 1965 | 1969 | rBV2 | 11860494 | 20694910 | 23.19%  | 1.971% |
| 72 | 13.944 | 1969 | 1975 | 1979 | rVV3 | 6870980  | 15880867 | 17.79%  | 1.512% |
| 73 | 14.017 | 1983 | 1987 | 1988 | rVV  | 6242345  | 9394317  | 10.52%  | 0.895% |
| 74 | 14.036 | 1988 | 1990 | 1995 | rVV  | 7770543  | 12328042 | 13.81%  | 1.174% |
| 75 | 14.097 | 1995 | 2000 | 2004 | rVB  | 9428398  | 14138818 | 15.84%  | 1.346% |
| 76 | 14.200 | 2012 | 2017 | 2022 | rVB4 | 6404090  | 10343253 | 11.59%  | 0.985% |

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 Sample : N1084-10  
 Misc : 6.57g/10mL/MSVOA\_W/SOIL  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_W  
 ClientSampleId :  
 BGLN3

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\_W\Method\SFAMWLM011122SMA.M  
 Title : SFAM01.0

|     |        |      |      |      |      |         |          |        |        |
|-----|--------|------|------|------|------|---------|----------|--------|--------|
| 77  | 14.261 | 2022 | 2027 | 2034 | rVV7 | 2846632 | 6417843  | 7.19%  | 0.611% |
| 78  | 14.341 | 2034 | 2040 | 2044 | rVV4 | 4321261 | 8498297  | 9.52%  | 0.809% |
| 79  | 14.389 | 2044 | 2048 | 2050 | rVV2 | 6695166 | 10297503 | 11.54% | 0.981% |
| 80  | 14.420 | 2050 | 2053 | 2056 | rVV2 | 6278698 | 9775251  | 10.95% | 0.931% |
| 81  | 14.456 | 2056 | 2059 | 2062 | rVV  | 5013497 | 6729341  | 7.54%  | 0.641% |
| 82  | 14.493 | 2062 | 2065 | 2069 | rVV2 | 4620093 | 7135439  | 7.99%  | 0.679% |
| 83  | 14.542 | 2070 | 2073 | 2080 | rVB3 | 2122638 | 4041157  | 4.53%  | 0.385% |
| 84  | 14.639 | 2086 | 2089 | 2093 | rVB  | 1245939 | 1740219  | 1.95%  | 0.166% |
| 85  | 14.688 | 2093 | 2097 | 2100 | rBV3 | 717592  | 1373180  | 1.54%  | 0.131% |
| 86  | 14.725 | 2100 | 2103 | 2108 | rVB  | 1248804 | 2000085  | 2.24%  | 0.190% |
| 87  | 14.792 | 2108 | 2114 | 2119 | rBV2 | 3614924 | 8262627  | 9.26%  | 0.787% |
| 88  | 14.956 | 2137 | 2141 | 2146 | rVB  | 1195570 | 1911191  | 2.14%  | 0.182% |
| 89  | 15.005 | 2146 | 2149 | 2153 | rVB5 | 315448  | 446133   | 0.50%  | 0.042% |
| 90  | 15.060 | 2154 | 2158 | 2166 | rBV3 | 713959  | 1433974  | 1.61%  | 0.137% |
| 91  | 15.164 | 2172 | 2175 | 2184 | rVB7 | 623612  | 1267139  | 1.42%  | 0.121% |
| 92  | 15.255 | 2185 | 2190 | 2195 | rVB5 | 461656  | 1134323  | 1.27%  | 0.108% |
| 93  | 15.316 | 2195 | 2200 | 2203 | rBV2 | 968422  | 1559320  | 1.75%  | 0.148% |
| 94  | 15.359 | 2203 | 2207 | 2213 | rVB  | 1455558 | 2380701  | 2.67%  | 0.227% |
| 95  | 15.651 | 2247 | 2255 | 2261 | rBV9 | 320450  | 950472   | 1.06%  | 0.091% |
| 96  | 15.718 | 2263 | 2266 | 2273 | rVB6 | 248084  | 476765   | 0.53%  | 0.045% |
| 97  | 15.859 | 2286 | 2289 | 2293 | rVB5 | 161503  | 232409   | 0.26%  | 0.022% |
| 98  | 16.273 | 2353 | 2357 | 2365 | rVB  | 299185  | 550503   | 0.62%  | 0.052% |
| 99  | 16.432 | 2377 | 2383 | 2396 | rVB  | 1444947 | 3208206  | 3.59%  | 0.305% |
| 100 | 16.633 | 2409 | 2416 | 2429 | rVB  | 680019  | 1821163  | 2.04%  | 0.173% |

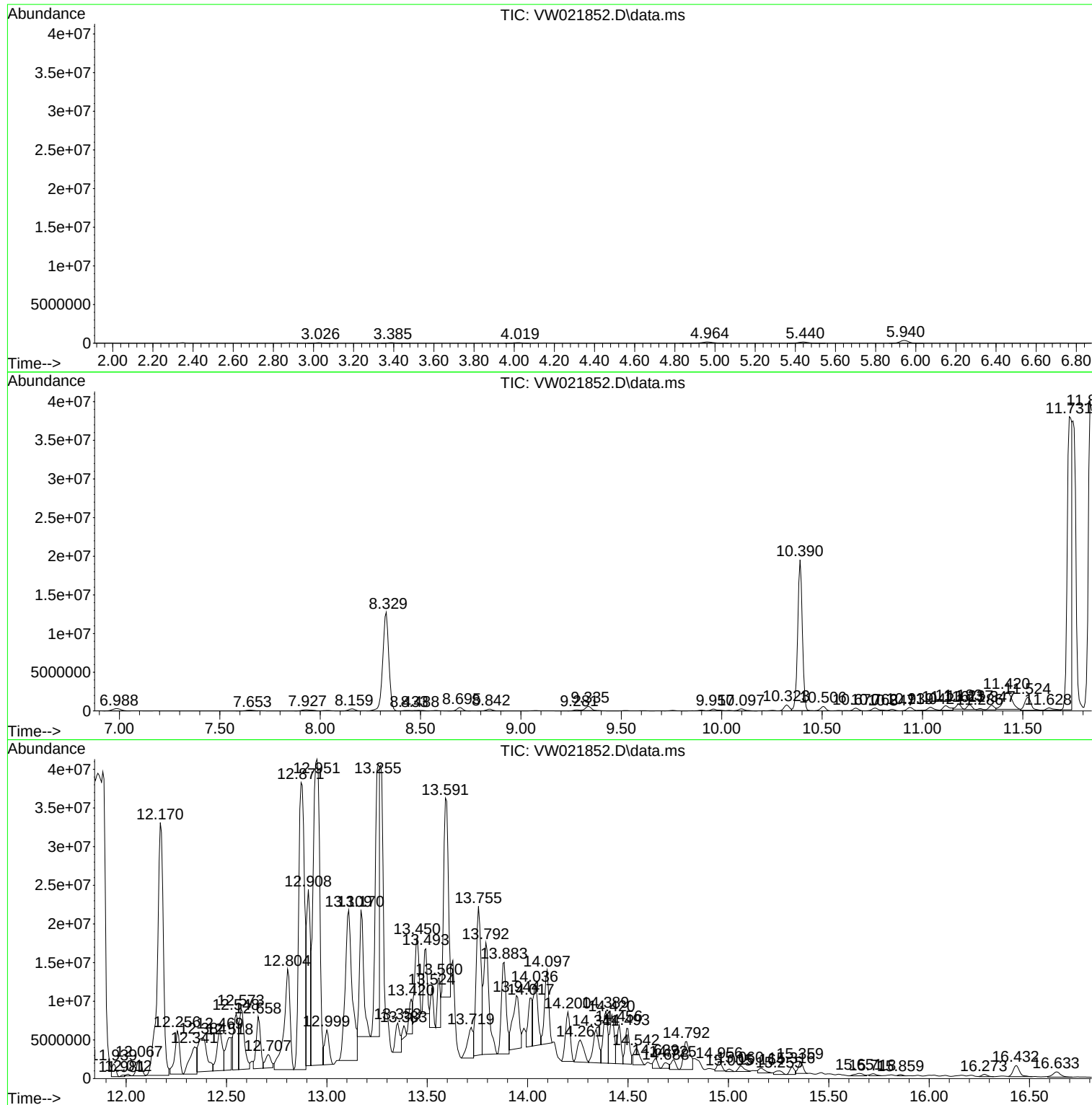
Sum of corrected areas: 1050217761

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
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Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
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Sample : N1084-10  
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ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
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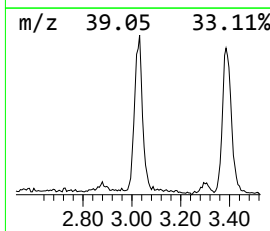
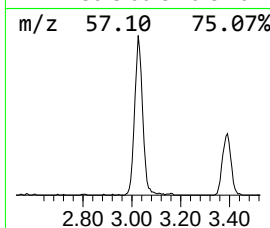
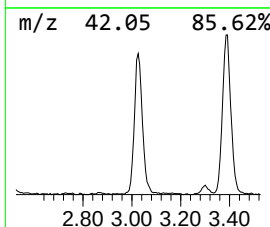
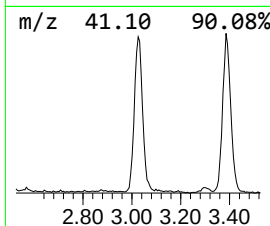
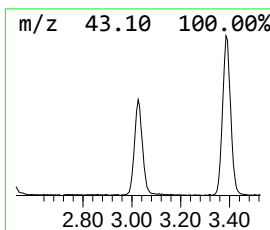
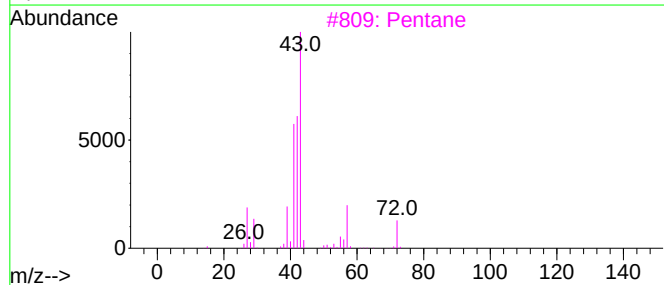
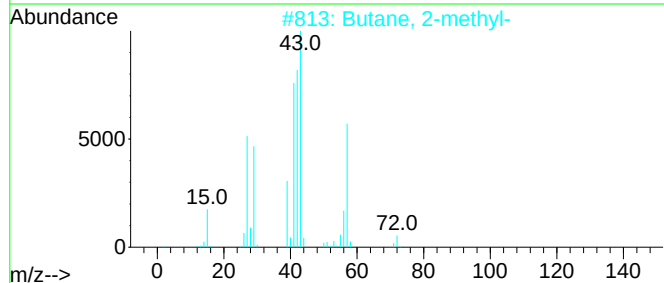
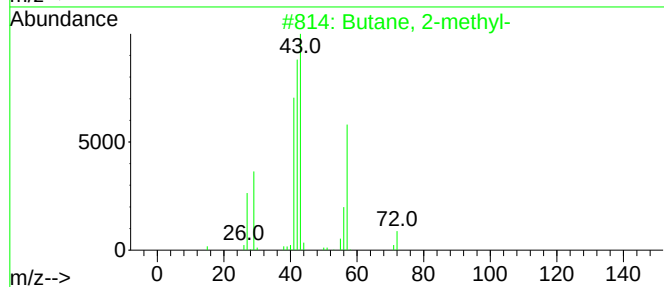
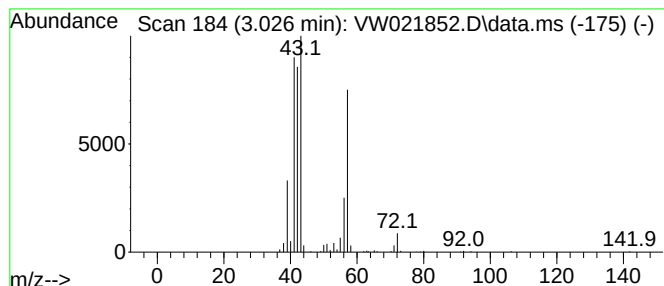
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

\*\*\*\*\*  
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 57

| R.T.  | EstConc   | Area   | Relative to ISTD    | R.T.  |
|-------|-----------|--------|---------------------|-------|
| 3.026 | 8.25 ug/L | 159459 | 1,4-Difluorobenzene | 8.842 |

| Hit# | of                         | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----------------------------|---|--------------|-----|---------|-------------|------|
| 1    | Butane, 2-methyl-          |   |              | 72  | C5H12   | 000078-78-4 | 86   |
| 2    | Butane, 2-methyl-          |   |              | 72  | C5H12   | 000078-78-4 | 86   |
| 3    | Pentane                    |   |              | 72  | C5H12   | 000109-66-0 | 33   |
| 4    | 3-Buten-1-ol               |   |              | 72  | C4H8O   | 000627-27-0 | 25   |
| 5    | Butane, 1-chloro-2-methyl- |   |              | 106 | C5H11Cl | 000616-13-7 | 22   |



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Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

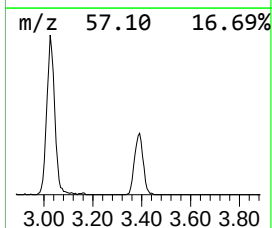
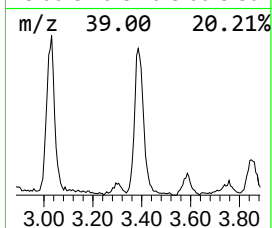
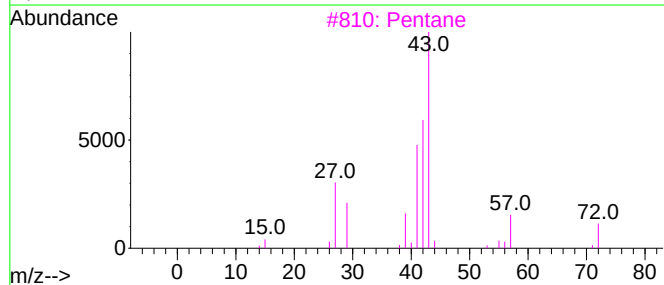
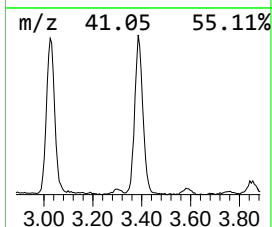
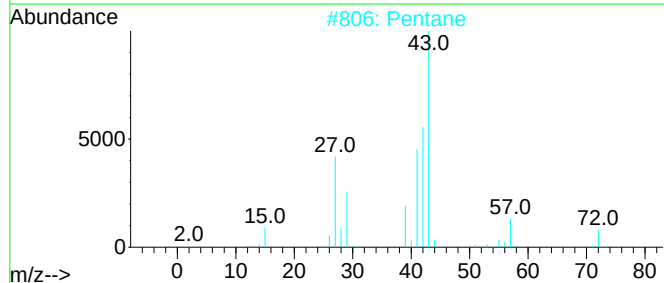
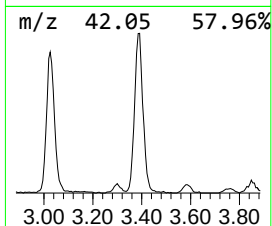
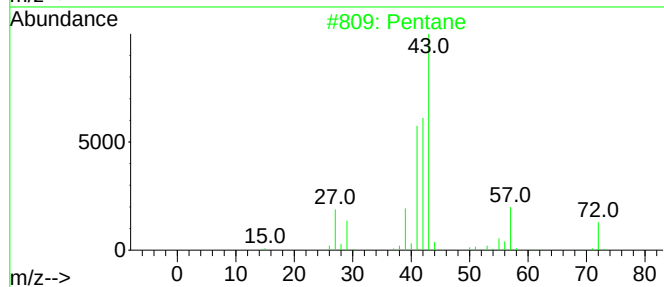
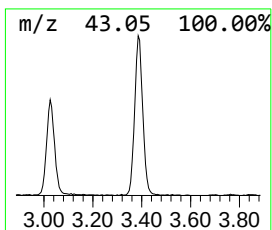
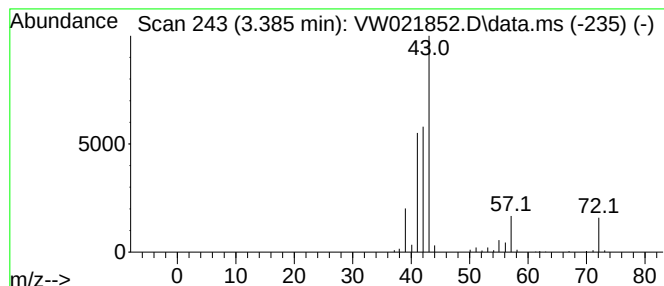
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 55

| R.T.  | EstConc   | Area   | Relative to ISTD    | R.T.  |
|-------|-----------|--------|---------------------|-------|
| 3.385 | 9.31 ug/L | 179967 | 1,4-Difluorobenzene | 8.842 |

| Hit# | of      | 5 | Tentative ID | MW | MolForm | CAS#        | Qual |
|------|---------|---|--------------|----|---------|-------------|------|
| 1    | Pentane |   |              | 72 | C5H12   | 000109-66-0 | 91   |
| 2    | Pentane |   |              | 72 | C5H12   | 000109-66-0 | 86   |
| 3    | Pentane |   |              | 72 | C5H12   | 000109-66-0 | 86   |
| 4    | Pentane |   |              | 72 | C5H12   | 000109-66-0 | 86   |
| 5    | Pentane |   |              | 72 | C5H12   | 000109-66-0 | 86   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

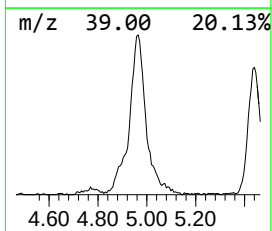
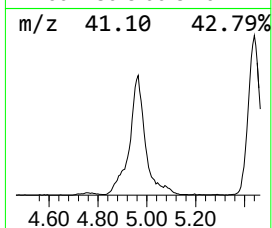
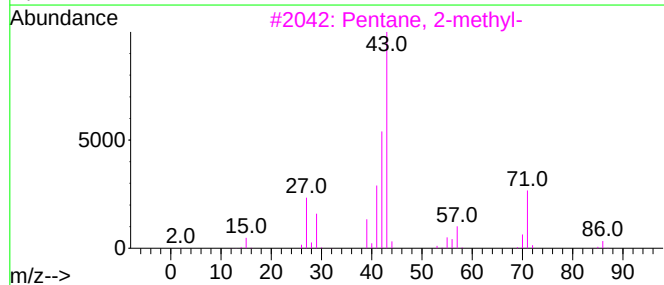
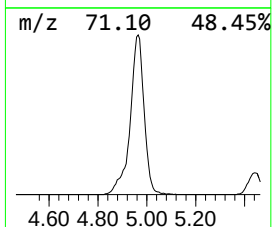
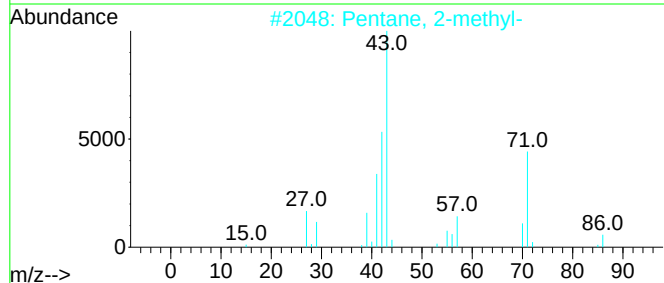
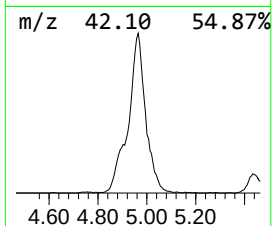
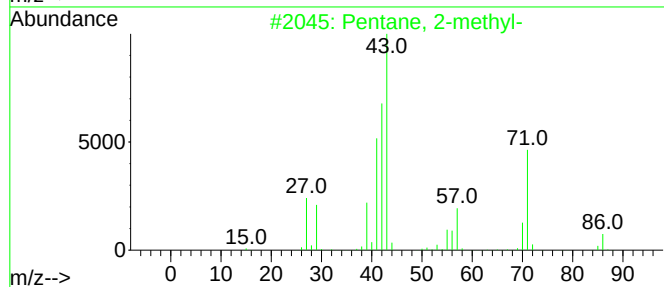
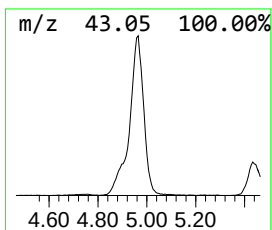
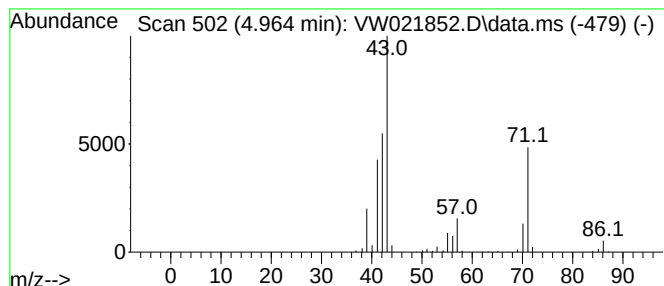
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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 4.964 | 37.02 ug/L | 715506 | 1,4-Difluorobenzene | 8.842 |

| Hit# | of 5 | Tentative ID       | MW | MolForm | CAS#        | Qual |
|------|------|--------------------|----|---------|-------------|------|
| 1    |      | Pentane, 2-methyl- | 86 | C6H14   | 000107-83-5 | 94   |
| 2    |      | Pentane, 2-methyl- | 86 | C6H14   | 000107-83-5 | 91   |
| 3    |      | Pentane, 2-methyl- | 86 | C6H14   | 000107-83-5 | 91   |
| 4    |      | Pentane, 2-methyl- | 86 | C6H14   | 000107-83-5 | 91   |
| 5    |      | Pentane, 2-methyl- | 86 | C6H14   | 000107-83-5 | 83   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

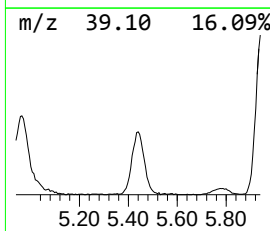
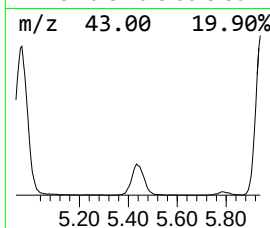
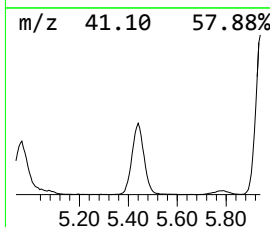
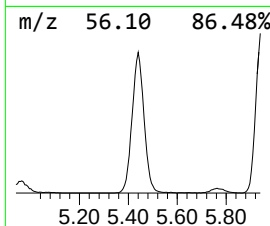
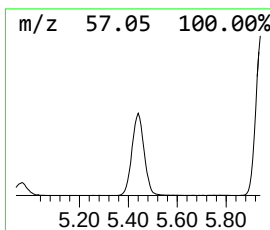
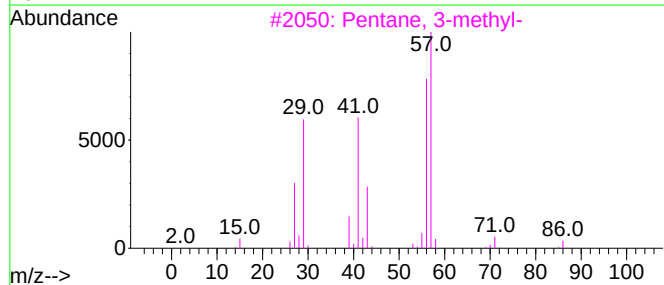
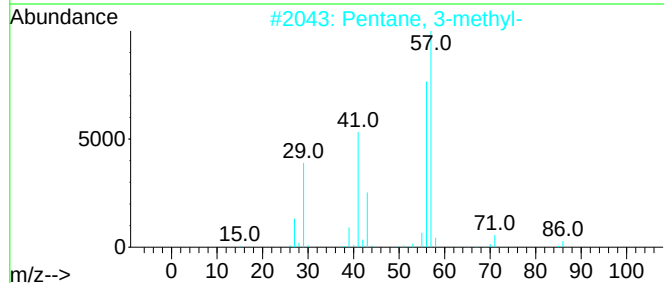
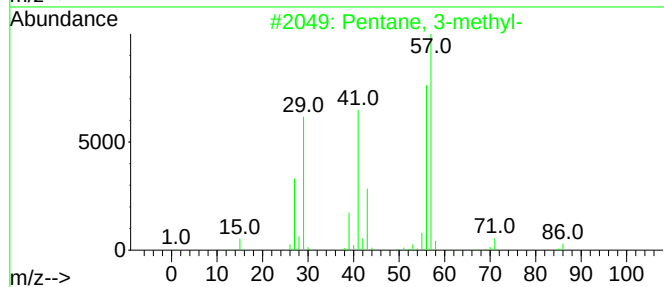
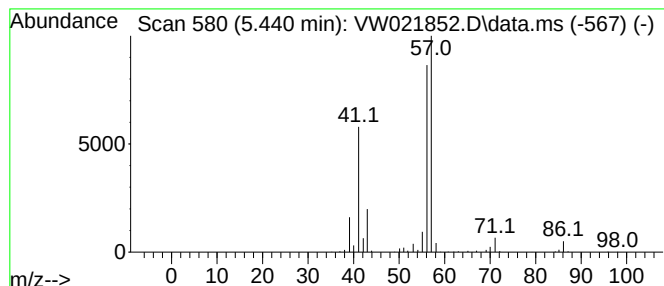
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Quant Title : SFAM01.0

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

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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 40

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 5.440 | 28.19 ug/L | 544829 | 1,4-Difluorobenzene | 8.842 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentane, 3-methyl-       | 86  | C6H14   | 000096-14-0 | 91   |
| 2    |    |   | Pentane, 3-methyl-       | 86  | C6H14   | 000096-14-0 | 90   |
| 3    |    |   | Pentane, 3-methyl-       | 86  | C6H14   | 000096-14-0 | 90   |
| 4    |    |   | Butane, 2,2,3-trimethyl- | 100 | C7H16   | 000464-06-2 | 43   |
| 5    |    |   | Hexane, 2,2,3-trimethyl- | 128 | C9H20   | 016747-25-4 | 43   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

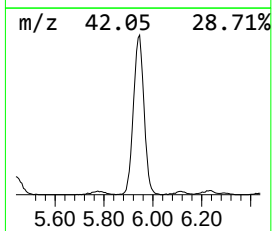
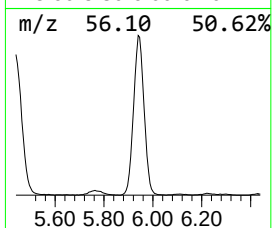
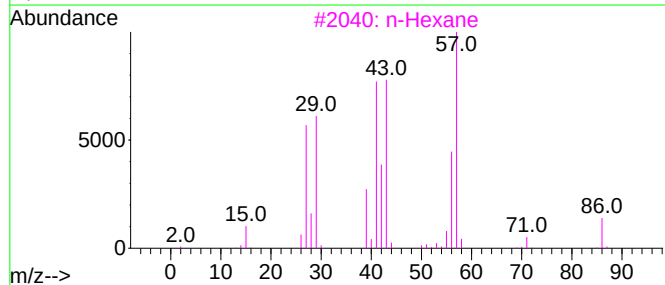
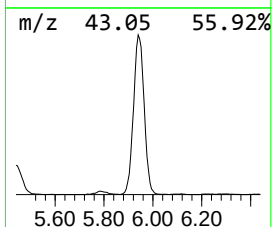
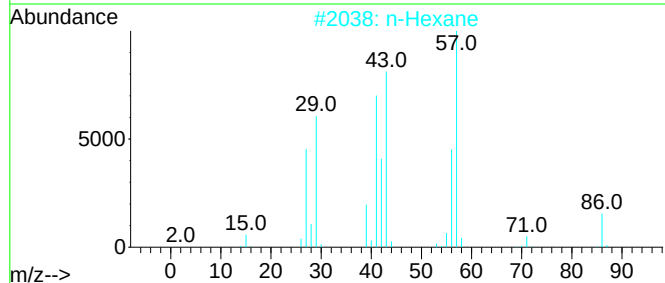
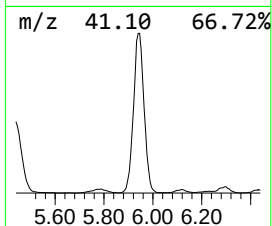
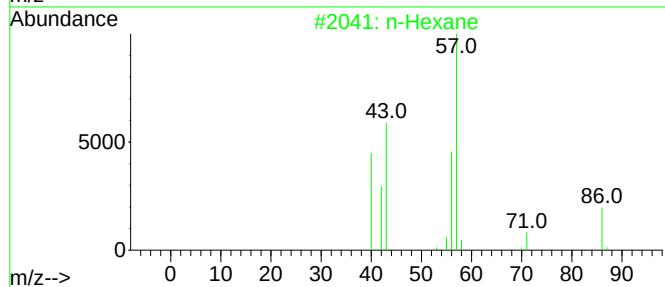
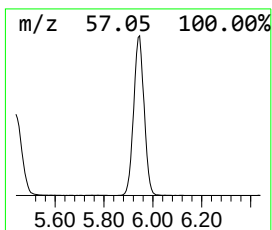
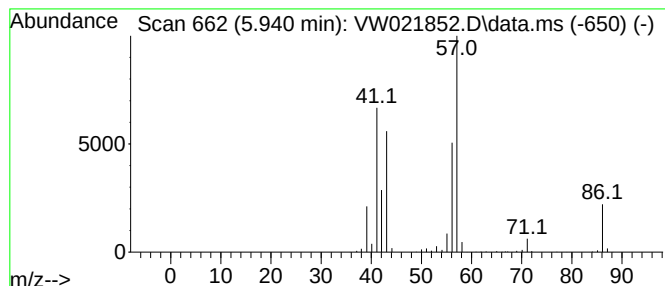
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T.  | EstConc    | Area    | Relative to ISTD    | R.T.  |
|-------|------------|---------|---------------------|-------|
| 5.940 | 55.49 ug/L | 1072410 | 1,4-Difluorobenzene | 8.842 |

| Hit# | of              | 5 | Tentative ID | MW | MolForm | CAS#         | Qual |
|------|-----------------|---|--------------|----|---------|--------------|------|
| 1    | n-Hexane        |   |              | 86 | C6H14   | 000110-54-3  | 87   |
| 2    | n-Hexane        |   |              | 86 | C6H14   | 000110-54-3  | 78   |
| 3    | n-Hexane        |   |              | 86 | C6H14   | 000110-54-3  | 78   |
| 4    | n-Hexane        |   |              | 86 | C6H14   | 000110-54-3  | 78   |
| 5    | 2-Ethyl-oxetane |   |              | 86 | C5H10O  | 1010386-40-2 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

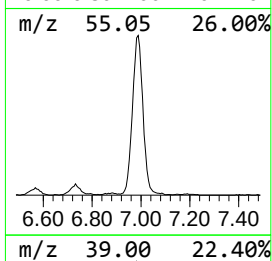
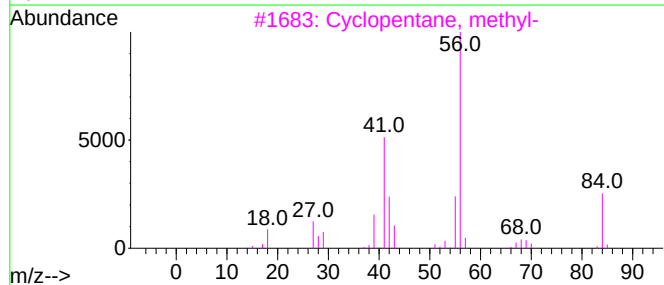
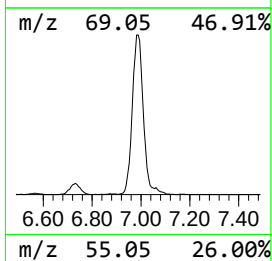
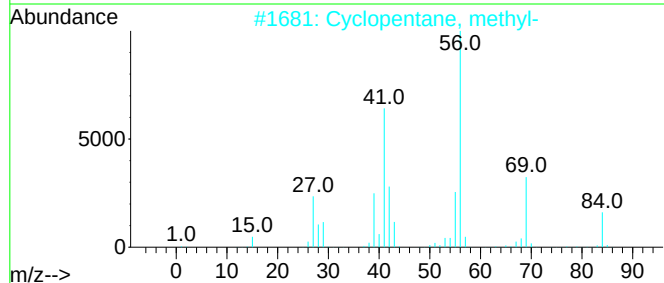
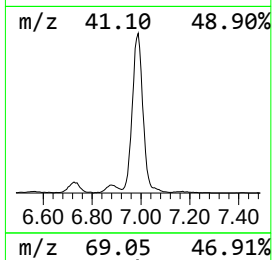
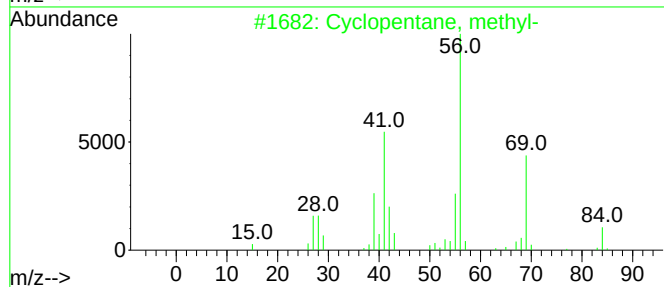
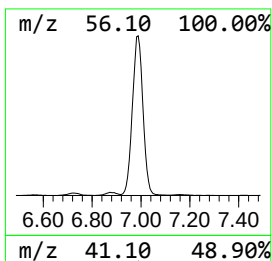
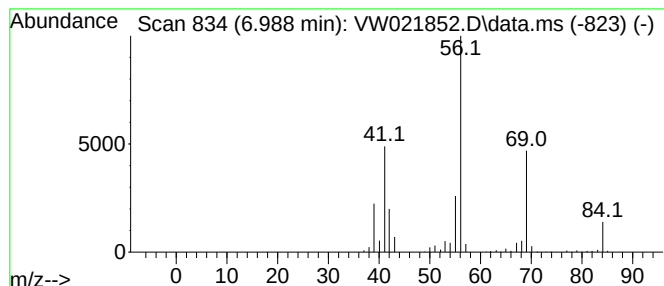
Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

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Peak Number 6 (DEL) Alkane: Cyclic6.988 Concentration Rank 21

| R.T.      | EstConc               | Area   | Relative to ISTD    | R.T.        |      |
|-----------|-----------------------|--------|---------------------|-------------|------|
| 6.988     | 44.67 ug/L            | 863296 | 1,4-Difluorobenzene | 8.842       |      |
| Hit# of 5 | Tentative ID          | MW     | MolForm             | CAS#        | Qual |
| 1         | Cyclopentane, methyl- | 84     | C6H12               | 000096-37-7 | 94   |
| 2         | Cyclopentane, methyl- | 84     | C6H12               | 000096-37-7 | 91   |
| 3         | Cyclopentane, methyl- | 84     | C6H12               | 000096-37-7 | 78   |
| 4         | Cyclohexane           | 84     | C6H12               | 000110-82-7 | 78   |
| 5         | 1-Pentene, 2-methyl-  | 84     | C6H12               | 000763-29-1 | 72   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L  
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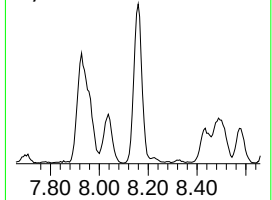
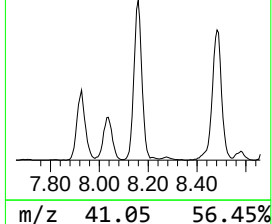
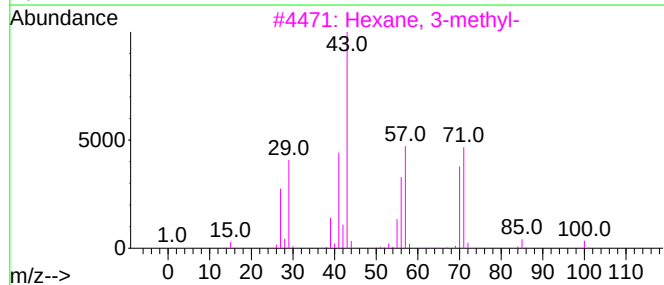
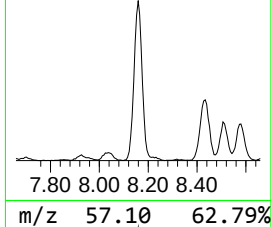
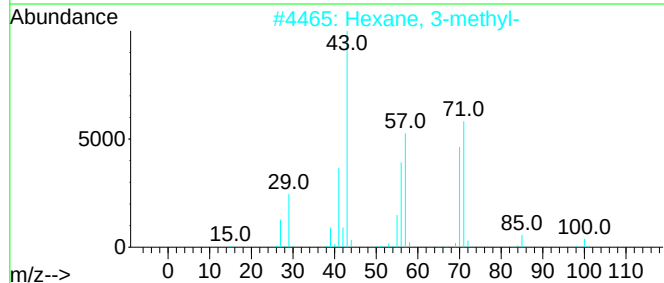
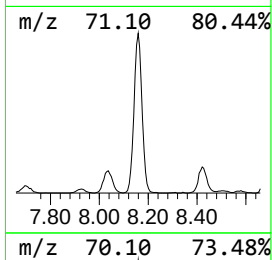
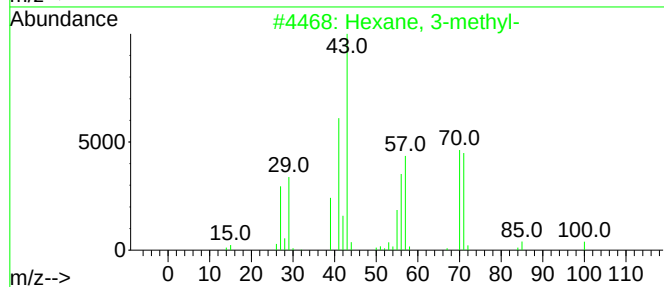
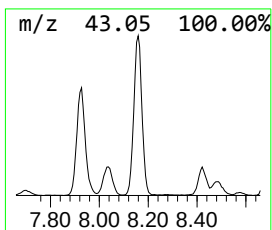
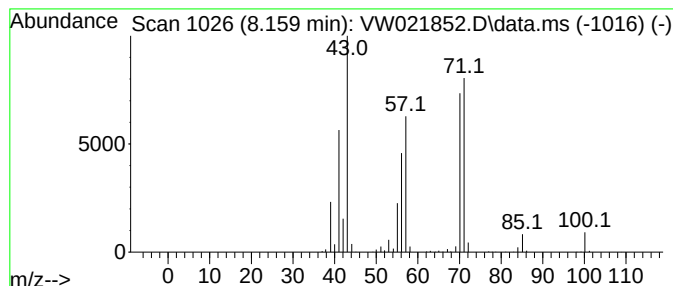
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Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 34

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 8.159 | 31.13 ug/L | 601585 | 1,4-Difluorobenzene | 8.842 |

| Hit# | of 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------------|-----|---------|-------------|------|
| 1    |      | Hexane, 3-methyl-         | 100 | C7H16   | 000589-34-4 | 95   |
| 2    |      | Hexane, 3-methyl-         | 100 | C7H16   | 000589-34-4 | 90   |
| 3    |      | Hexane, 3-methyl-         | 100 | C7H16   | 000589-34-4 | 76   |
| 4    |      | Pentane, 2,3,4-trimethyl- | 114 | C8H18   | 000565-75-3 | 64   |
| 5    |      | Decane, 3,3,4-trimethyl-  | 184 | C13H28  | 049622-18-6 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

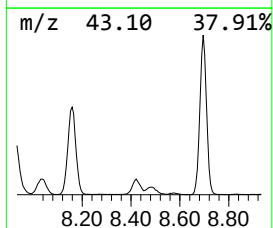
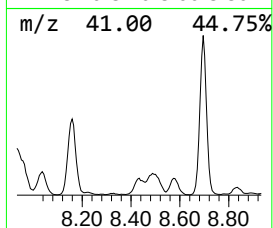
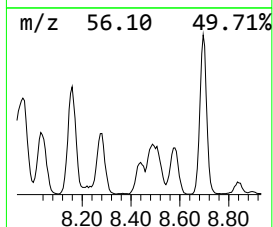
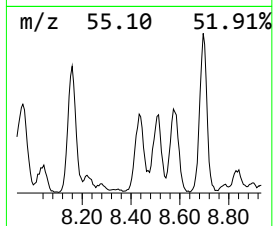
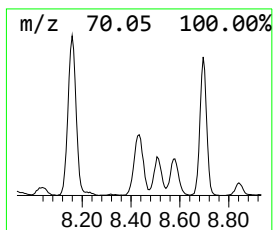
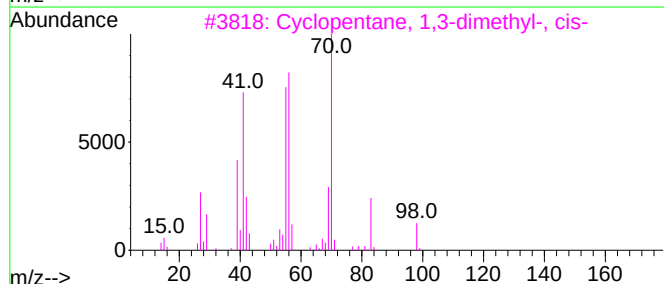
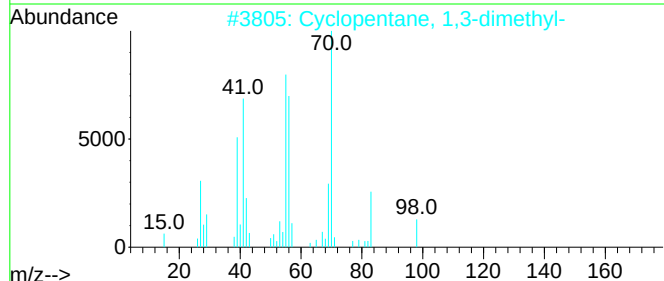
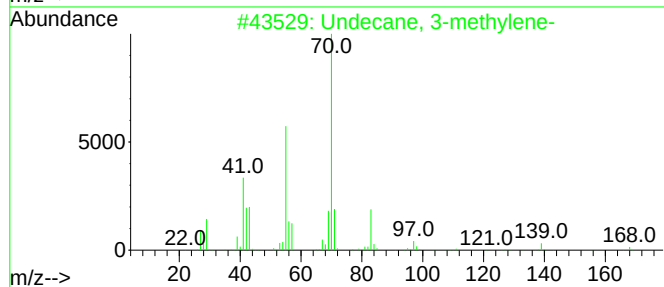
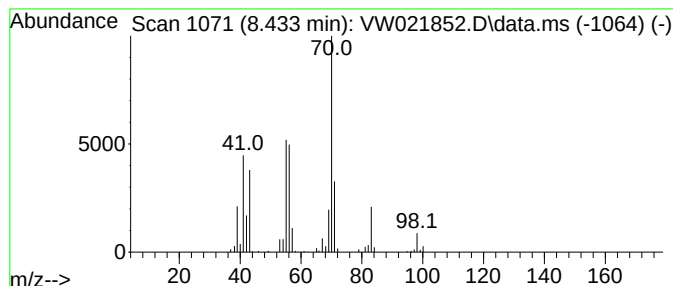
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Quant Title : SFAM01.0

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

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Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 54

| R.T.      | EstConc                           | Area   | Relative to ISTD    | R.T.        |      |
|-----------|-----------------------------------|--------|---------------------|-------------|------|
| 8.433     | 9.53 ug/L                         | 184204 | 1,4-Difluorobenzene | 8.842       |      |
| Hit# of 5 | Tentative ID                      | MW     | MolForm             | CAS#        | Qual |
| 1         | Undecane, 3-methylene-            | 168    | C12H24              | 071138-64-2 | 59   |
| 2         | Cyclopentane, 1,3-dimethyl-       | 98     | C7H14               | 002453-00-1 | 58   |
| 3         | Cyclopentane, 1,3-dimethyl-, cis- | 98     | C7H14               | 002532-58-3 | 53   |
| 4         | Cyclopentane, 1,2-dimethyl-, cis- | 98     | C7H14               | 001192-18-3 | 53   |
| 5         | Cyclopentane, 1,3-dimethyl-, cis- | 98     | C7H14               | 002532-58-3 | 52   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

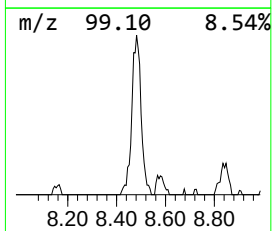
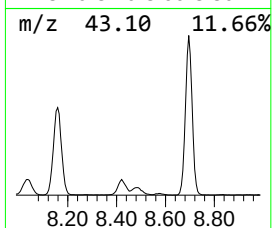
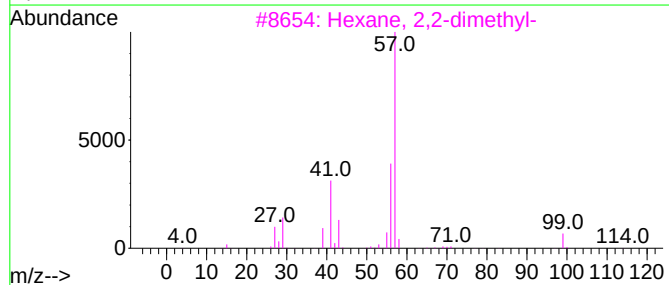
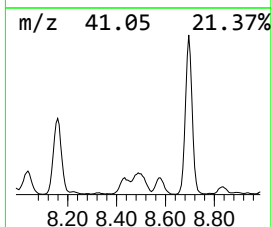
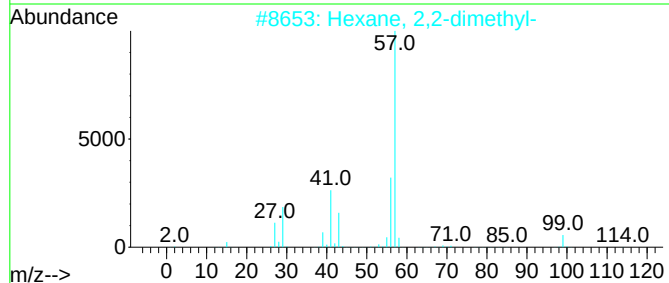
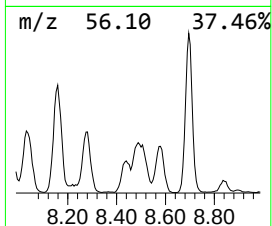
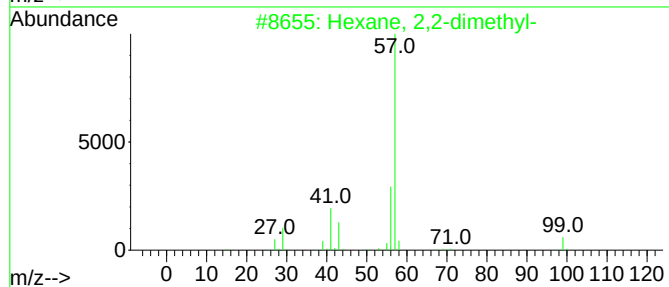
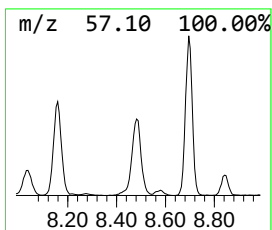
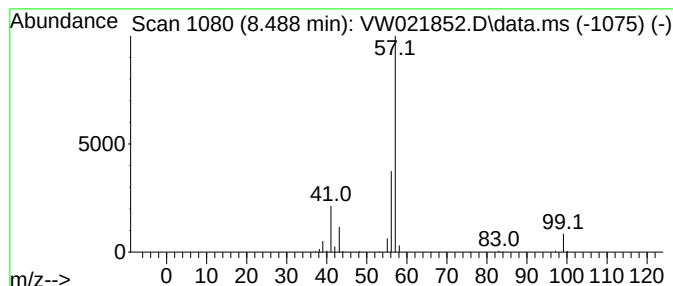
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Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 51

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 8.488 | 12.39 ug/L | 239438 | 1,4-Difluorobenzene | 8.842 |

| Hit# | of 5                      | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|---------------------------|--------------|-----|---------|-------------|------|
| 1    | Hexane, 2,2-dimethyl-     |              | 114 | C8H18   | 000590-73-8 | 72   |
| 2    | Hexane, 2,2-dimethyl-     |              | 114 | C8H18   | 000590-73-8 | 72   |
| 3    | Hexane, 2,2-dimethyl-     |              | 114 | C8H18   | 000590-73-8 | 72   |
| 4    | Pentane, 2,2,4-trimethyl- |              | 114 | C8H18   | 000540-84-1 | 56   |
| 5    | Methane, isocyanato-      |              | 57  | C2H3NO  | 000624-83-9 | 45   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

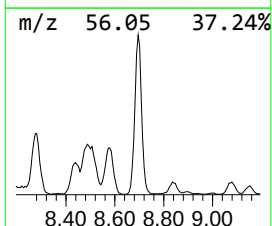
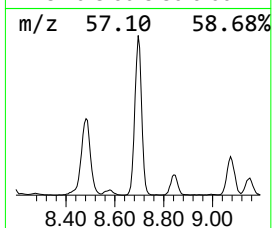
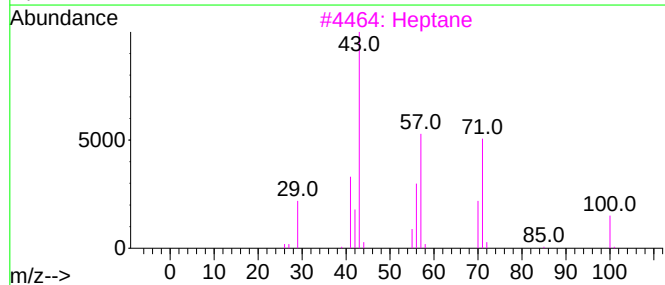
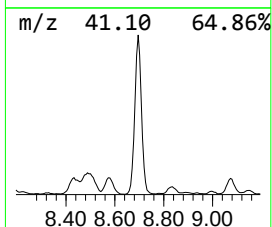
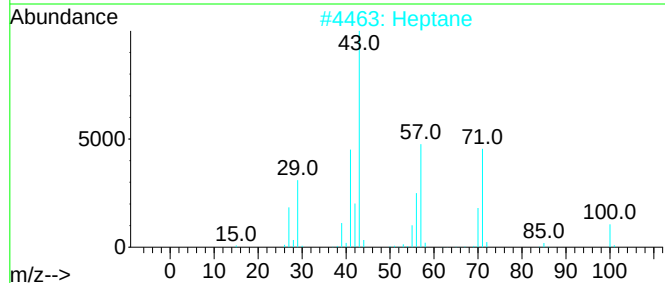
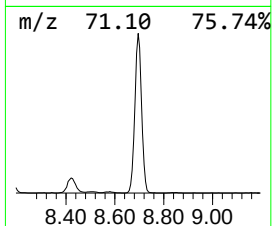
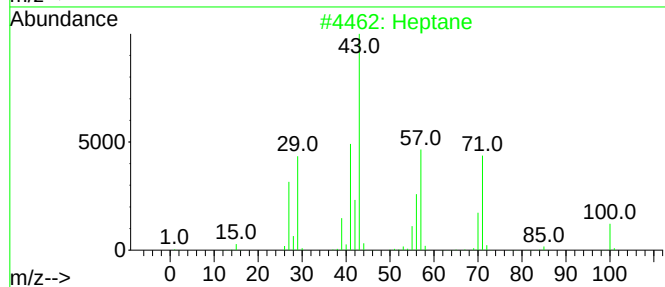
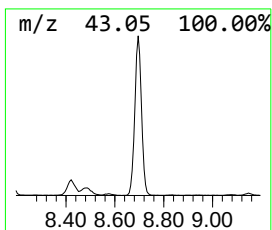
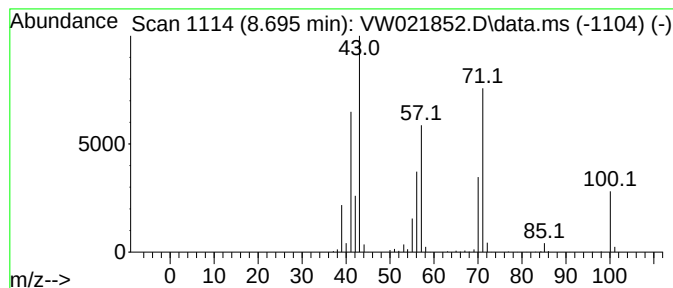
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TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 8.695 | 44.70 ug/L | 863839 | 1,4-Difluorobenzene | 8.842 |

| Hit# | of 5 | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------|-----|---------|-------------|------|
| 1    |      | Heptane            | 100 | C7H16   | 000142-82-5 | 87   |
| 2    |      | Heptane            | 100 | C7H16   | 000142-82-5 | 87   |
| 3    |      | Heptane            | 100 | C7H16   | 000142-82-5 | 72   |
| 4    |      | Heptane            | 100 | C7H16   | 000142-82-5 | 53   |
| 5    |      | Pentane, 2-methyl- | 86  | C6H14   | 000107-83-5 | 53   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

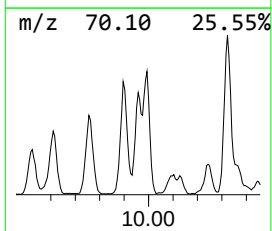
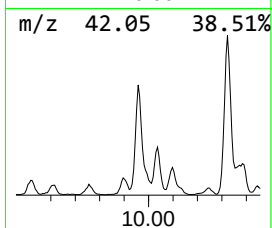
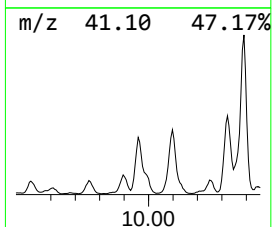
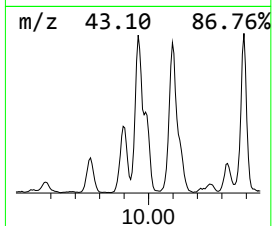
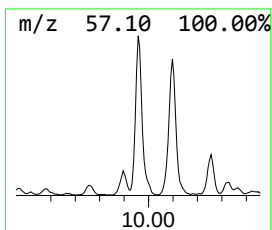
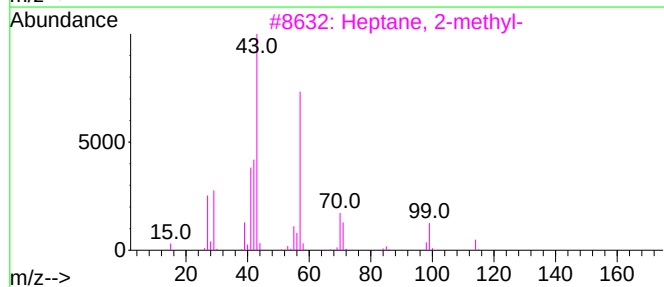
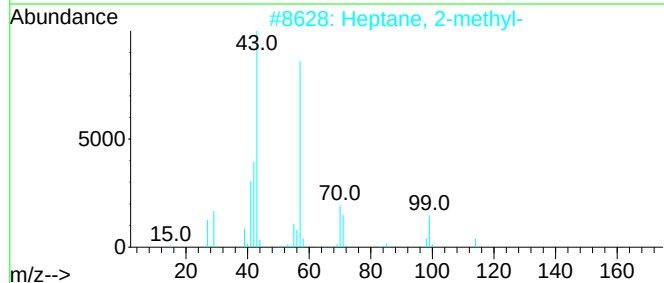
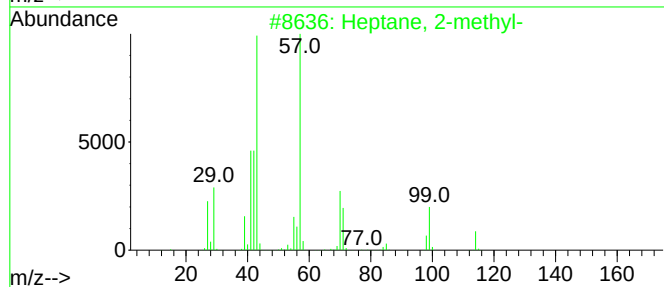
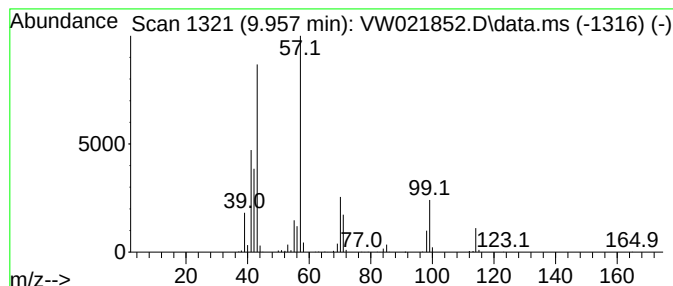
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Quant Title : SFAM01.0

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

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Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 47

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 9.957 | 19.66 ug/L | 379886 | 1,4-Difluorobenzene | 8.842 |

| 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 | 91   |
| 3    |    |   | Heptane, 2-methyl-    | 114 | C8H18   | 000592-27-8 | 91   |
| 4    |    |   | 2-Piperidinone        | 99  | C5H9NO  | 000675-20-7 | 58   |
| 5    |    |   | Hexane, 2,5-dimethyl- | 114 | C8H18   | 000592-13-2 | 47   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

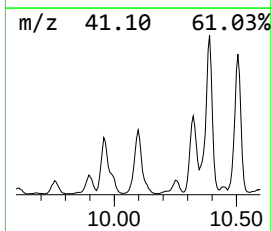
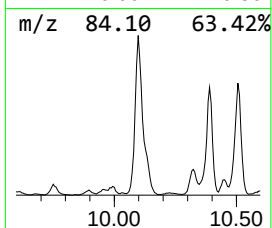
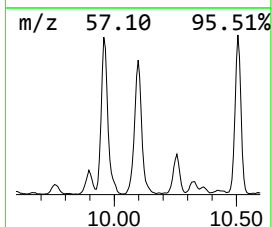
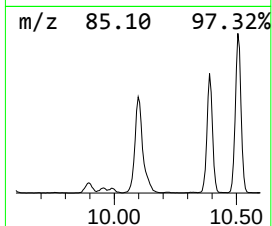
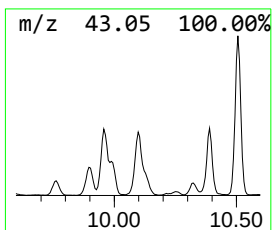
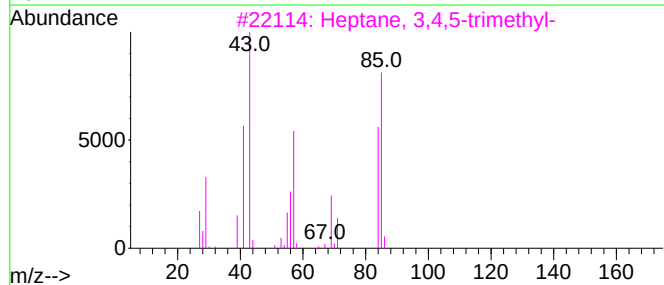
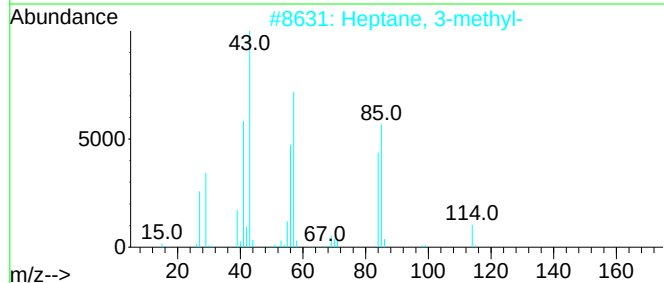
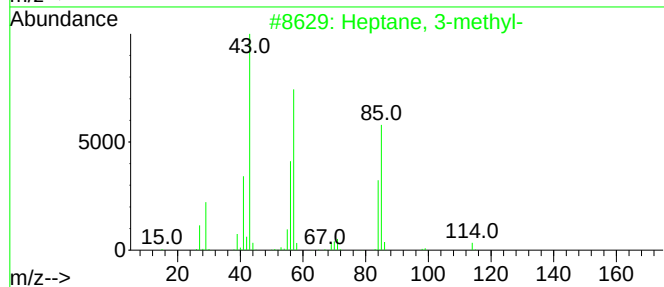
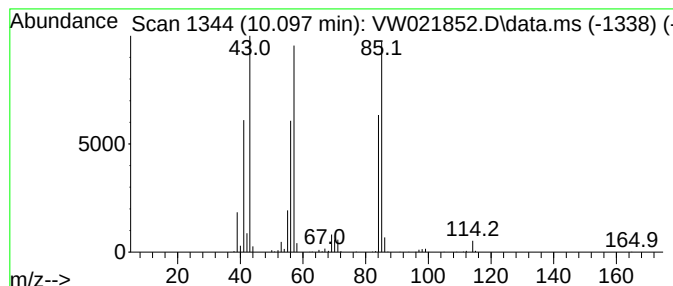
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Quant Title : SFAM01.0

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

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

| R.T.      | EstConc                             | Area   | Relative to ISTD    | R.T.        |      |
|-----------|-------------------------------------|--------|---------------------|-------------|------|
| 10.097    | 30.63 ug/L                          | 591973 | 1,4-Difluorobenzene | 8.842       |      |
| 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 | 83   |
| 3         | Heptane, 3,4,5-trimethyl-           | 142    | C10H22              | 020278-89-1 | 72   |
| 4         | Butanoic acid, 2-methyl-, 2-meth... | 186    | C11H22O2            | 083783-88-4 | 72   |
| 5         | Pentane, 3-ethyl-2,4-dimethyl-      | 128    | C9H20               | 001068-87-7 | 59   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

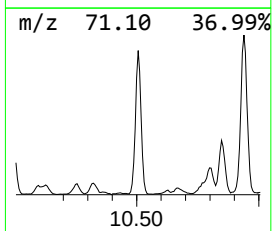
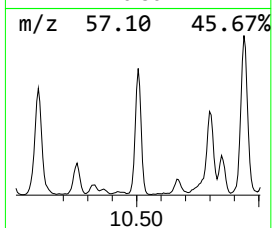
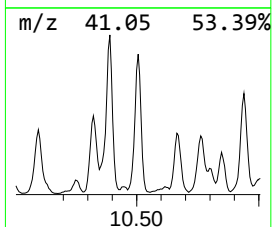
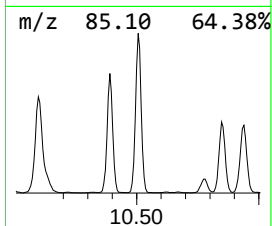
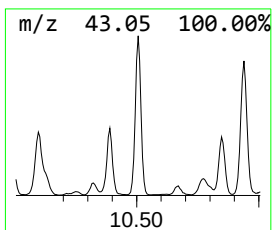
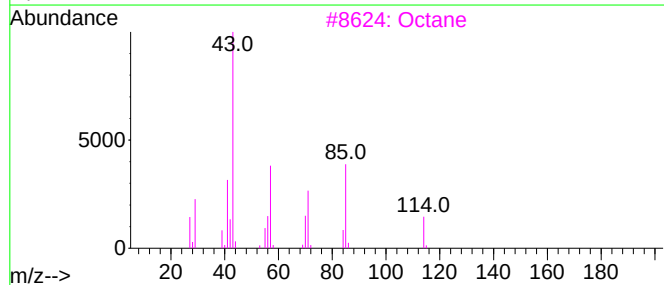
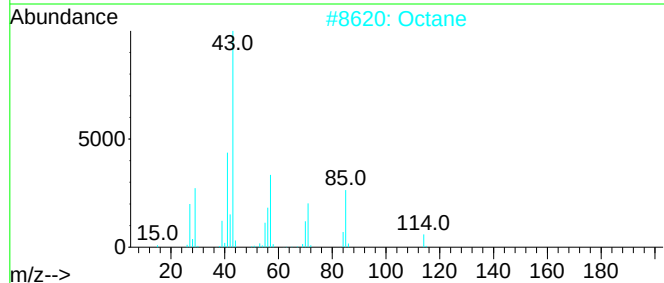
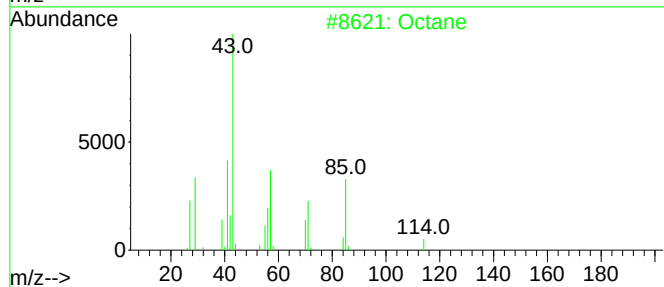
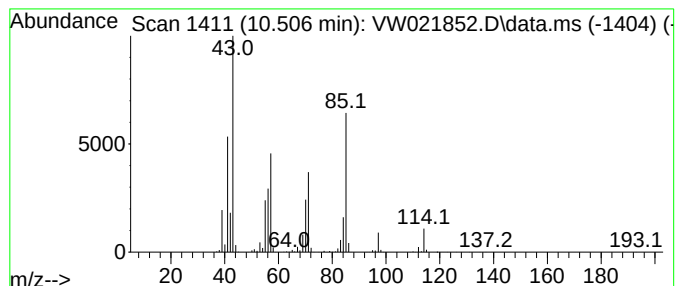
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TIC Integration Parameters: LSCINT.P

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Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 10.506 | 31.29 ug/L | 1034150 | Chlorobenzene-d5 | 11.634 |

| Hit# | of     | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|--------|---|--------------|-----|---------|-------------|------|
| 1    | Octane |   |              | 114 | C8H18   | 000111-65-9 | 90   |
| 2    | Octane |   |              | 114 | C8H18   | 000111-65-9 | 90   |
| 3    | Octane |   |              | 114 | C8H18   | 000111-65-9 | 87   |
| 4    | Octane |   |              | 114 | C8H18   | 000111-65-9 | 87   |
| 5    | Octane |   |              | 114 | C8H18   | 000111-65-9 | 87   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

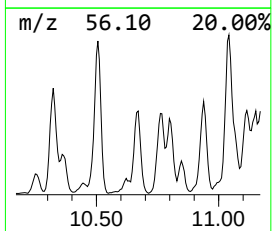
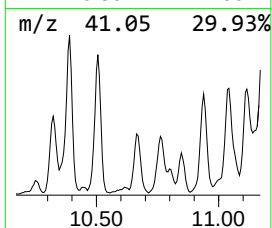
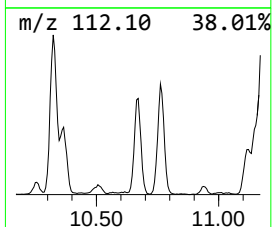
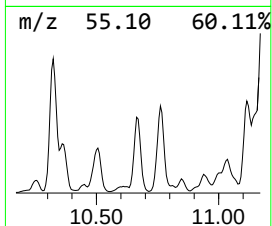
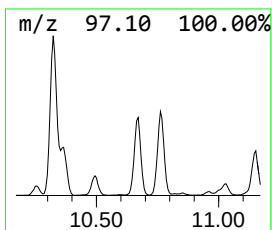
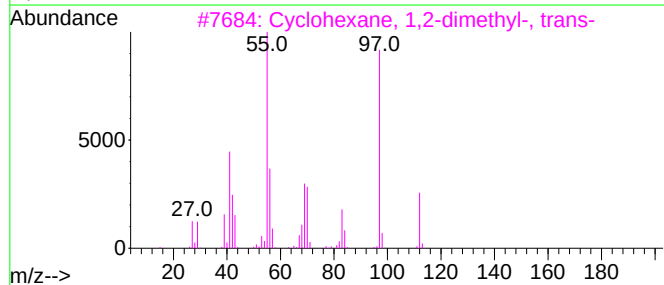
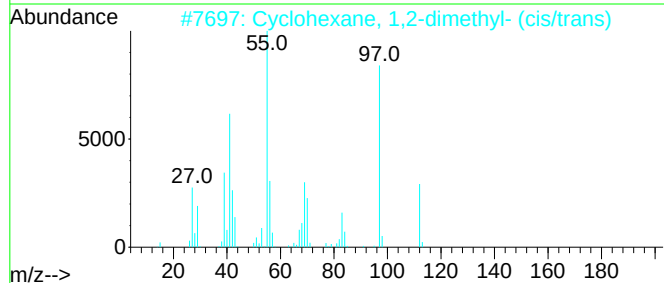
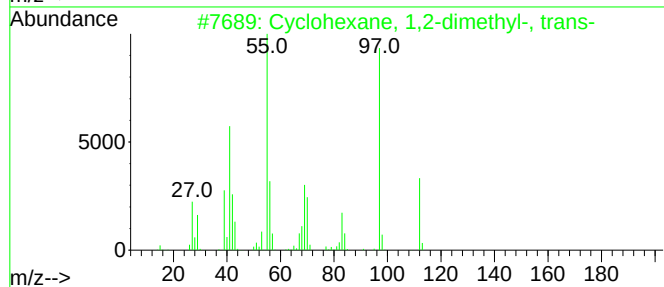
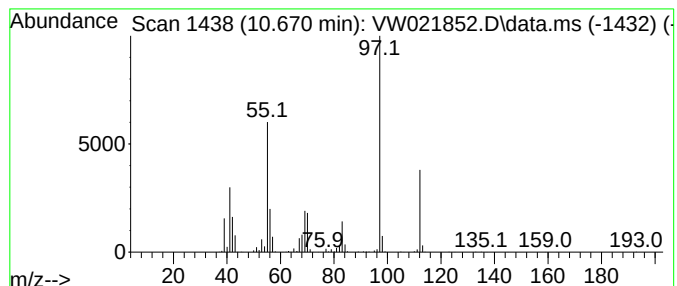
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Quant Title : SFAM01.0

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

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Peak Number 14 (DEL) Alkane: Cyclic10.671 Concentration Rank 49

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.671 | 19.32 ug/L | 638517 | Chlorobenzene-d5 | 11.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,2-dimethyl-, trans-  | 112 | C8H16   | 006876-23-9 | 91   |
| 2    |    |   | Cyclohexane, 1,2-dimethyl- (cis/... | 112 | C8H16   | 000583-57-3 | 91   |
| 3    |    |   | Cyclohexane, 1,2-dimethyl-, trans-  | 112 | C8H16   | 006876-23-9 | 91   |
| 4    |    |   | Cyclohexane, 1,2-dimethyl- (cis/... | 112 | C8H16   | 000583-57-3 | 91   |
| 5    |    |   | Cyclohexane, 1,2-dimethyl-, trans-  | 112 | C8H16   | 006876-23-9 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

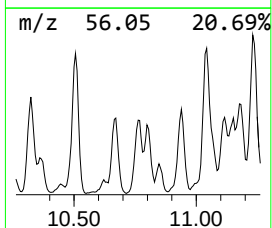
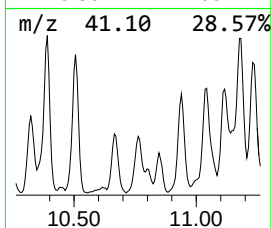
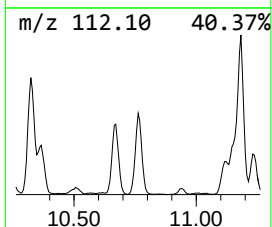
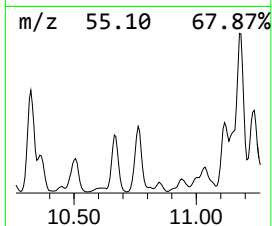
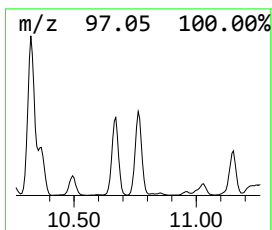
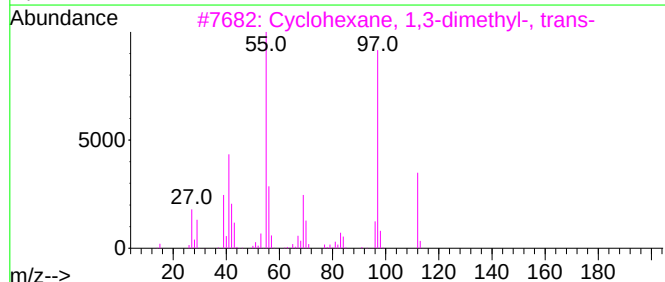
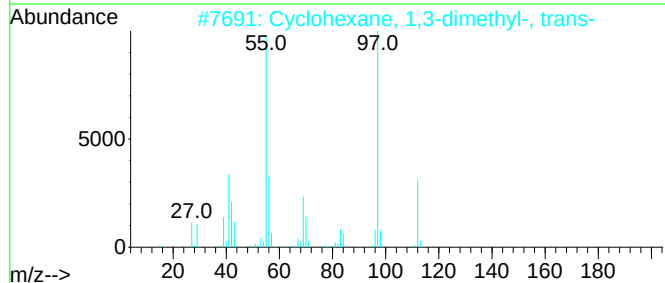
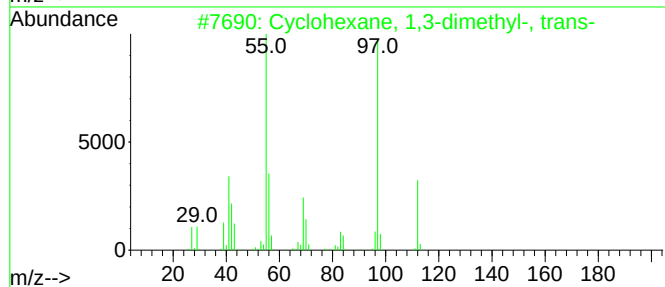
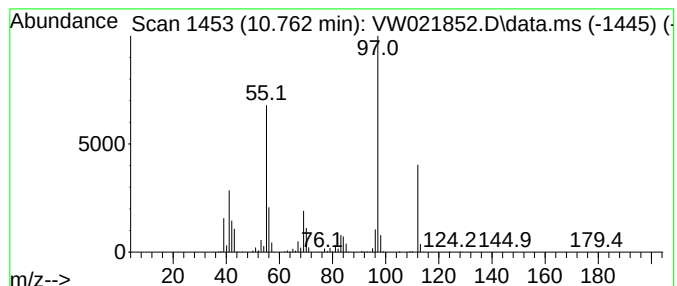
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Quant Title : SFAM01.0

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

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Peak Number 15 (DEL) Alkane: Cyclic10.762 Concentration Rank 44

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.762 | 22.77 ug/L | 752599 | Chlorobenzene-d5 | 11.634 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16   | 002207-03-6 | 94   |
| 2    |    |   | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16   | 002207-03-6 | 94   |
| 3    |    |   | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16   | 002207-03-6 | 94   |
| 4    |    |   | Cyclohexane, 1,3-dimethyl-, trans- | 112 | C8H16   | 002207-03-6 | 94   |
| 5    |    |   | Cyclohexane, 1,4-dimethyl-, trans- | 112 | C8H16   | 002207-04-7 | 93   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

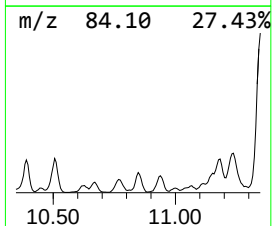
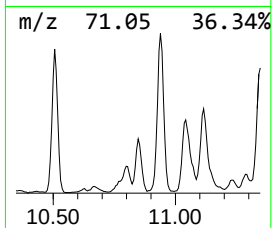
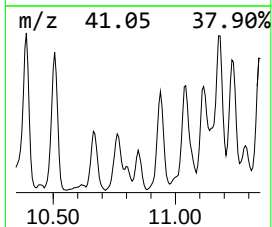
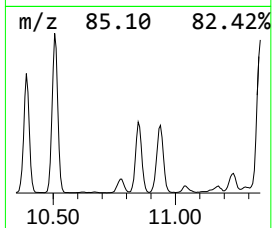
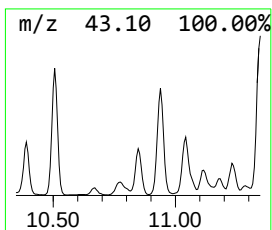
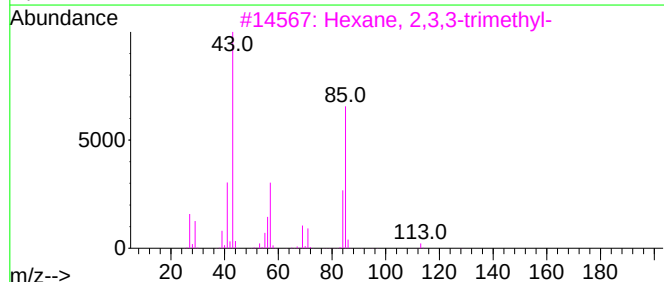
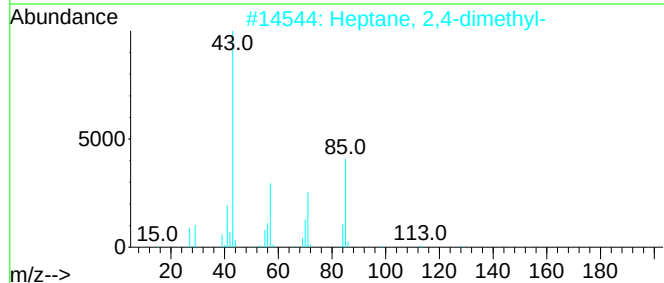
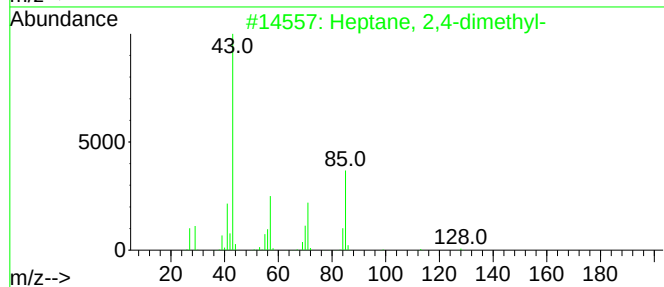
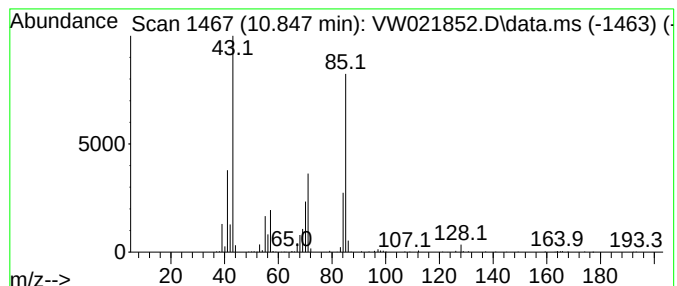
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 53

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.847 | 10.20 ug/L | 337028 | Chlorobenzene-d5 | 11.634 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptane, 2,4-dimethyl-        | 128 | C9H20   | 002213-23-2 | 91   |
| 2    |    |   | Heptane, 2,4-dimethyl-        | 128 | C9H20   | 002213-23-2 | 90   |
| 3    |    |   | Hexane, 2,3,3-trimethyl-      | 128 | C9H20   | 016747-28-7 | 59   |
| 4    |    |   | Pentane, 2,3,3,4-tetramethyl- | 128 | C9H20   | 016747-38-9 | 59   |
| 5    |    |   | Hexane, 2,3,4-trimethyl-      | 128 | C9H20   | 000921-47-1 | 58   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

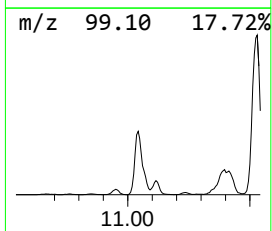
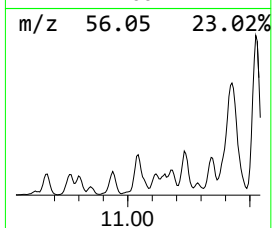
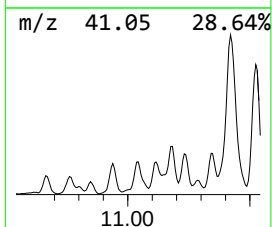
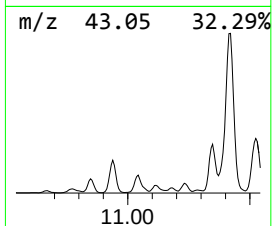
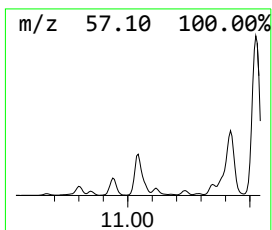
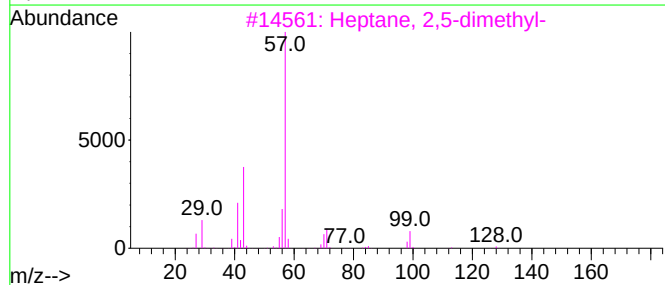
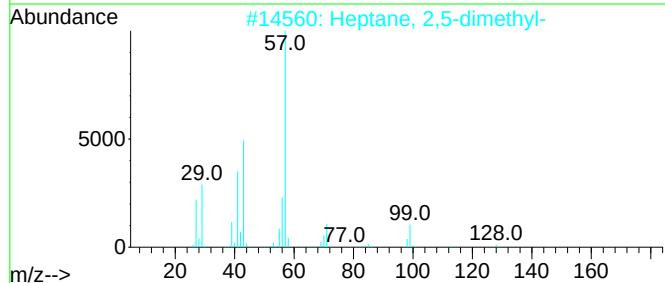
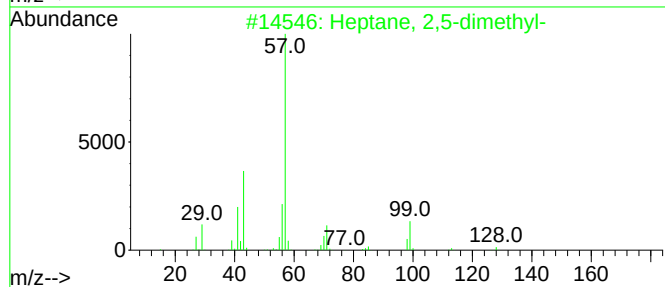
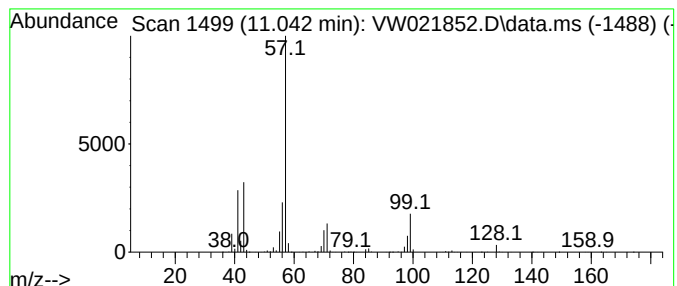
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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.042 | 30.05 ug/L | 993184 | Chlorobenzene-d5 | 11.634 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------|-----|---------|-------------|------|
| 1    |      | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 91   |
| 2    |      | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 87   |
| 3    |      | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 87   |
| 4    |      | Heptane, 2,5-dimethyl- | 128 | C9H20   | 002216-30-0 | 81   |
| 5    |      | Heptane, 3,5-dimethyl- | 128 | C9H20   | 000926-82-9 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

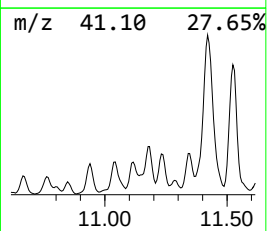
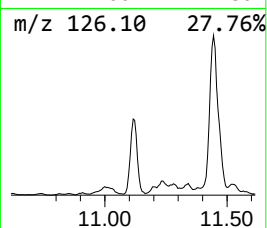
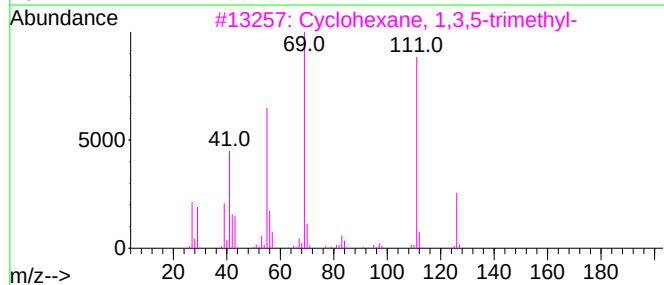
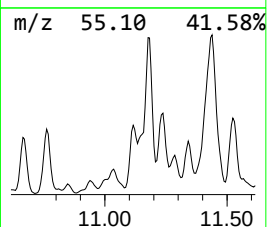
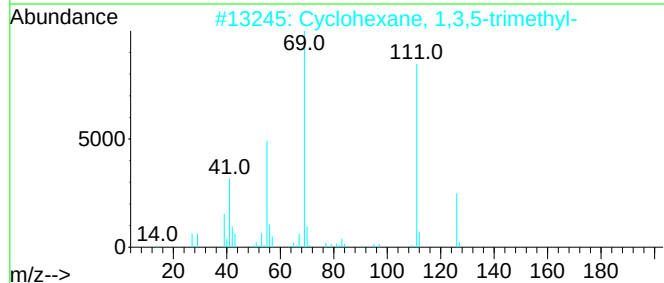
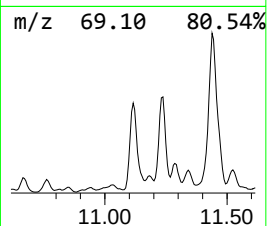
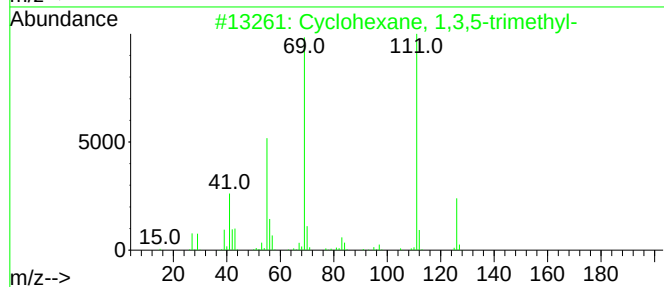
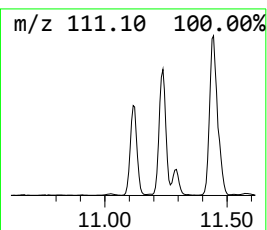
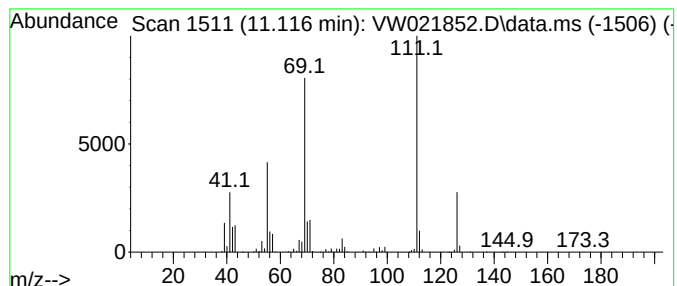
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 (DEL) Alkane: Cyclic11.116 Concentration Rank 23

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 11.116 | 40.98 ug/L | 1354570 | Chlorobenzene-d5 | 11.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,3,5-trimethyl-       | 126 | C9H18   | 001839-63-0 | 96   |
| 2    |    |   | Cyclohexane, 1,3,5-trimethyl-       | 126 | C9H18   | 001839-63-0 | 96   |
| 3    |    |   | Cyclohexane, 1,3,5-trimethyl-       | 126 | C9H18   | 001839-63-0 | 91   |
| 4    |    |   | Cyclohexane, 1,3,5-trimethyl-, (... | 126 | C9H18   | 001795-27-3 | 91   |
| 5    |    |   | Cyclohexane, 1,3,5-trimethyl-, (... | 126 | C9H18   | 001795-27-3 | 81   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

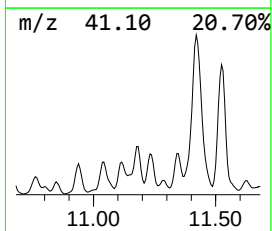
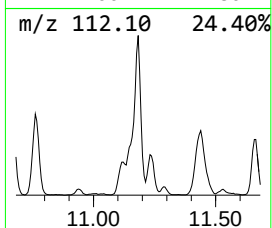
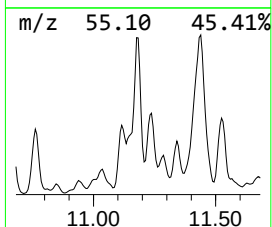
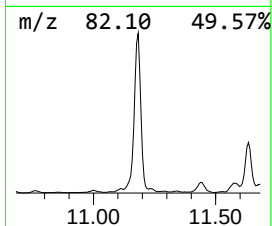
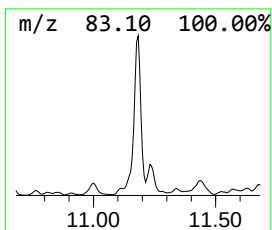
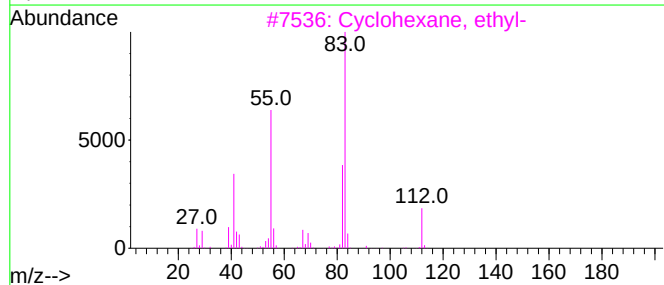
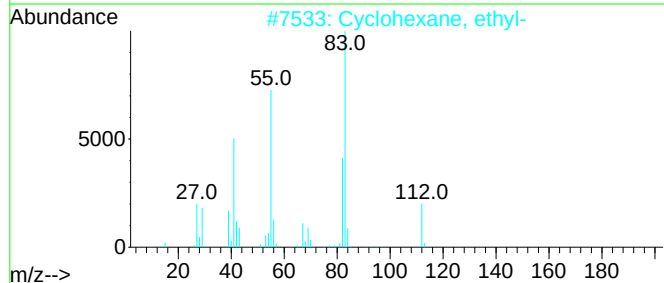
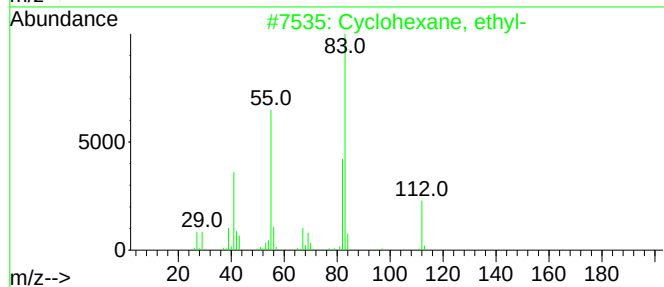
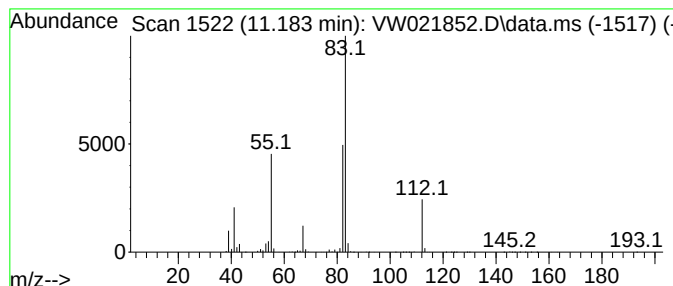
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Quant Title : SFAM01.0

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

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Peak Number 19 (DEL) Alkane: Cyclic11.183 Concentration Rank 30

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 11.183 | 32.71 ug/L | 1081180 | Chlorobenzene-d5 | 11.634 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, ethyl- | 112 | C8H16   | 001678-91-7 | 83   |
| 2    |    |   | Cyclohexane, ethyl- | 112 | C8H16   | 001678-91-7 | 83   |
| 3    |    |   | Cyclohexane, ethyl- | 112 | C8H16   | 001678-91-7 | 72   |
| 4    |    |   | Cyclohexane, ethyl- | 112 | C8H16   | 001678-91-7 | 72   |
| 5    |    |   | Cyclohexane, butyl- | 140 | C10H20  | 001678-93-9 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

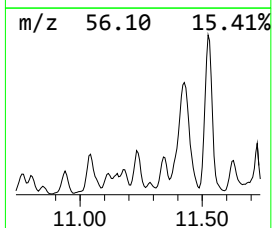
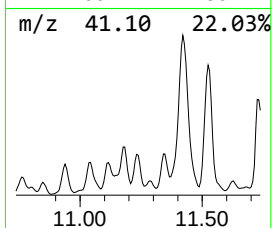
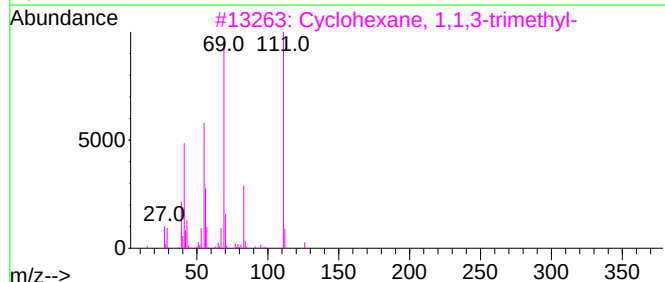
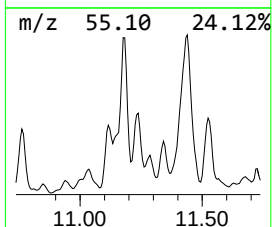
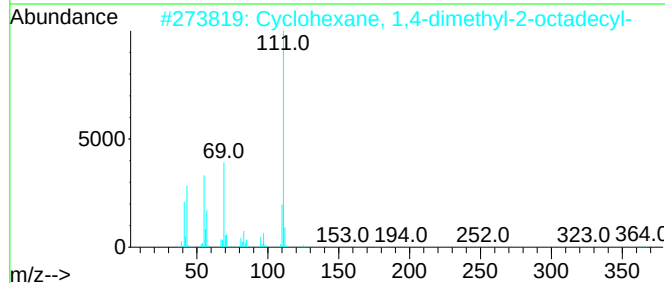
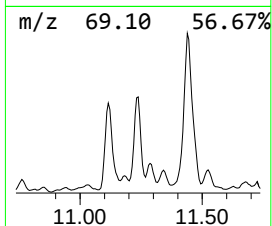
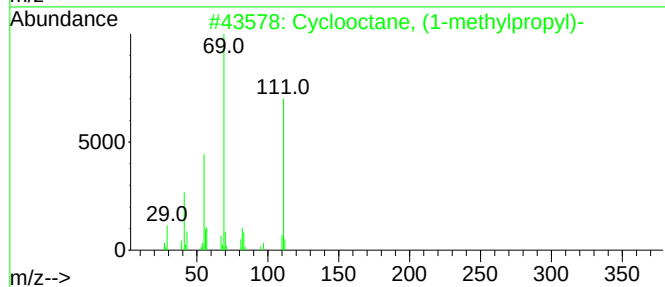
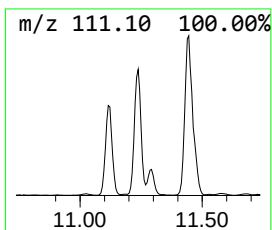
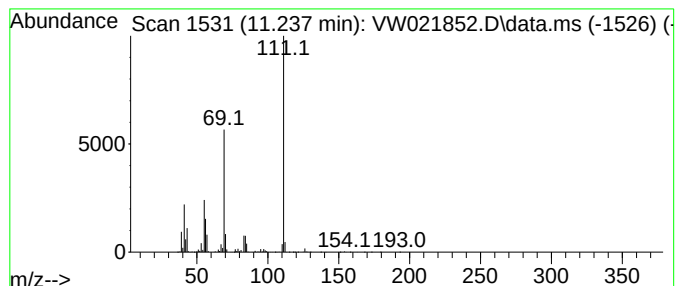
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Peak Number 20 (DEL) Alkane: Cyclic11.238 Concentration Rank 31

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 11.238 | 32.06 ug/L | 1059770 | Chlorobenzene-d5 | 11.634 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Cyclooctane, (1-methylpropyl)-      | 168 | C12H24  | 016538-89-9 | 59   |
| 2    |      | Cyclohexane, 1,4-dimethyl-2-octa... | 364 | C26H52  | 055282-02-5 | 56   |
| 3    |      | Cyclohexane, 1,1,3-trimethyl-       | 126 | C9H18   | 003073-66-3 | 46   |
| 4    |      | Cyclohexane, 1,1,3-trimethyl-       | 126 | C9H18   | 003073-66-3 | 46   |
| 5    |      | Pyrazol-4-amine, 1,5-dimethyl-      | 111 | C5H9N3  | 121983-36-6 | 45   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

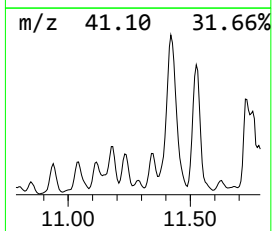
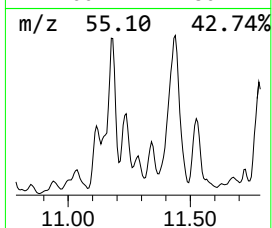
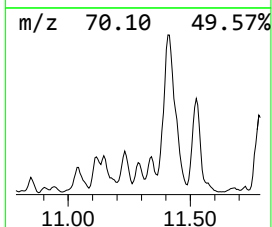
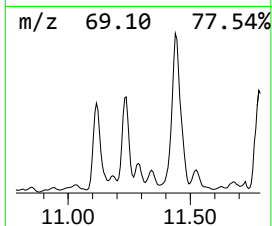
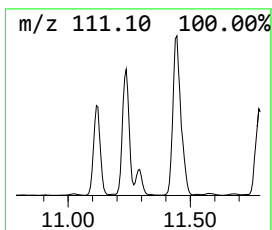
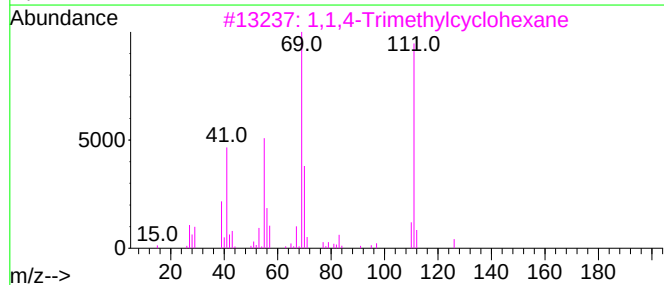
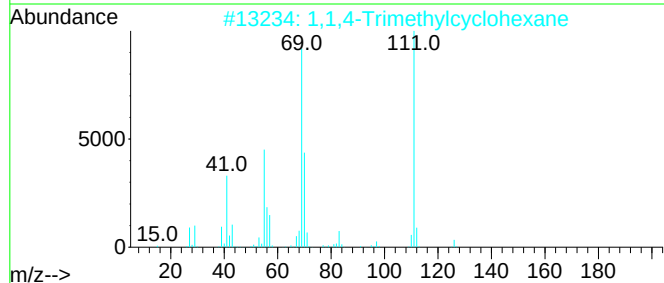
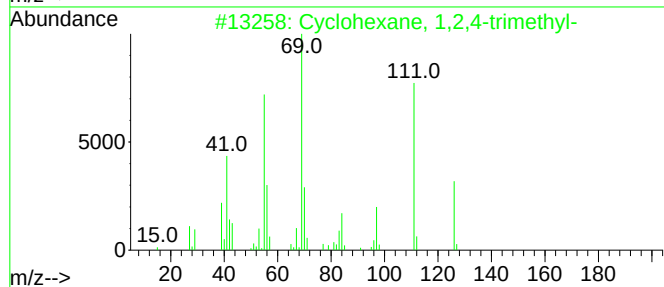
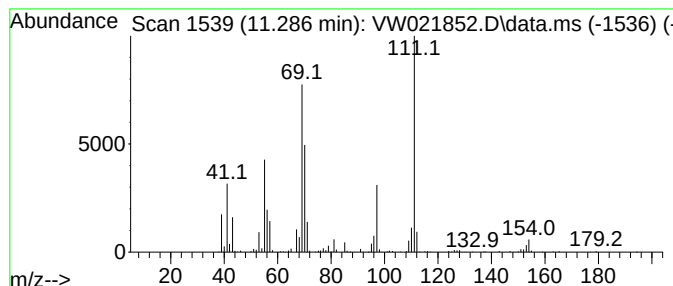
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Quant Title : SFAM01.0

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

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Peak Number 21 (DEL) Alkane: Cyclic11.286 Concentration Rank 62

| R.T.      | EstConc                        | Area   | Relative to ISTD |         |             | R.T.   |
|-----------|--------------------------------|--------|------------------|---------|-------------|--------|
| 11.286    | 6.50 ug/L                      | 214810 | Chlorobenzene-d5 |         |             | 11.634 |
| Hit# of 5 | Tentative ID                   |        | MW               | MolForm | CAS#        | Qual   |
| 1         | Cyclohexane, 1,2,4-trimethyl-  |        | 126              | C9H18   | 002234-75-5 | 72     |
| 2         | 1,1,4-Trimethylcyclohexane     |        | 126              | C9H18   | 007094-27-1 | 72     |
| 3         | 1,1,4-Trimethylcyclohexane     |        | 126              | C9H18   | 007094-27-1 | 59     |
| 4         | Cyclooctane, (1-methylpropyl)- |        | 168              | C12H24  | 016538-89-9 | 53     |
| 5         | Cyclooctane, tetradecyl-       |        | 308              | C22H44  | 149003-36-1 | 53     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

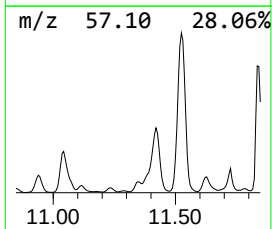
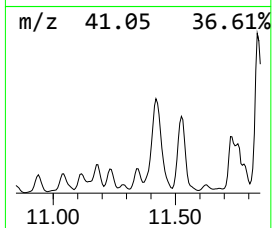
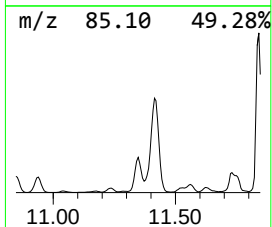
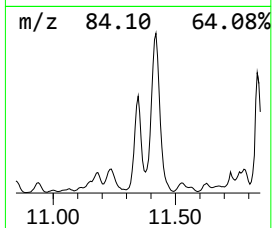
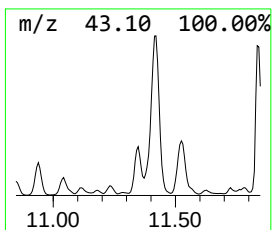
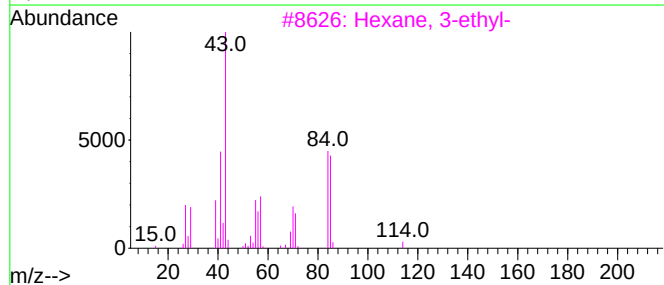
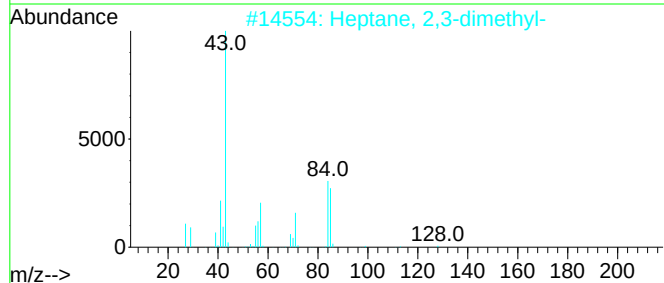
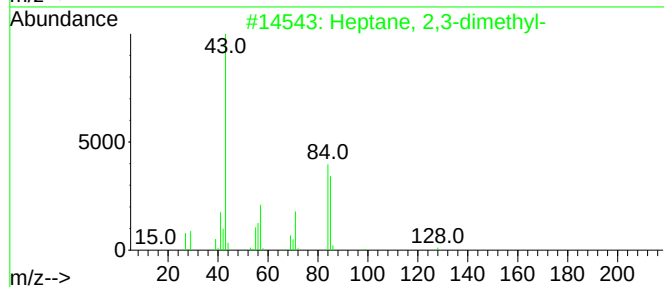
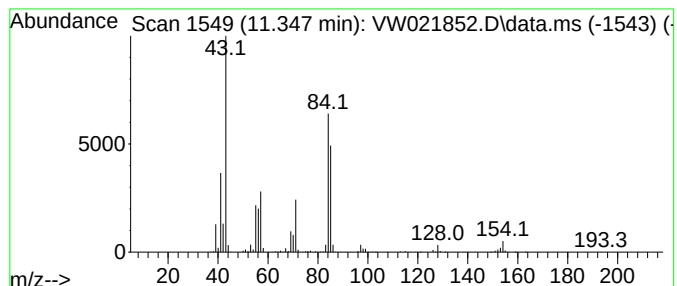
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Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 11.347 | 31.67 ug/L | 1046640 | Chlorobenzene-d5 | 11.634 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptane, 2,3-dimethyl- | 128 | C9H20   | 003074-71-3 | 87   |
| 2    |    |   | Heptane, 2,3-dimethyl- | 128 | C9H20   | 003074-71-3 | 83   |
| 3    |    |   | Hexane, 3-ethyl-       | 114 | C8H18   | 000619-99-8 | 83   |
| 4    |    |   | Heptane, 2,3-dimethyl- | 128 | C9H20   | 003074-71-3 | 76   |
| 5    |    |   | 1-Octene, 4-methyl-    | 126 | C9H18   | 013151-12-7 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

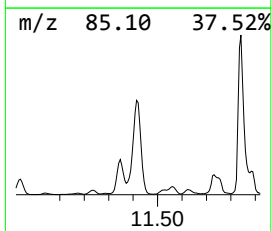
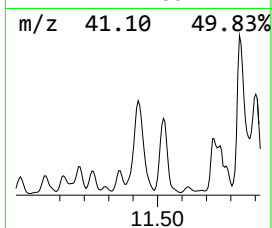
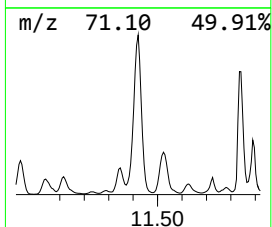
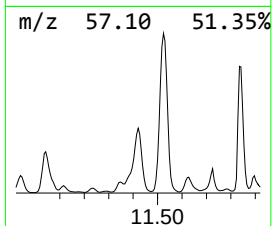
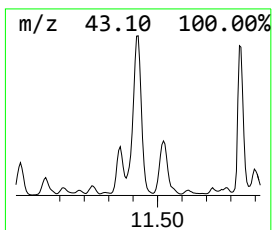
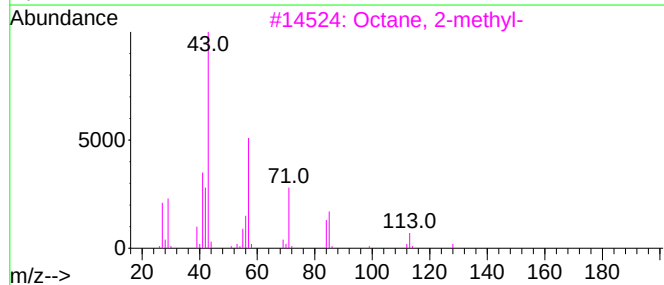
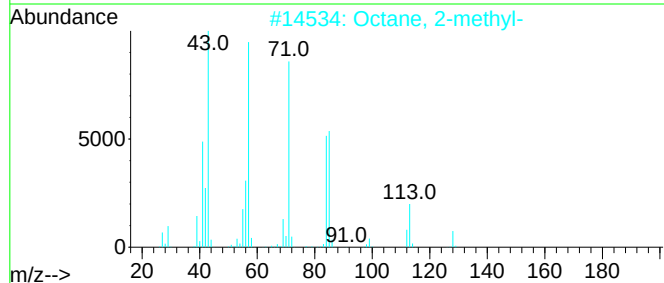
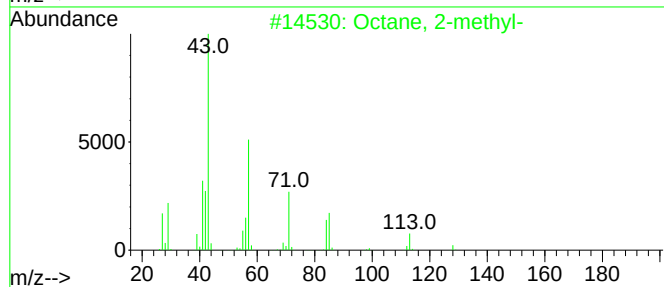
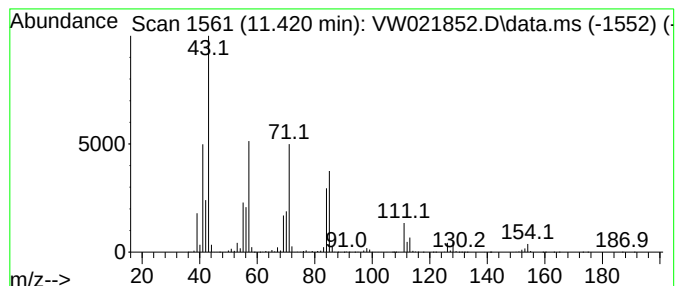
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Quant Title : SFAM01.0

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

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Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 11.420 | 206.29 ug/L | 6818500 | Chlorobenzene-d5 | 11.634 |

| Hit# | of                       | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|--------------------------|---|--------------|-----|---------|-------------|------|
| 1    | Octane, 2-methyl-        |   |              | 128 | C9H20   | 003221-61-2 | 74   |
| 2    | Octane, 2-methyl-        |   |              | 128 | C9H20   | 003221-61-2 | 64   |
| 3    | Octane, 2-methyl-        |   |              | 128 | C9H20   | 003221-61-2 | 59   |
| 4    | Octane, 4-methyl-        |   |              | 128 | C9H20   | 002216-34-4 | 52   |
| 5    | Hexane, 2,3,4-trimethyl- |   |              | 128 | C9H20   | 000921-47-1 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

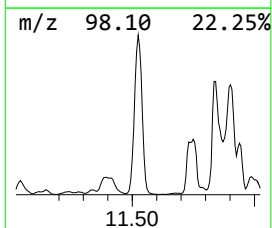
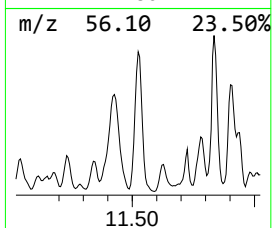
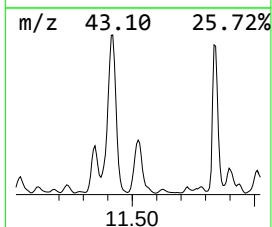
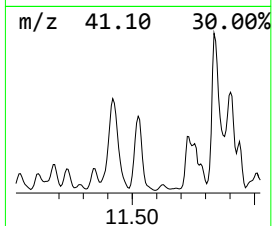
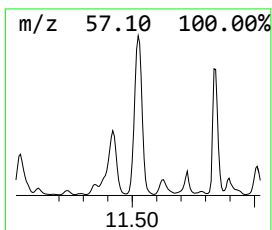
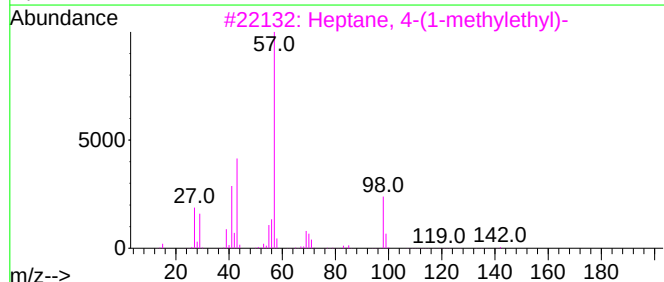
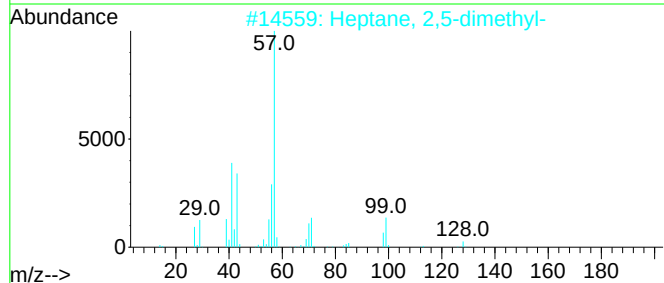
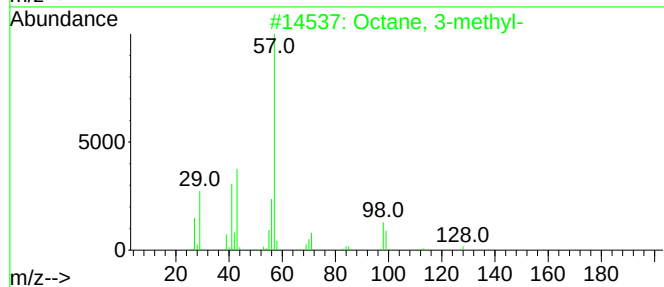
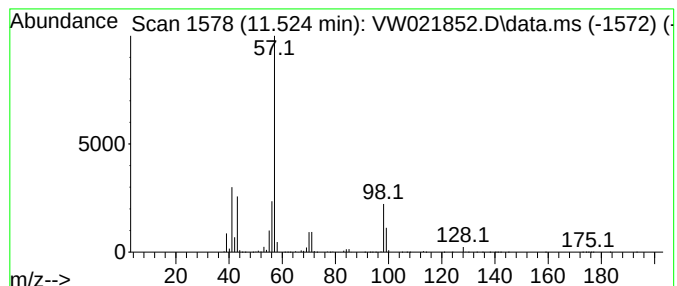
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 11.524 | 112.43 ug/L | 3716140 | Chlorobenzene-d5 | 11.634 |

| Hit# | of 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------------|-----|---------|-------------|------|
| 1    |      | Octane, 3-methyl-           | 128 | C9H20   | 002216-33-3 | 83   |
| 2    |      | Heptane, 2,5-dimethyl-      | 128 | C9H20   | 002216-30-0 | 76   |
| 3    |      | Heptane, 4-(1-methylethyl)- | 142 | C10H22  | 052896-87-4 | 72   |
| 4    |      | Octane, 3-methyl-           | 128 | C9H20   | 002216-33-3 | 72   |
| 5    |      | Heptane, 4-propyl-          | 142 | C10H22  | 003178-29-8 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

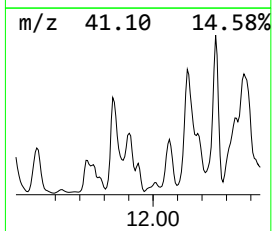
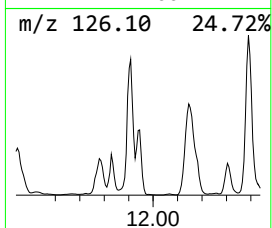
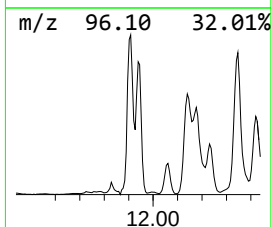
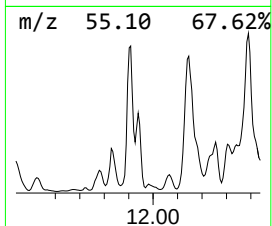
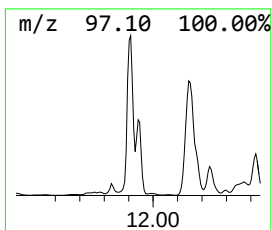
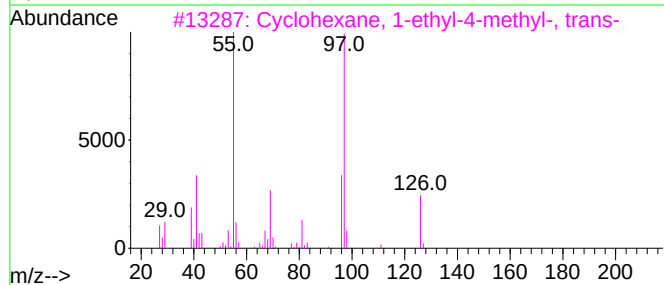
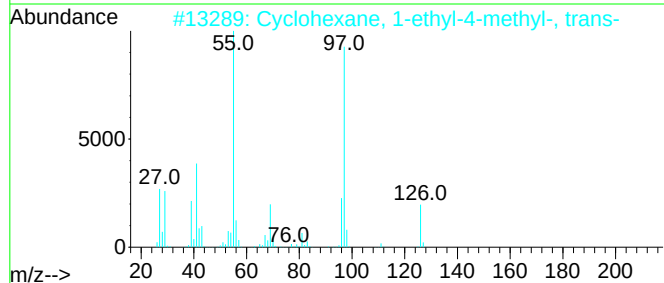
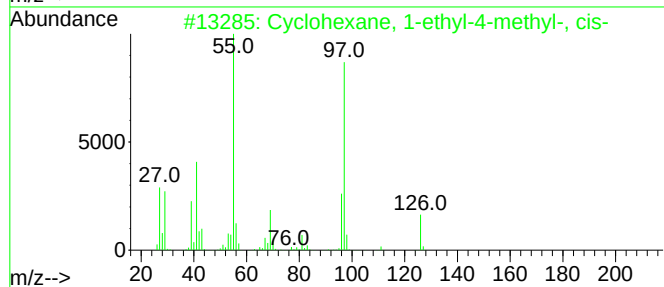
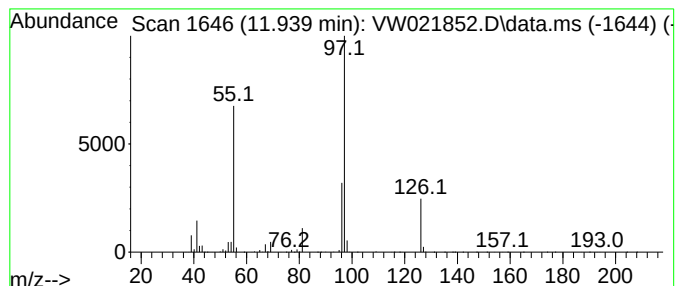
Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

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Peak Number 25 (DEL) Alkane: Cyclic11.939 Concentration Rank 15

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 11.939    | 60.66 ug/L                          | 2004810 | Chlorobenzene-d5 | 11.634      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | Cyclohexane, 1-ethyl-4-methyl-, ... | 126     | C9H18            | 004926-78-7 | 78   |
| 2         | Cyclohexane, 1-ethyl-4-methyl-, ... | 126     | C9H18            | 006236-88-0 | 78   |
| 3         | Cyclohexane, 1-ethyl-4-methyl-, ... | 126     | C9H18            | 006236-88-0 | 72   |
| 4         | 1-Ethyl-4-methylcyclohexane         | 126     | C9H18            | 003728-56-1 | 56   |
| 5         | Furan, 2,3-dihydro-4-(1-methylpr... | 126     | C8H14O           | 034379-54-9 | 53   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

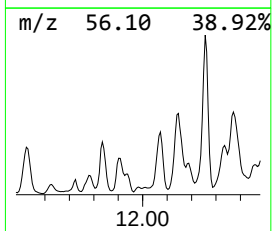
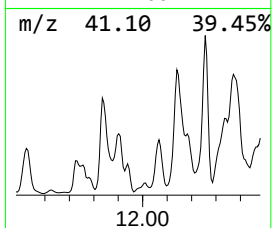
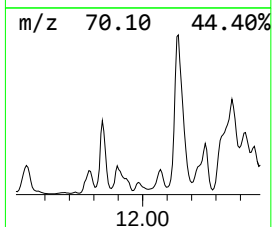
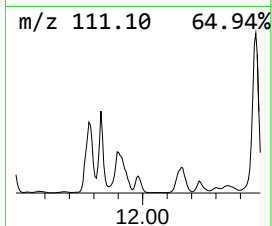
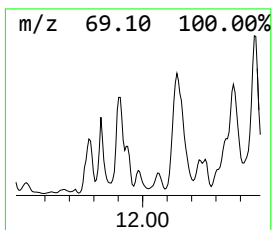
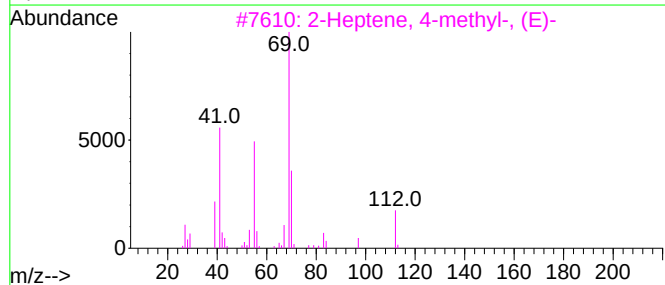
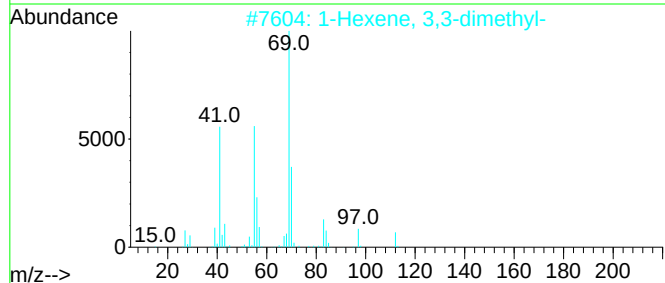
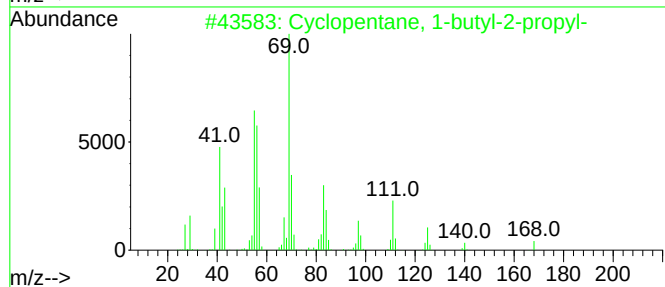
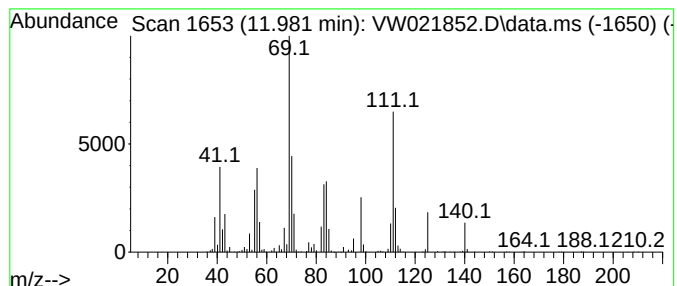
Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

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Peak Number 26 (DEL) Alkane: Cyclic11.981 Concentration Rank 58

| R.T.      | EstConc                         | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|---------------------------------|--------|------------------|---------|-------------|------|
| 11.981    | 8.20 ug/L                       | 271132 | Chlorobenzene-d5 |         | 11.634      |      |
| Hit# of 5 | Tentative ID                    |        | MW               | MolForm | CAS#        | Qual |
| 1         | Cyclopentane, 1-butyl-2-propyl- |        | 168              | C12H24  | 062199-50-2 | 53   |
| 2         | 1-Hexene, 3,3-dimethyl-         |        | 112              | C8H16   | 003404-77-1 | 47   |
| 3         | 2-Heptene, 4-methyl-, (E)-      |        | 112              | C8H16   | 066225-17-0 | 43   |
| 4         | 3-Hexene, 2-methyl-, (E)-       |        | 98               | C7H14   | 000692-24-0 | 43   |
| 5         | 2-Methyl-2-heptene              |        | 112              | C8H16   | 000627-97-4 | 43   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

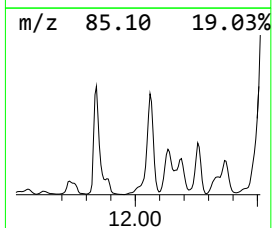
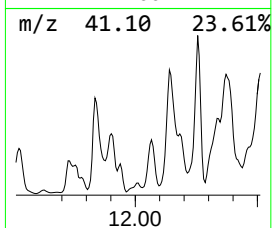
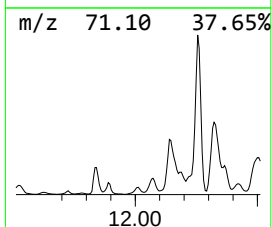
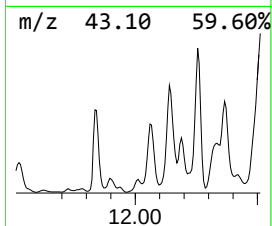
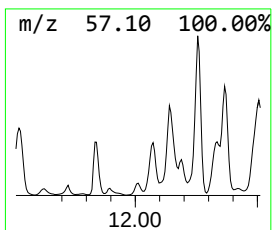
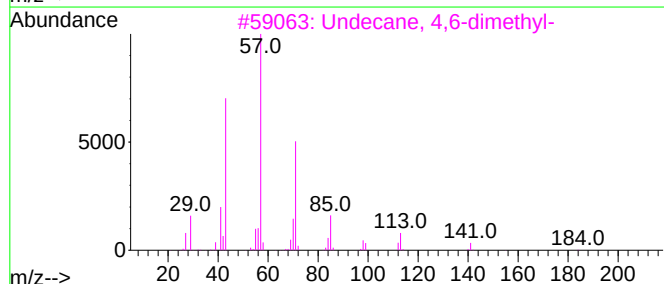
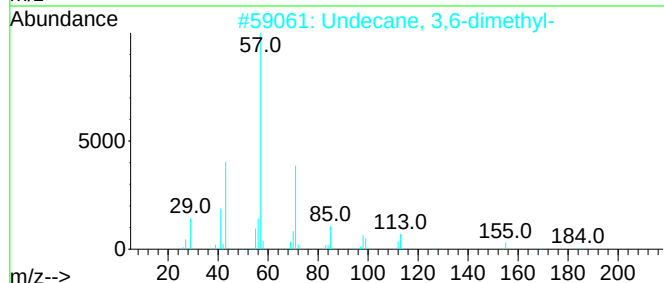
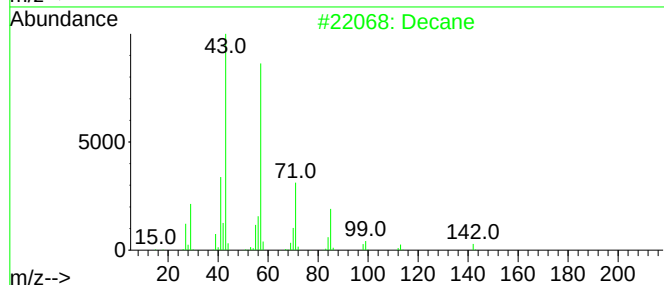
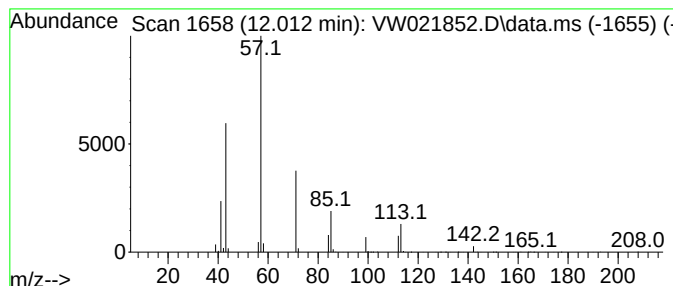
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 59

| R.T.   | EstConc   | Area   | Relative to ISTD | R.T.   |
|--------|-----------|--------|------------------|--------|
| 12.012 | 7.47 ug/L | 246736 | Chlorobenzene-d5 | 11.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Decane                              | 142 | C10H22    | 000124-18-5  | 72   |
| 2    |    |   | Undecane, 3,6-dimethyl-             | 184 | C13H28    | 017301-28-9  | 72   |
| 3    |    |   | Undecane, 4,6-dimethyl-             | 184 | C13H28    | 017312-82-2  | 64   |
| 4    |    |   | Decane, 2-methyl-                   | 156 | C11H24    | 006975-98-0  | 64   |
| 5    |    |   | Sulfurous acid, decyl 2-ethylhex... | 334 | C18H38O3S | 1000309-19-3 | 53   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

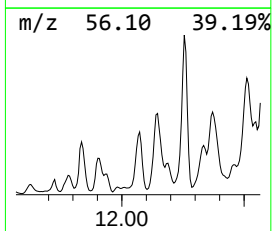
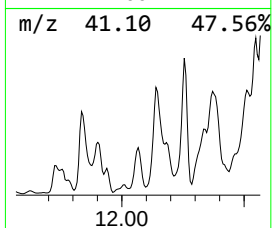
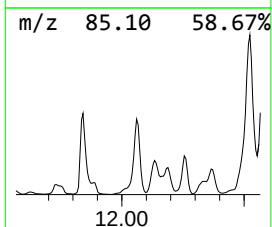
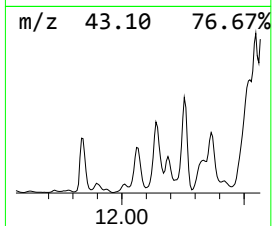
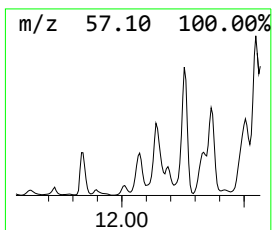
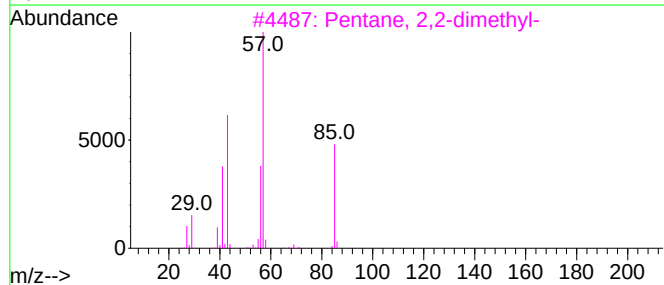
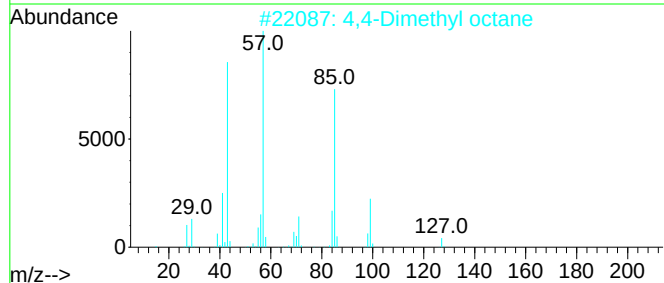
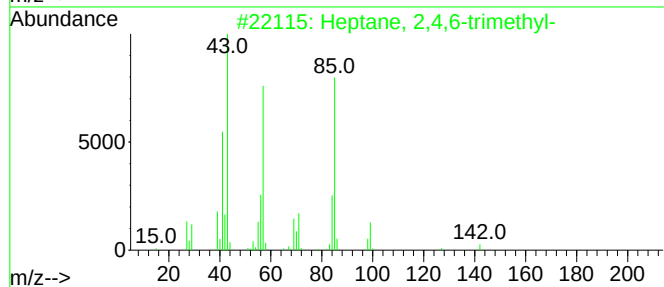
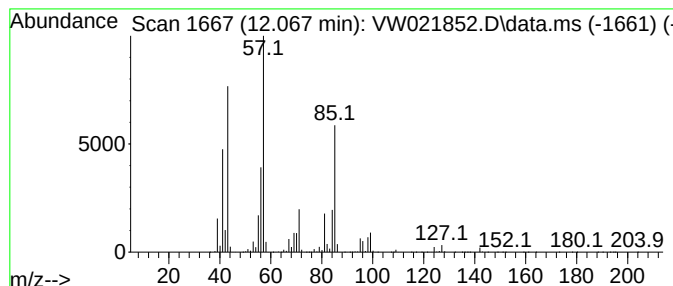
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Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 12.067 | 112.35 ug/L | 3713440 | Chlorobenzene-d5 | 11.634 |

| Hit# | of 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------------|-----|---------|-------------|------|
| 1    |      | Heptane, 2,4,6-trimethyl-    | 142 | C10H22  | 002613-61-8 | 64   |
| 2    |      | 4,4-Dimethyl octane          | 142 | C10H22  | 015869-95-1 | 47   |
| 3    |      | Pentane, 2,2-dimethyl-       | 100 | C7H16   | 000590-35-2 | 47   |
| 4    |      | Hexane, 2,2,3,3-tetramethyl- | 142 | C10H22  | 013475-81-5 | 43   |
| 5    |      | Octane, 2,2,6-trimethyl-     | 156 | C11H24  | 062016-28-8 | 43   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

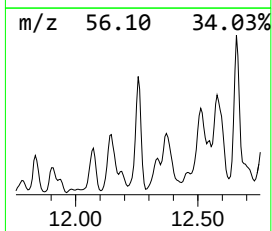
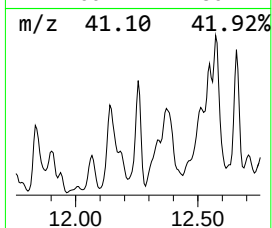
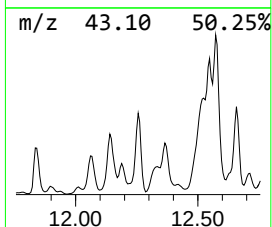
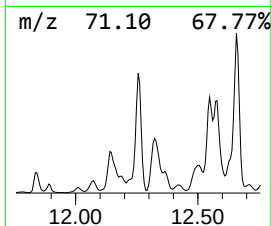
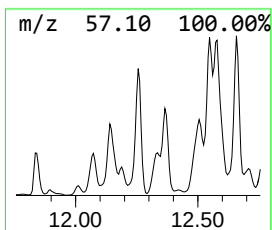
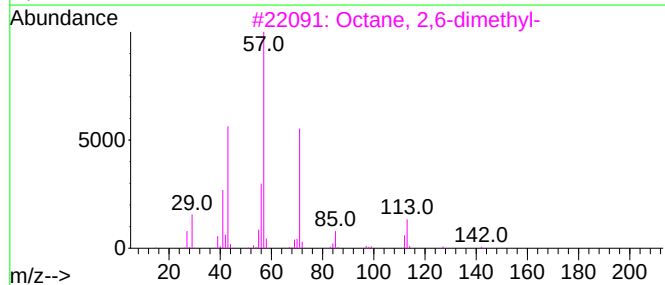
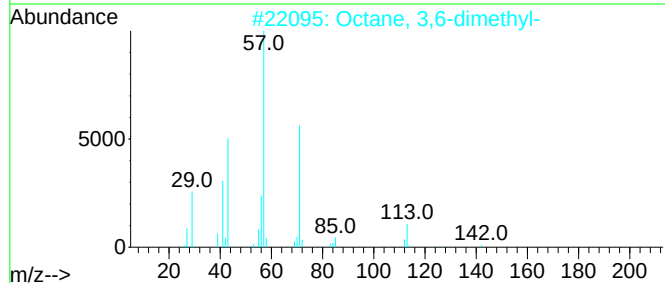
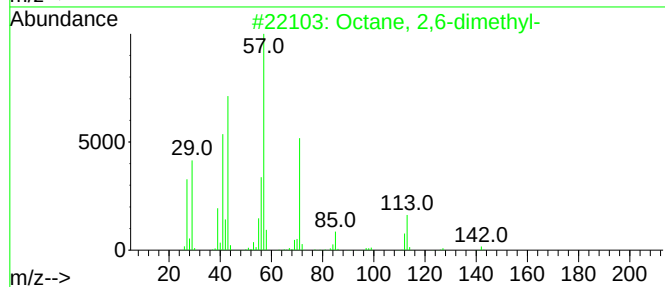
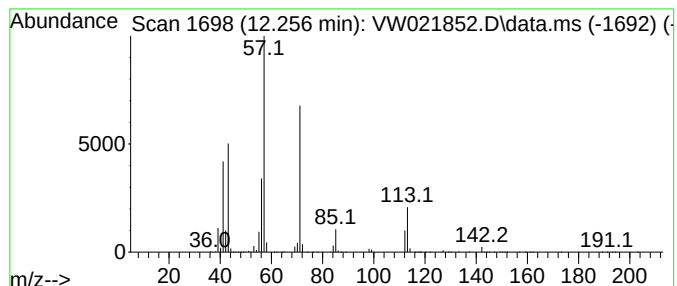
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Quant Title : SFAM01.0

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

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Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 12.256 | 277.95 ug/L | 9187030 | Chlorobenzene-d5 | 11.634 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 90   |
| 2    |    |   | Octane, 3,6-dimethyl- | 142 | C10H22  | 015869-94-0 | 90   |
| 3    |    |   | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 87   |
| 4    |    |   | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 87   |
| 5    |    |   | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 80   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

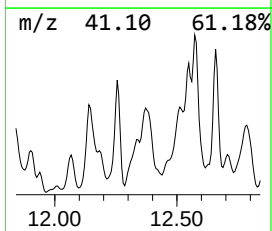
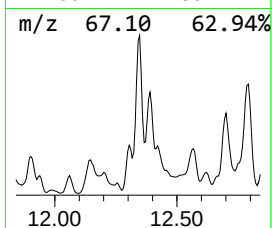
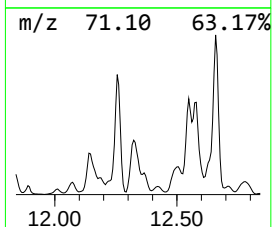
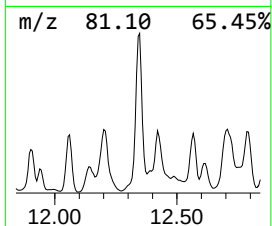
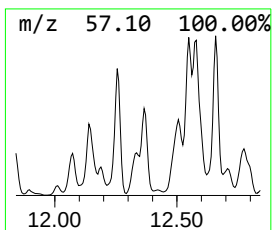
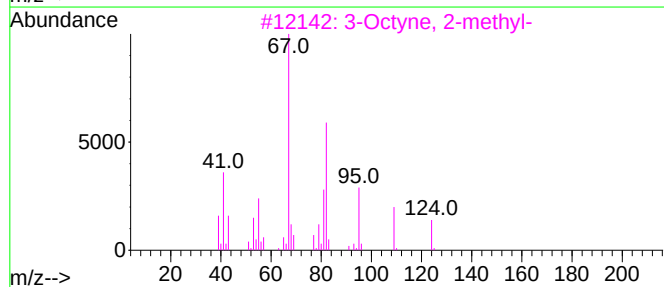
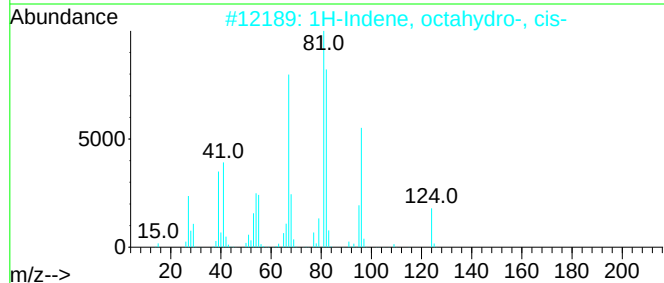
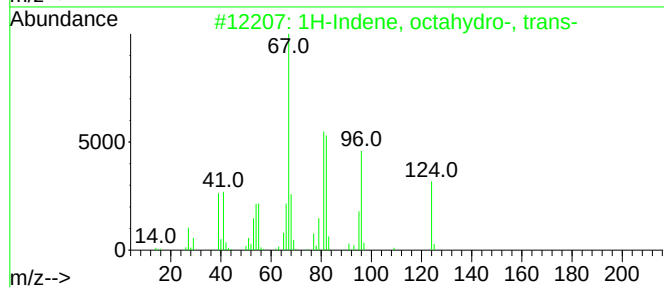
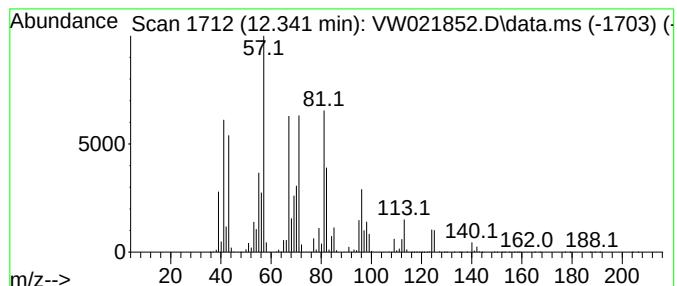
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Quant Title : SFAM01.0

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

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Peak Number 30 unknown-01 Concentration Rank 5

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 12.341 | 275.05 ug/L | 9090940 | Chlorobenzene-d5 | 11.634 |

| Hit# | of 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------|-----|---------|-------------|------|
| 1    |      | 1H-Indene, octahydro-, trans- | 124 | C9H16   | 003296-50-2 | 49   |
| 2    |      | 1H-Indene, octahydro-, cis-   | 124 | C9H16   | 004551-51-3 | 45   |
| 3    |      | 3-Octyne, 2-methyl-           | 124 | C9H16   | 055402-15-8 | 43   |
| 4    |      | 1H-Indene, octahydro-, trans- | 124 | C9H16   | 003296-50-2 | 43   |
| 5    |      | 3-Hexen-1-ol                  | 100 | C6H12O  | 000544-12-7 | 38   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

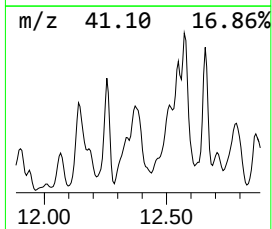
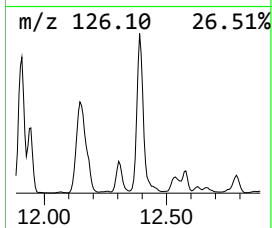
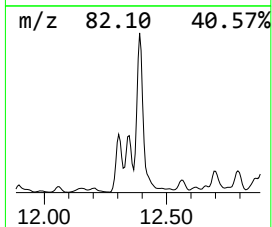
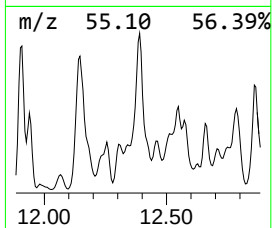
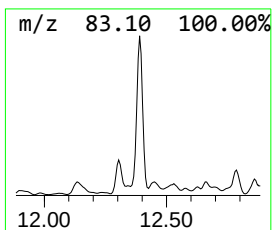
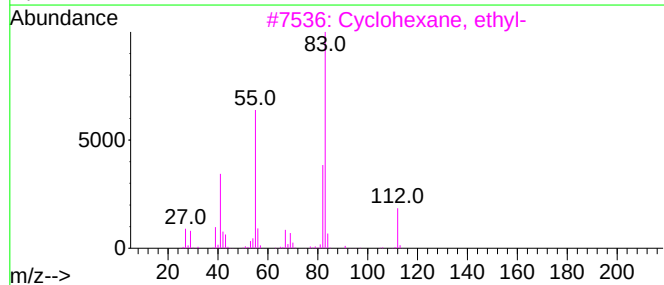
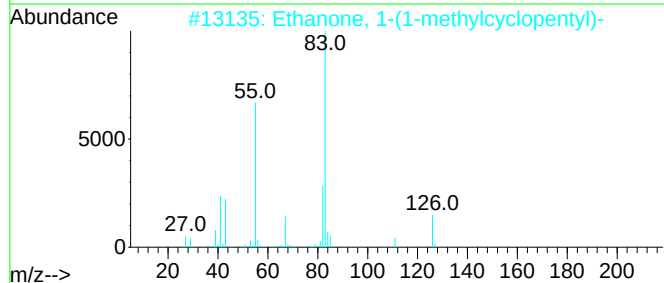
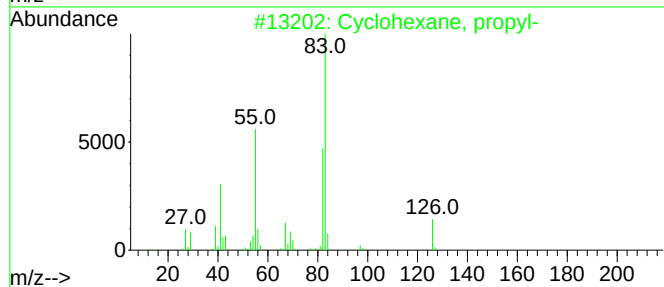
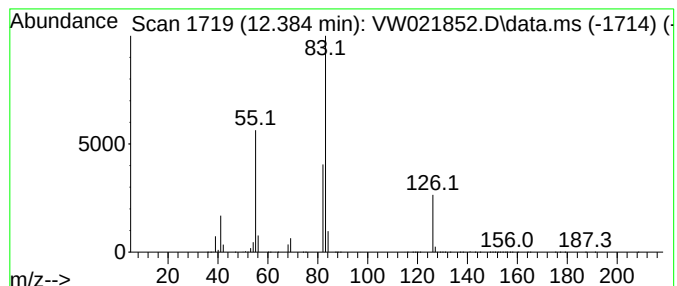
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Quant Title : SFAM01.0

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

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Peak Number 31 (DEL) Alkane: Cyclic12.384 Concentration Rank 1

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 12.384 | 395.20 ug/L | 13062200 | Chlorobenzene-d5 | 11.634 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Cyclohexane, propyl-                | 126 | C9H18    | 001678-92-8  | 90   |
| 2    |    |   | Ethanone, 1-(1-methylcyclopentyl)-  | 126 | C8H14O   | 013388-93-7  | 78   |
| 3    |    |   | Cyclohexane, ethyl-                 | 112 | C8H16    | 001678-91-7  | 72   |
| 4    |    |   | Cyclohexane, propyl-                | 126 | C9H18    | 001678-92-8  | 59   |
| 5    |    |   | Acetaldehyde 2E-hexenyl 3Z-hexen... | 226 | C14H26O2 | 1000431-45-5 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

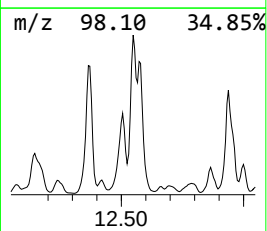
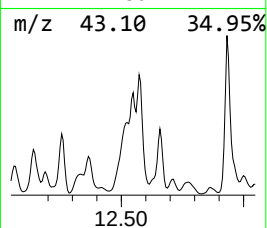
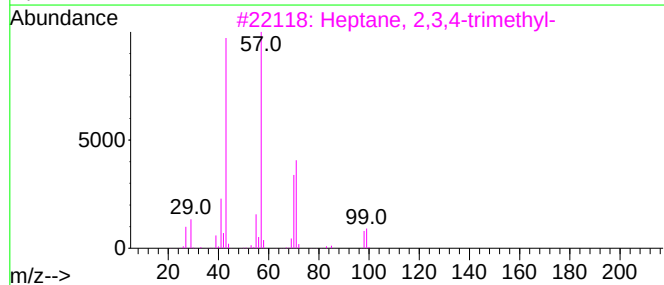
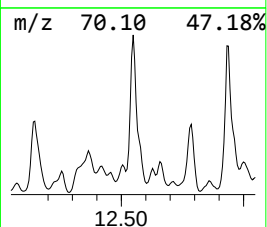
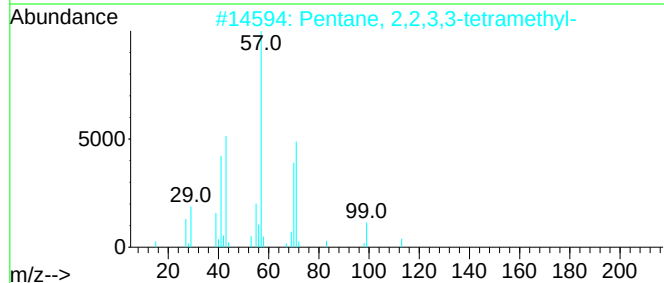
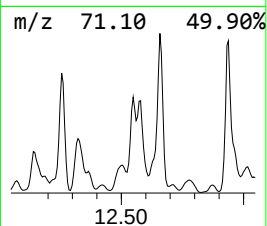
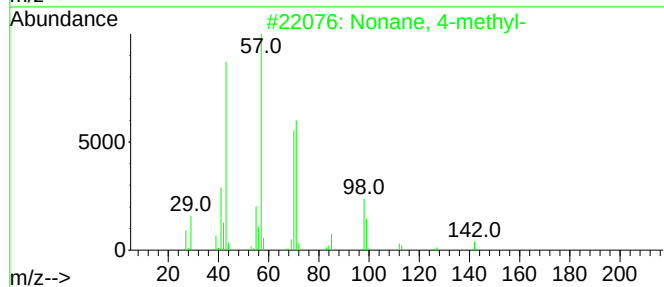
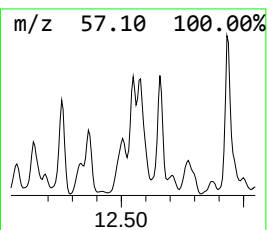
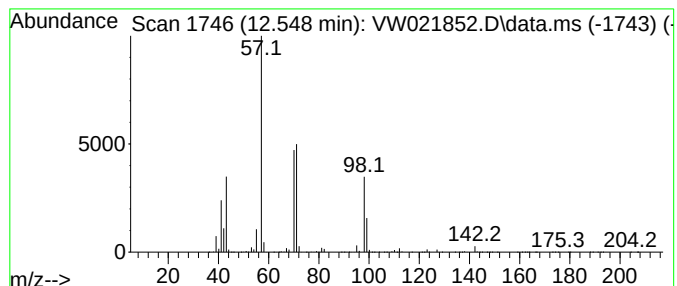
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Quant Title : SFAM01.0

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

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Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 12.548 | 346.36 ug/L | 11448100 | Chlorobenzene-d5 | 11.634 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 74   |
| 2    |    |   | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 64   |
| 3    |    |   | Heptane, 2,3,4-trimethyl-     | 142 | C10H22  | 052896-95-4 | 64   |
| 4    |    |   | Octane, 2,5-dimethyl-         | 142 | C10H22  | 015869-89-3 | 53   |
| 5    |    |   | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

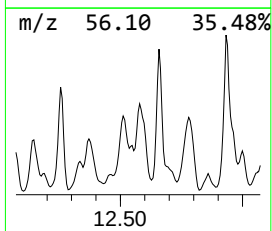
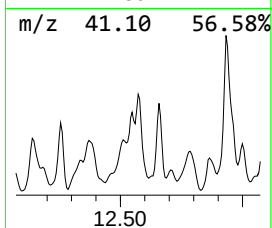
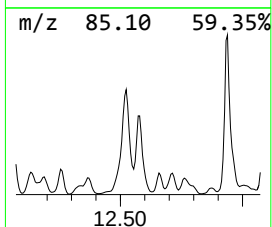
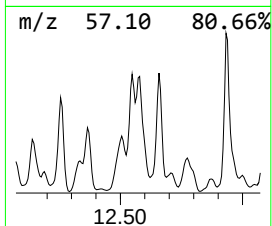
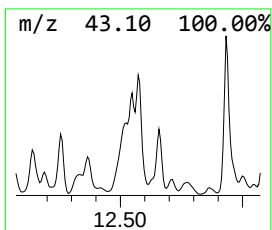
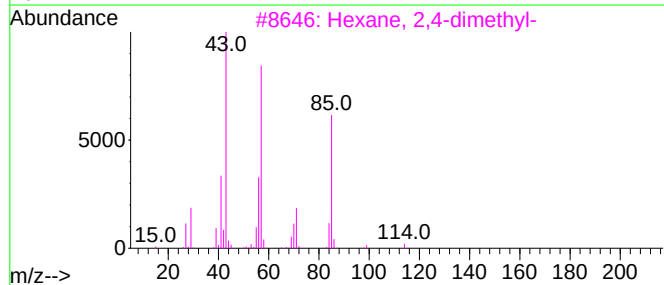
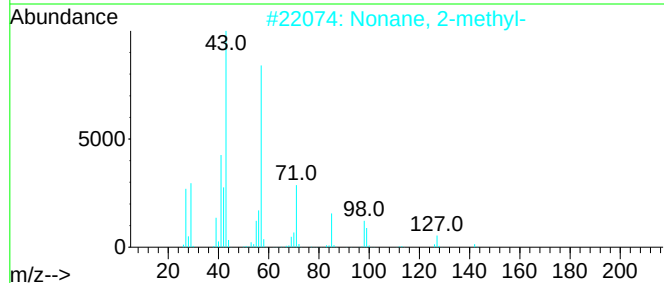
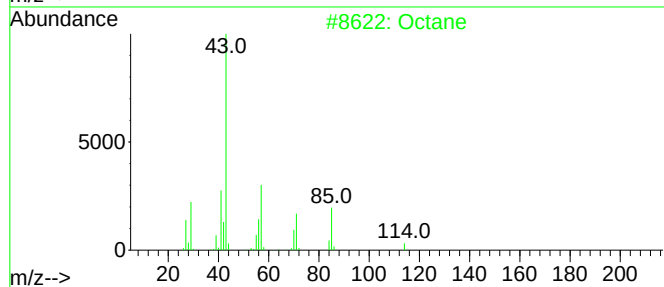
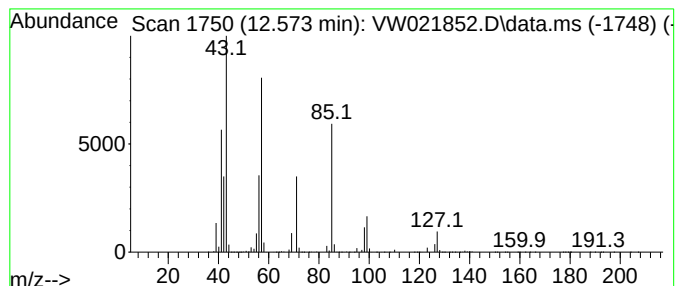
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TIC Integration Parameters: LSCINT.P

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Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 12.573 | 391.40 ug/L | 12936600 | Chlorobenzene-d5 | 11.634 |

| Hit# | of                      | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|-------------------------|---|--------------|-----|---------|-------------|------|
| 1    | Octane                  |   |              | 114 | C8H18   | 000111-65-9 | 59   |
| 2    | Nonane, 2-methyl-       |   |              | 142 | C10H22  | 000871-83-0 | 53   |
| 3    | Hexane, 2,4-dimethyl-   |   |              | 114 | C8H18   | 000589-43-5 | 50   |
| 4    | 4,4-Dimethyl octane     |   |              | 142 | C10H22  | 015869-95-1 | 50   |
| 5    | Undecane, 5,7-dimethyl- |   |              | 184 | C13H28  | 017312-83-3 | 47   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

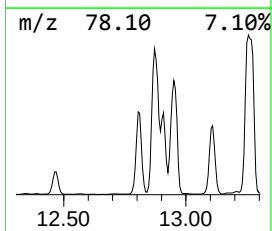
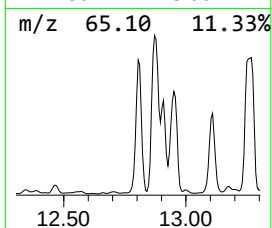
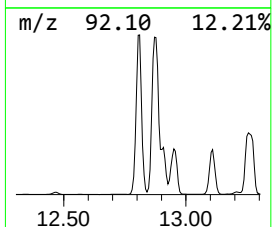
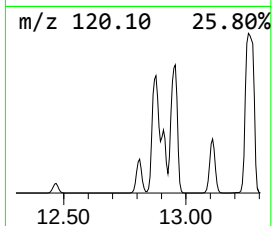
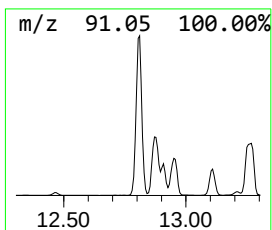
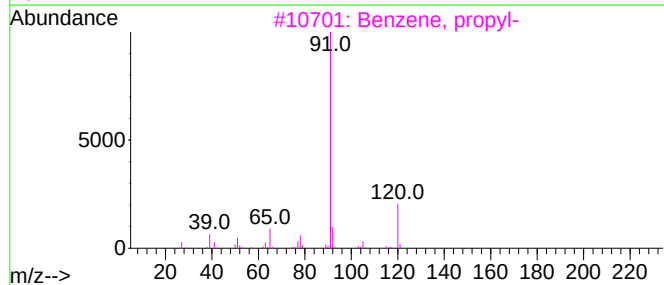
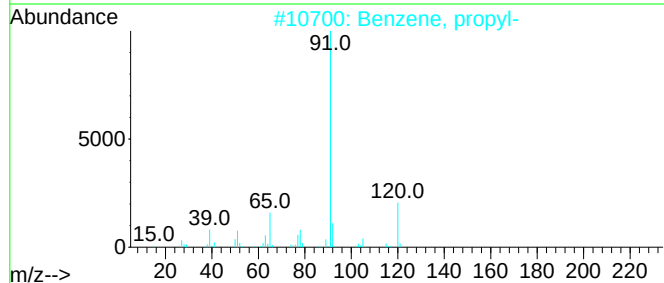
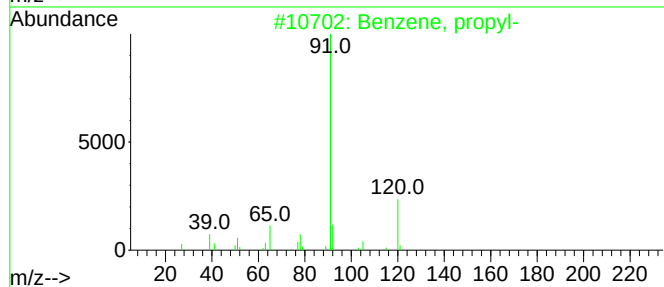
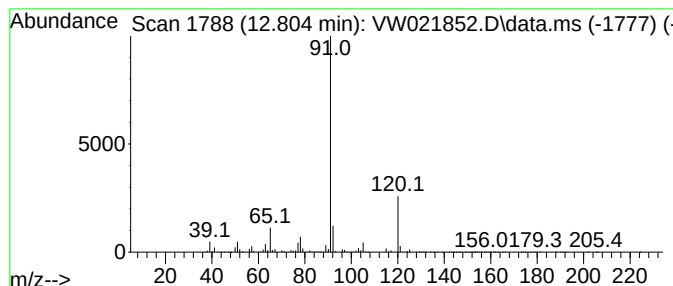
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Quant Title : SFAM01.0

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

\*\*\*\*\*  
Peak Number 34 Benzene, propyl- Concentration Rank 12

| R.T.   | EstConc    | Area     | Relative to ISTD       | R.T.   |
|--------|------------|----------|------------------------|--------|
| 12.804 | 96.76 ug/L | 27538700 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, propyl-          | 120 | C9H12   | 000103-65-1 | 93   |
| 2    |    |   | Benzene, propyl-          | 120 | C9H12   | 000103-65-1 | 93   |
| 3    |    |   | Benzene, propyl-          | 120 | C9H12   | 000103-65-1 | 87   |
| 4    |    |   | Benzene, propyl-          | 120 | C9H12   | 000103-65-1 | 87   |
| 5    |    |   | N-Benzyl-2-phenethylamine | 211 | C15H17N | 003647-71-0 | 72   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

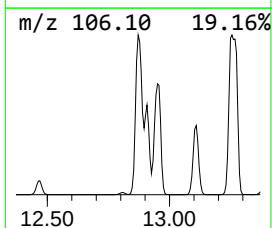
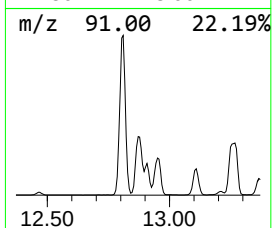
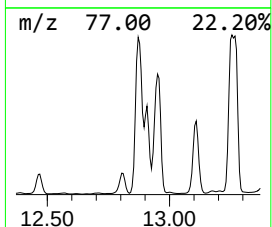
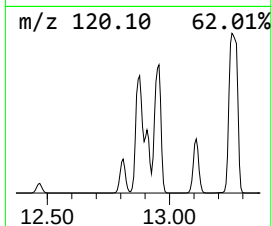
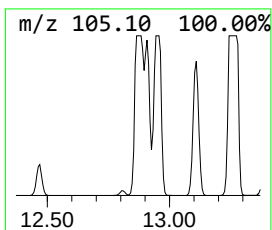
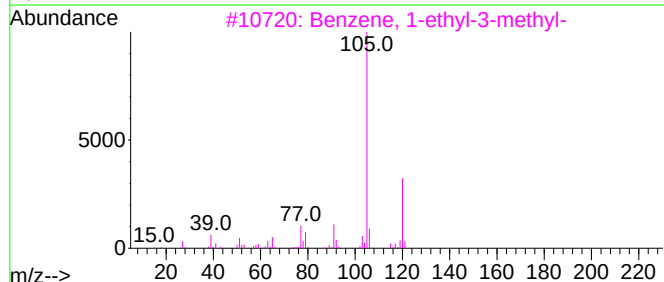
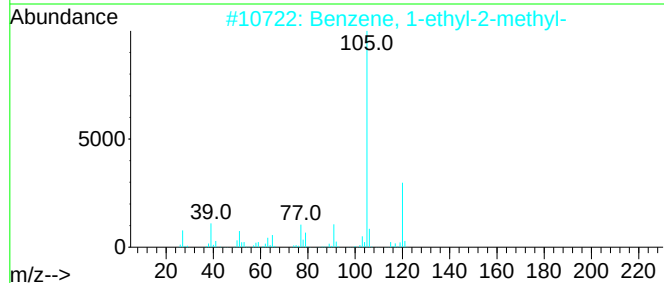
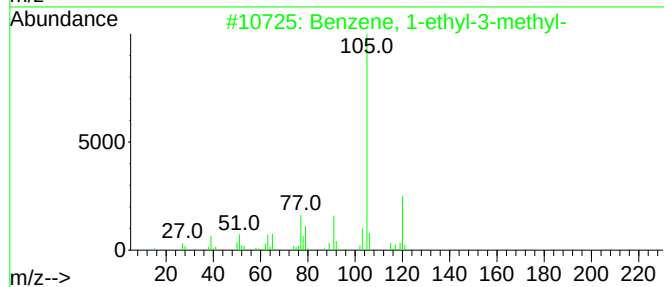
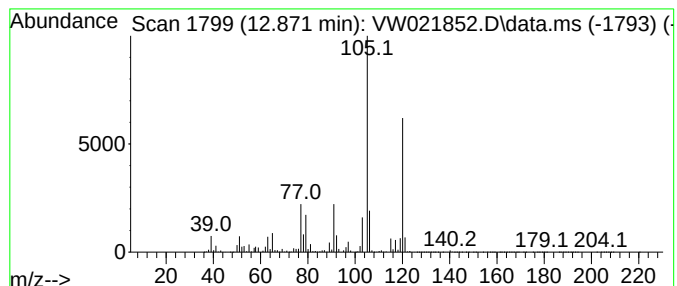
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Quant Title : SFAM01.0

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

\*\*\*\*\*  
Peak Number 35 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

| R.T.   | EstConc     | Area     | Relative to ISTD       | R.T.   |
|--------|-------------|----------|------------------------|--------|
| 12.871 | 269.78 ug/L | 76783300 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 81   |
| 2    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 81   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 81   |
| 4    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 81   |
| 5    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 81   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

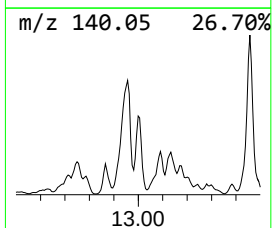
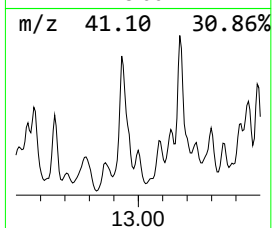
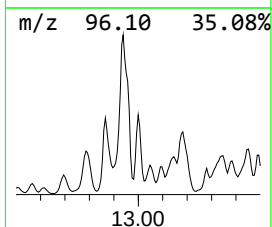
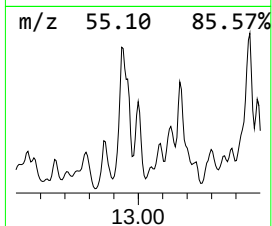
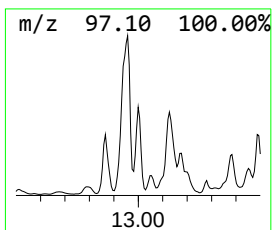
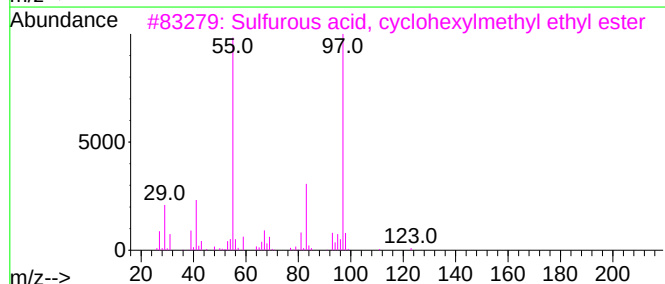
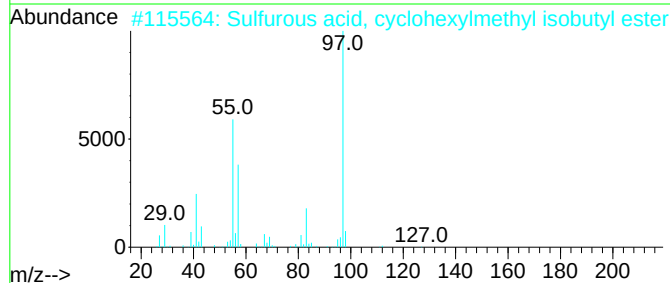
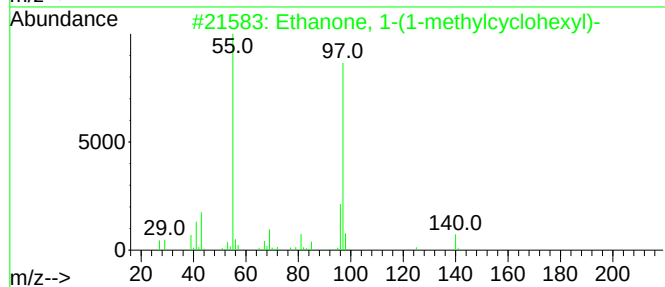
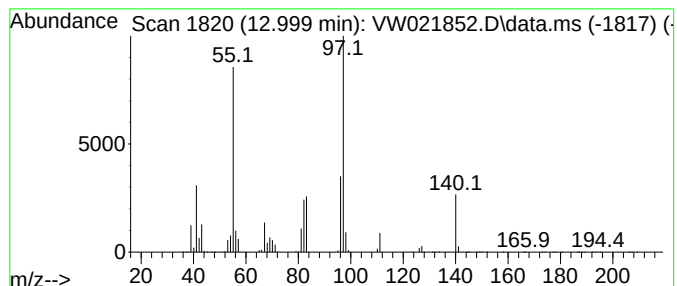
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Quant Title : SFAM01.0

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

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Peak Number 36 Ethanone, 1-(1-methylcyclohexyl) Concentration Rank 42

| R.T.    | EstConc    | Area                                | Relative to ISTD       |           |              | R.T.   |
|---------|------------|-------------------------------------|------------------------|-----------|--------------|--------|
| 12.999  | 23.71 ug/L | 6747850                             | 1,4-Dichlorobenzene-d4 |           |              | 13.560 |
| Hit# of | 5          | Tentative ID                        | MW                     | MolForm   | CAS#         | Qual   |
| 1       |            | Ethanone, 1-(1-methylcyclohexyl)-   | 140                    | C9H16O    | 002890-62-2  | 59     |
| 2       |            | Sulfurous acid, cyclohexylmethyl... | 234                    | C11H22O3S | 1000309-21-3 | 53     |
| 3       |            | Sulfurous acid, cyclohexylmethyl... | 206                    | C9H18O3S  | 1000309-21-2 | 53     |
| 4       |            | Cyclohexane, 1-ethyl-4-methyl-, ... | 126                    | C9H18     | 006236-88-0  | 53     |
| 5       |            | Cyclohexane, 1-isopropyl-1-methyl-  | 140                    | C10H20    | 016580-26-0  | 50     |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

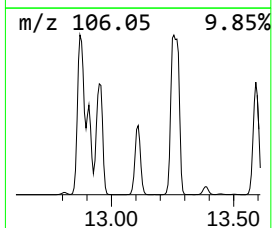
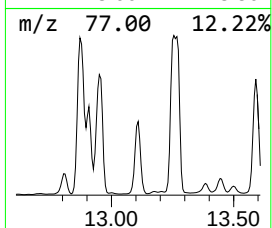
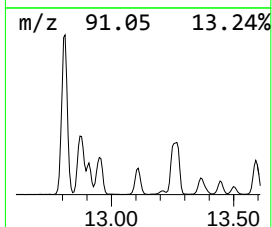
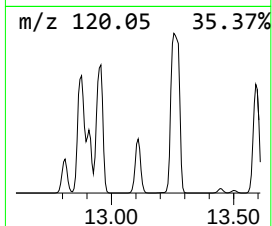
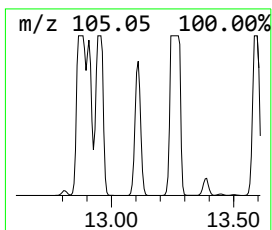
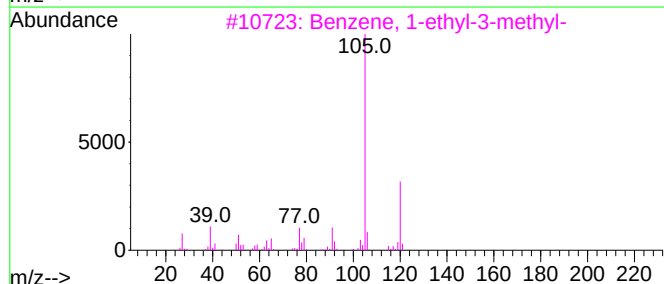
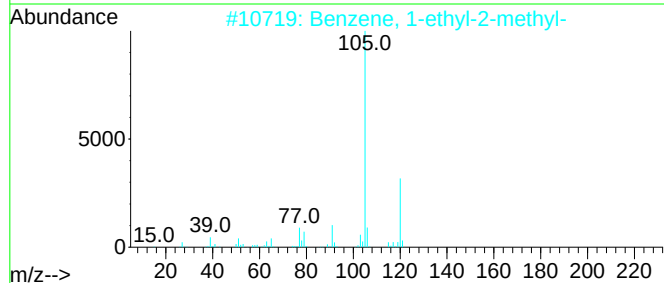
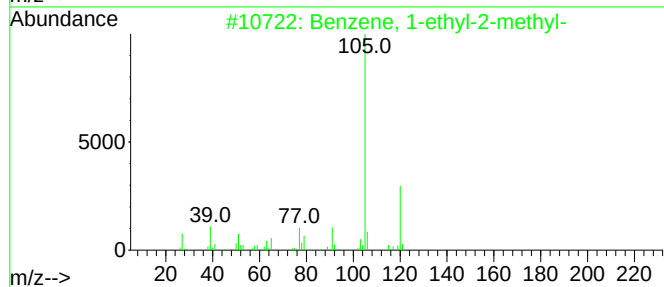
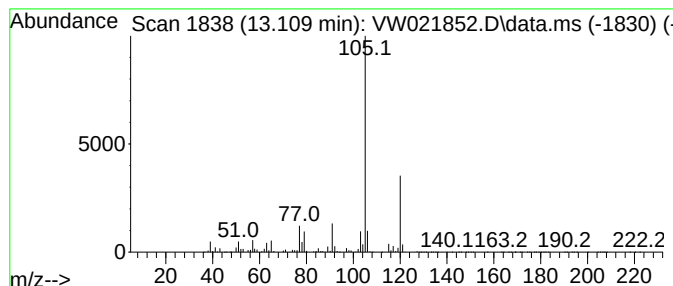
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Quant Title : SFAM01.0

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

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Peak Number 37 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

| R.T.   | EstConc     | Area     | Relative to ISTD       | R.T.   |
|--------|-------------|----------|------------------------|--------|
| 13.109 | 165.83 ug/L | 47195800 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 2    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 4    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |
| 5    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

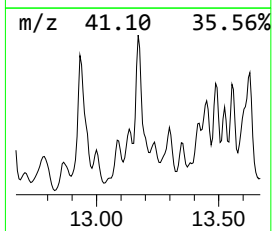
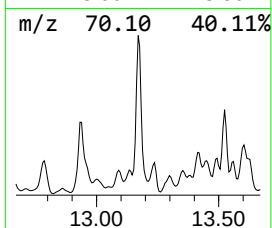
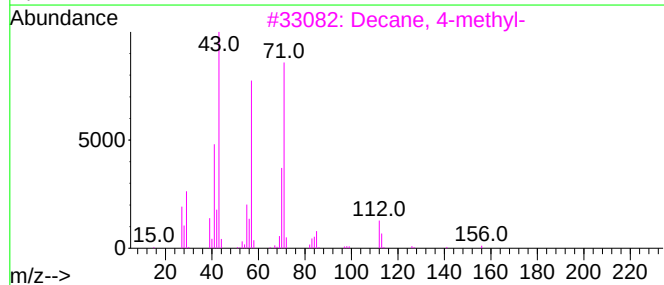
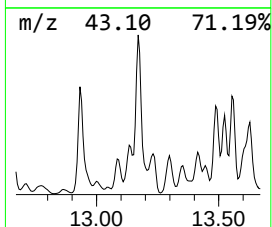
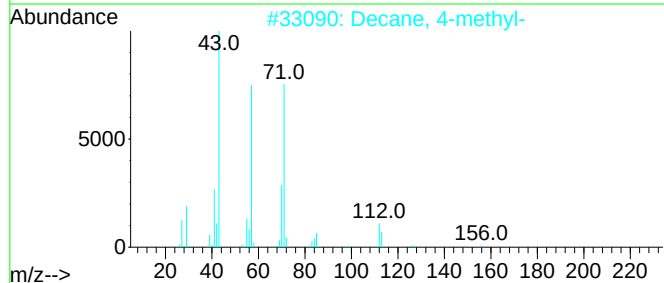
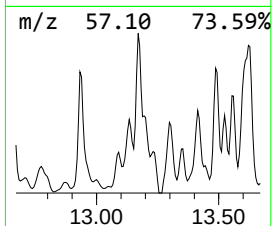
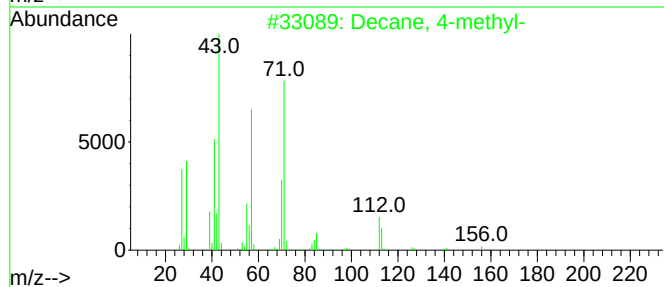
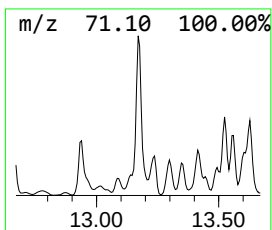
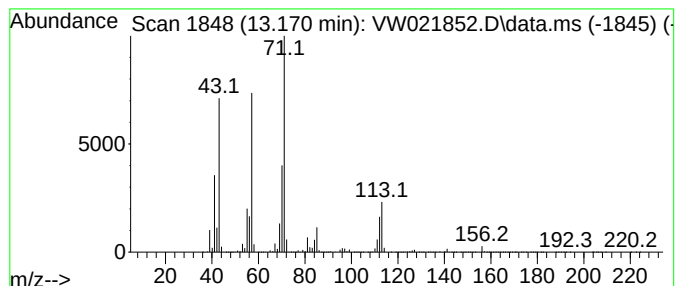
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Quant Title : SFAM01.0

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

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Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.      | EstConc                   | Area     | Relative to ISTD       |         |             | R.T.   |
|-----------|---------------------------|----------|------------------------|---------|-------------|--------|
| 13.170    | 83.23 ug/L                | 23688600 | 1,4-Dichlorobenzene-d4 |         |             | 13.560 |
| Hit# of 5 | Tentative ID              |          | MW                     | MolForm | CAS#        | Qual   |
| 1         | Decane, 4-methyl-         |          | 156                    | C11H24  | 002847-72-5 | 90     |
| 2         | Decane, 4-methyl-         |          | 156                    | C11H24  | 002847-72-5 | 87     |
| 3         | Decane, 4-methyl-         |          | 156                    | C11H24  | 002847-72-5 | 83     |
| 4         | Heptane, 3,3,5-trimethyl- |          | 142                    | C10H22  | 007154-80-5 | 72     |
| 5         | Heptane, 3,3,5-trimethyl- |          | 142                    | C10H22  | 007154-80-5 | 72     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

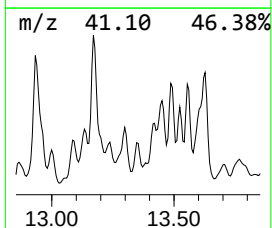
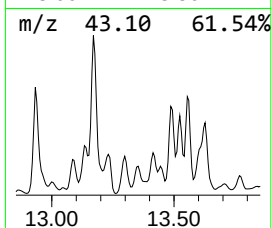
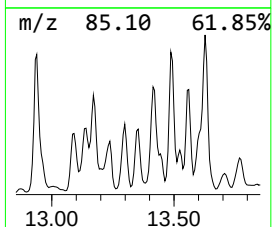
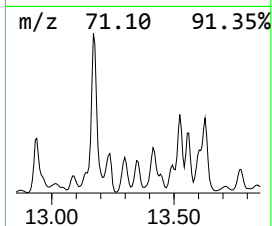
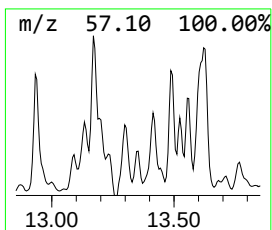
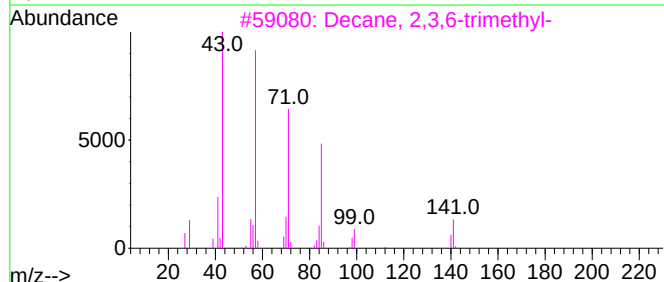
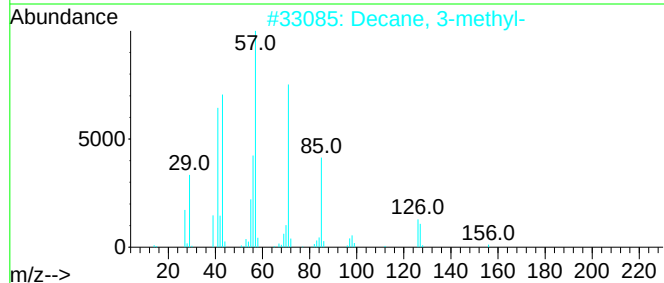
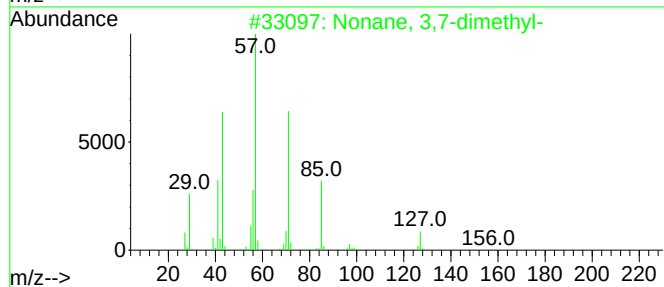
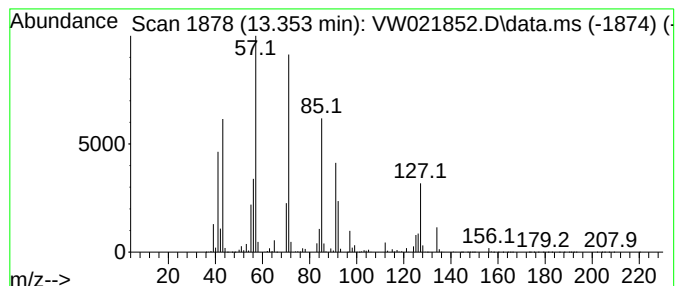
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MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

\*\*\*\*\*  
Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 46

| R.T.      | EstConc                    | Area    | Relative to ISTD       |         |             | R.T.   |
|-----------|----------------------------|---------|------------------------|---------|-------------|--------|
| 13.353    | 22.35 ug/L                 | 6360010 | 1,4-Dichlorobenzene-d4 |         |             | 13.560 |
| Hit# of 5 | Tentative ID               |         | MW                     | MolForm | CAS#        | Qual   |
| 1         | Nonane, 3,7-dimethyl-      |         | 156                    | C11H24  | 017302-32-8 | 93     |
| 2         | Decane, 3-methyl-          |         | 156                    | C11H24  | 013151-34-3 | 74     |
| 3         | Decane, 2,3,6-trimethyl-   |         | 184                    | C13H28  | 062238-12-4 | 53     |
| 4         | Octane, 2-methyl-          |         | 128                    | C9H20   | 003221-61-2 | 52     |
| 5         | Nonane, 2-methyl-5-propyl- |         | 184                    | C13H28  | 031081-17-1 | 50     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

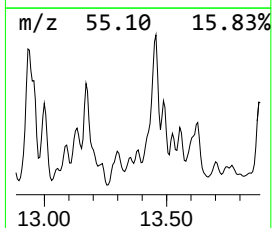
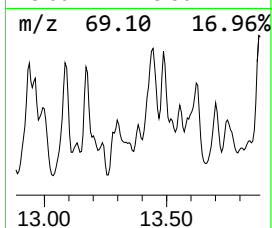
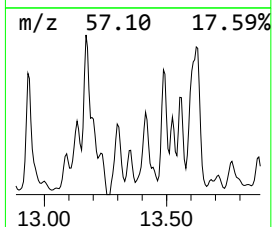
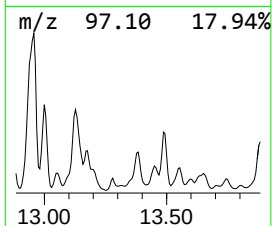
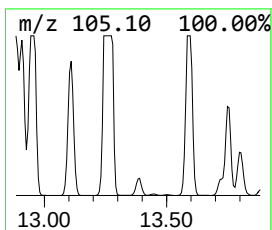
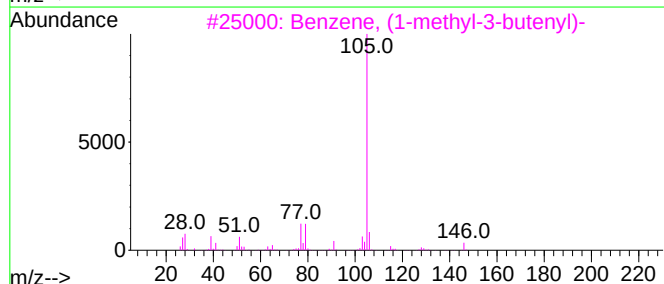
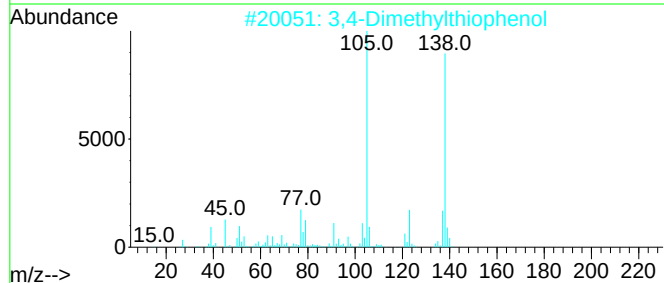
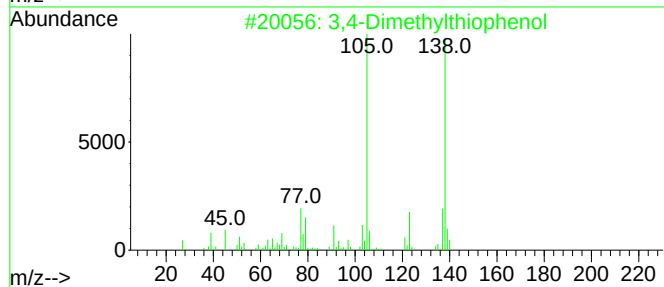
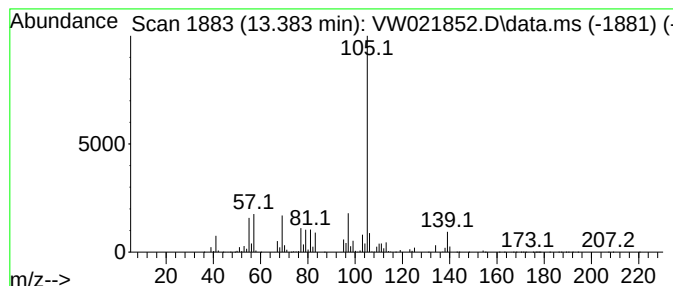
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Quant Title : SFAM01.0

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

\*\*\*\*\*  
Peak Number 40 unknown-02 Concentration Rank 65

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 13.383 | 5.52 ug/L | 1570110 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 3,4-Dimethylthiophenol              | 138 | C8H10S    | 018800-53-8  | 38   |
| 2    |    |   | 3,4-Dimethylthiophenol              | 138 | C8H10S    | 018800-53-8  | 38   |
| 3    |    |   | Benzene, (1-methyl-3-butenyl)-      | 146 | C11H14    | 010340-49-5  | 35   |
| 4    |    |   | 1-Phenylethanethiol, S-pentafluo... | 284 | C11H9F5OS | 1000469-38-3 | 35   |
| 5    |    |   | Disulfide, bis(p-methylbenzyl)      | 274 | C16H18S2  | 020193-94-6  | 35   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

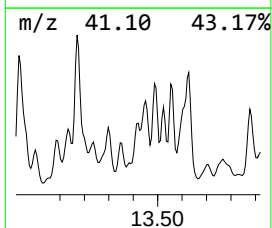
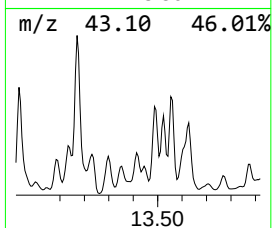
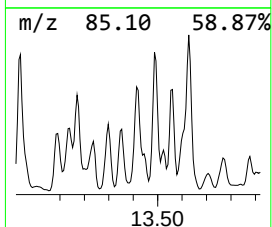
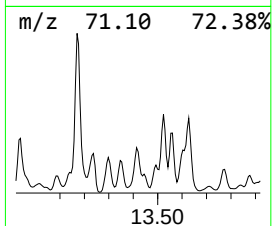
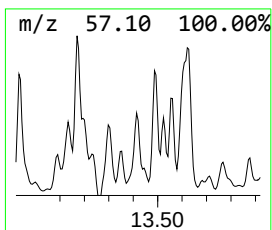
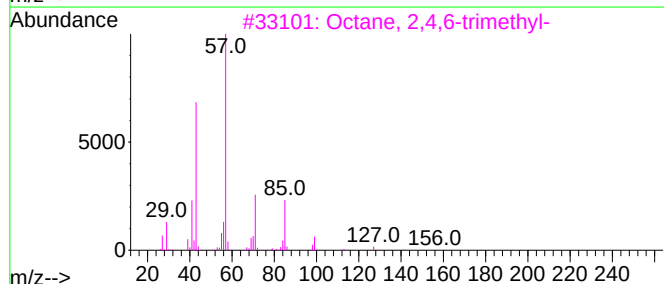
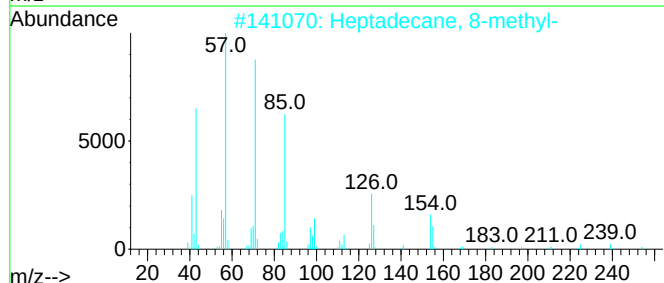
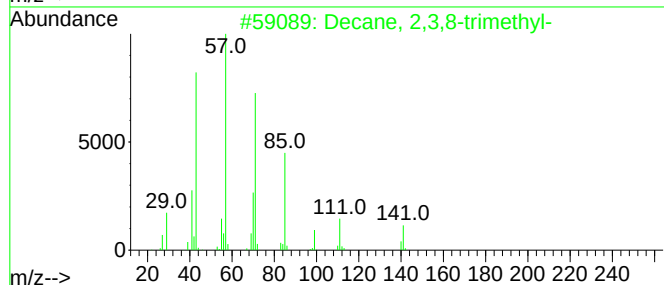
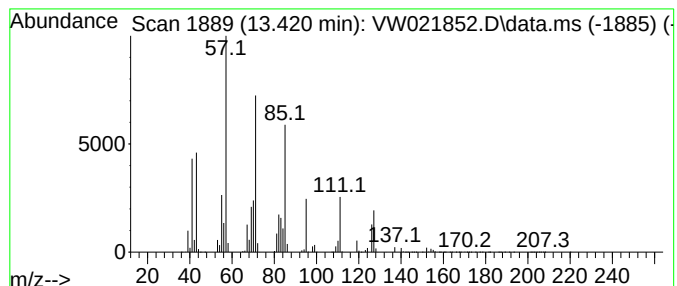
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TIC Integration Parameters: LSCINT.P

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Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 48

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 13.420 | 19.49 ug/L | 5547820 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Decane, 2,3,8-trimethyl-            | 184 | C13H28    | 062238-14-6  | 50   |
| 2    |    |   | Heptadecane, 8-methyl-              | 254 | C18H38    | 013287-23-5  | 50   |
| 3    |    |   | Octane, 2,4,6-trimethyl-            | 156 | C11H24    | 062016-37-9  | 46   |
| 4    |    |   | Sulfurous acid, isohexyl 2-penty... | 236 | C11H24O3S | 1000309-15-5 | 43   |
| 5    |    |   | Nonane, 5-butyl-                    | 184 | C13H28    | 017312-63-9  | 43   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

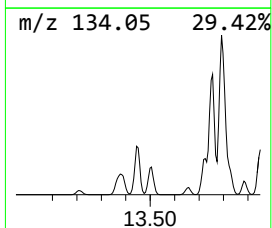
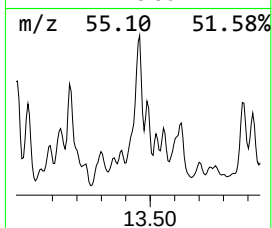
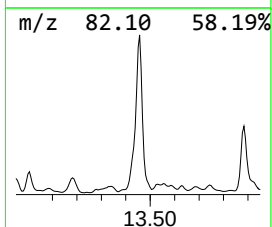
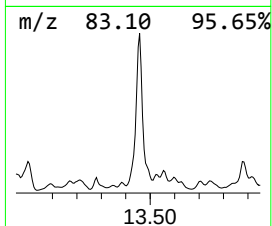
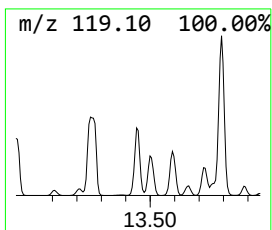
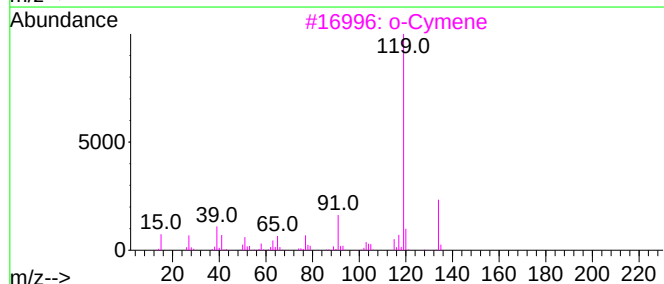
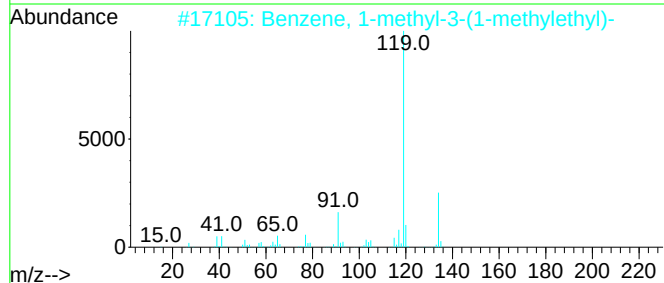
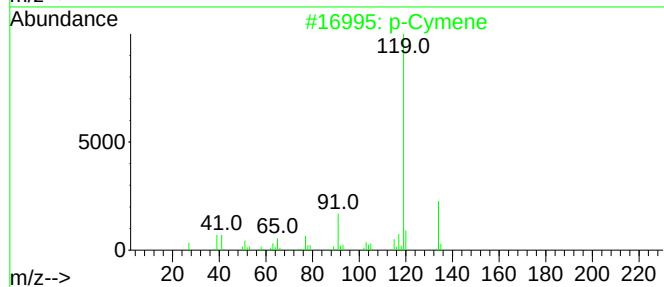
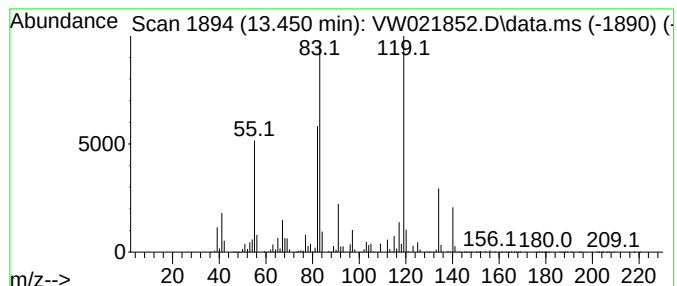
Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

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Peak Number 42 p-Cymene Concentration Rank 18

| R.T.      | EstConc                             | Area     | Relative to ISTD       |         |             | R.T.   |
|-----------|-------------------------------------|----------|------------------------|---------|-------------|--------|
| 13.451    | 53.56 ug/L                          | 15243300 | 1,4-Dichlorobenzene-d4 |         |             | 13.560 |
| Hit# of 5 | Tentative ID                        |          | MW                     | MolForm | CAS#        | Qual   |
| 1         | p-Cymene                            |          | 134                    | C10H14  | 000099-87-6 | 86     |
| 2         | Benzene, 1-methyl-3-(1-methyleth... |          | 134                    | C10H14  | 000535-77-3 | 86     |
| 3         | o-Cymene                            |          | 134                    | C10H14  | 000527-84-4 | 83     |
| 4         | Benzene, 1-methyl-3-(1-methyleth... |          | 134                    | C10H14  | 000535-77-3 | 83     |
| 5         | p-Cymene                            |          | 134                    | C10H14  | 000099-87-6 | 81     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

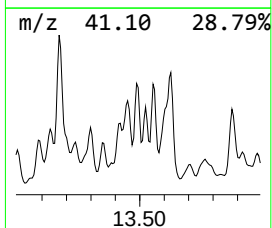
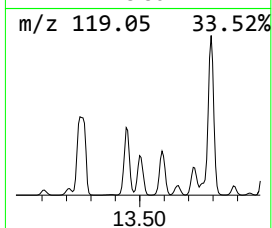
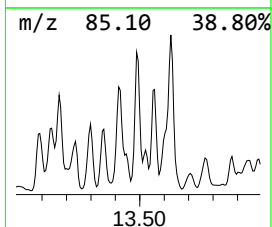
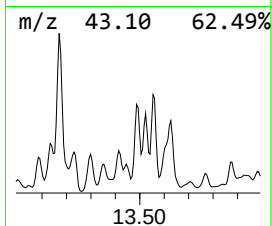
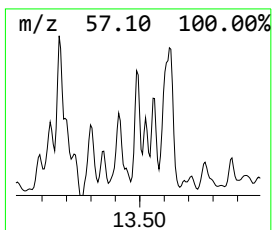
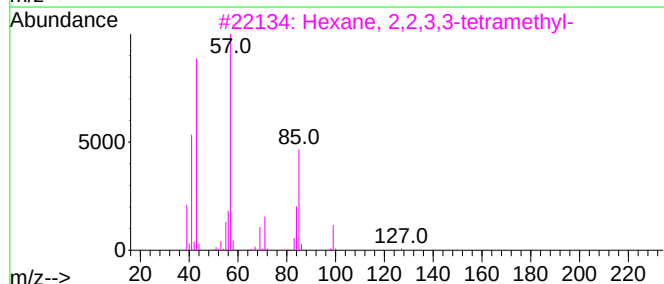
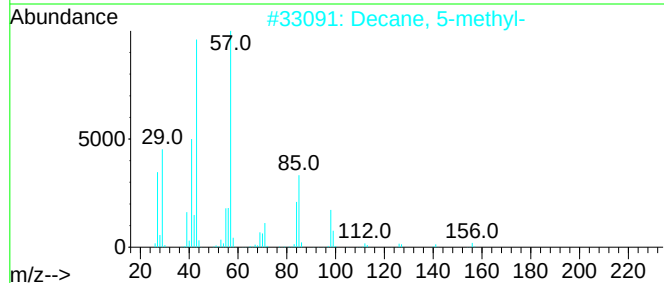
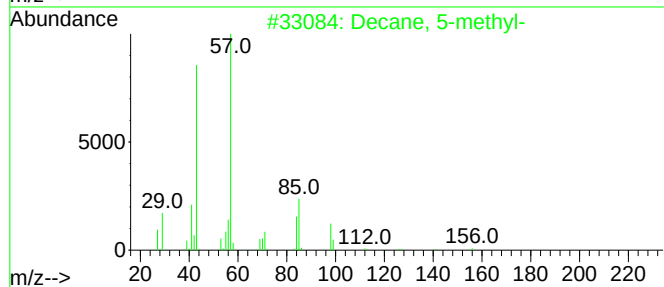
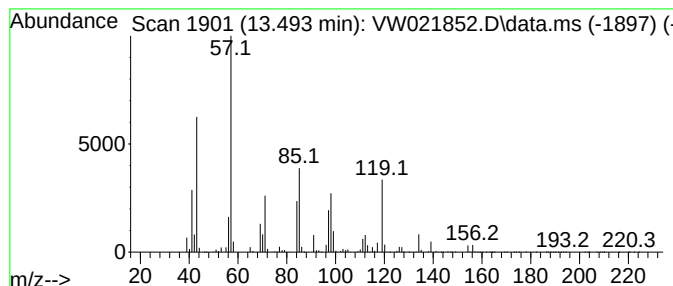
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Quant Title : SFAM01.0

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

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Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T.      | EstConc                      | Area     | Relative to ISTD       |         |             | R.T.   |
|-----------|------------------------------|----------|------------------------|---------|-------------|--------|
| 13.493    | 35.36 ug/L                   | 10064200 | 1,4-Dichlorobenzene-d4 |         |             | 13.560 |
| Hit# of 5 | Tentative ID                 |          | MW                     | MolForm | CAS#        | Qual   |
| 1         | Decane, 5-methyl-            |          | 156                    | C11H24  | 013151-35-4 | 70     |
| 2         | Decane, 5-methyl-            |          | 156                    | C11H24  | 013151-35-4 | 49     |
| 3         | Hexane, 2,2,3,3-tetramethyl- |          | 142                    | C10H22  | 013475-81-5 | 43     |
| 4         | Nonane                       |          | 128                    | C9H20   | 000111-84-2 | 43     |
| 5         | Heptane, 4-ethyl-            |          | 128                    | C9H20   | 002216-32-2 | 43     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

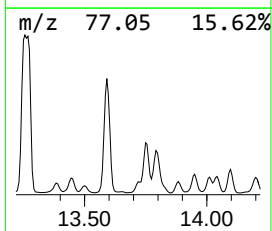
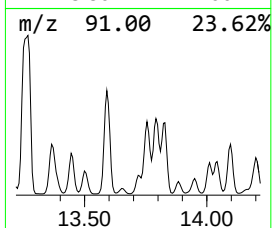
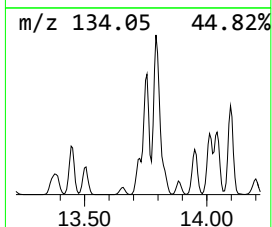
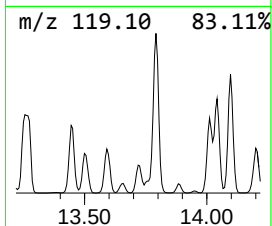
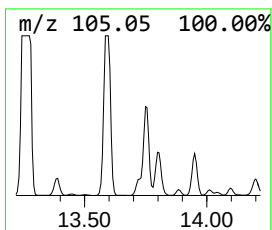
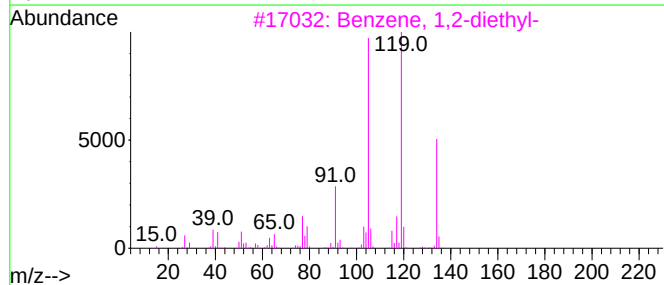
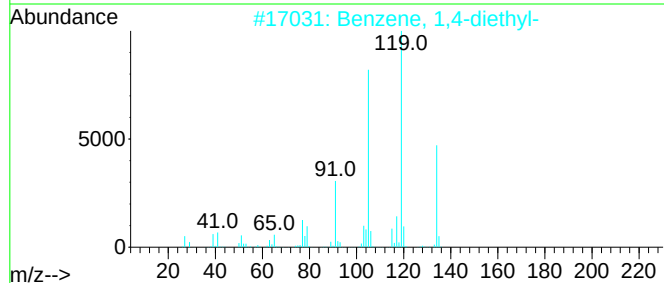
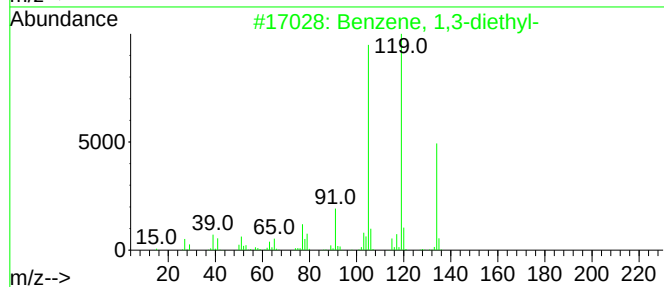
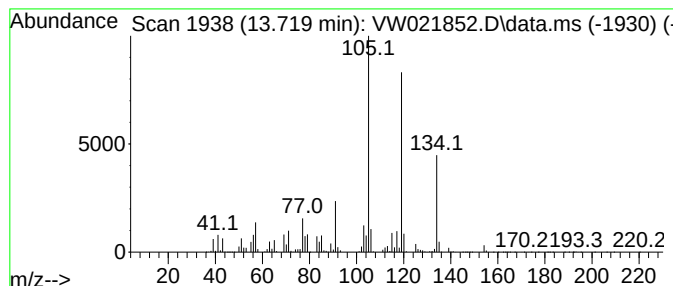
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Quant Title : SFAM01.0

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

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Peak Number 44 Benzene, 1,3-diethyl- Concentration Rank 39

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 13.719 | 28.91 ug/L | 8226670 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,3-diethyl- | 134 | C10H14  | 000141-93-5 | 95   |
| 2    |    |   | Benzene, 1,4-diethyl- | 134 | C10H14  | 000105-05-5 | 95   |
| 3    |    |   | Benzene, 1,2-diethyl- | 134 | C10H14  | 000135-01-3 | 94   |
| 4    |    |   | Benzene, 1,3-diethyl- | 134 | C10H14  | 000141-93-5 | 94   |
| 5    |    |   | Benzene, 1,4-diethyl- | 134 | C10H14  | 000105-05-5 | 90   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

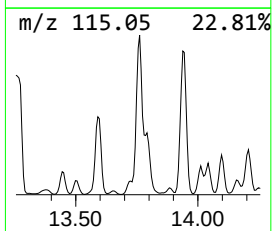
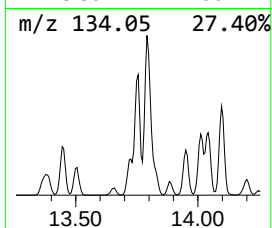
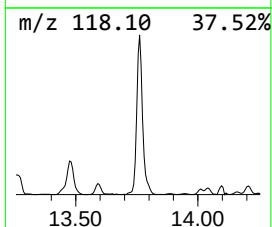
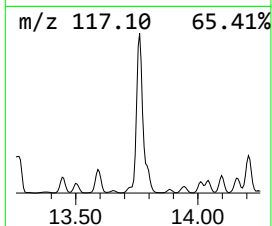
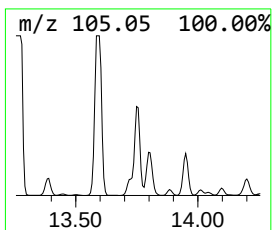
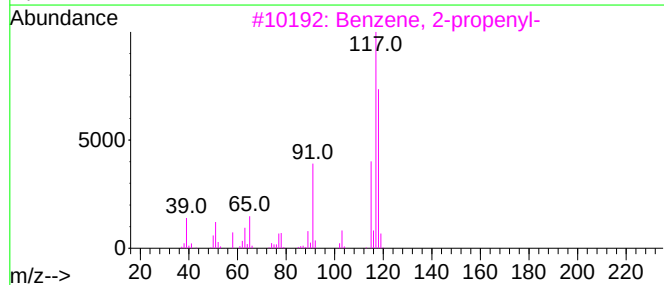
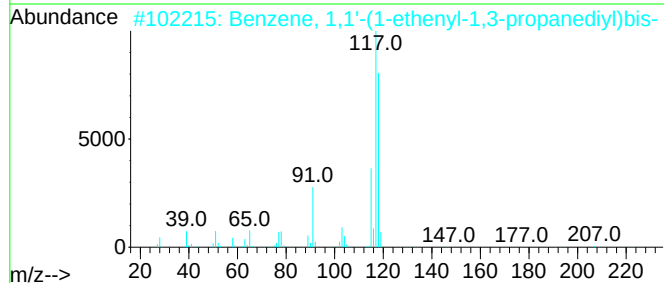
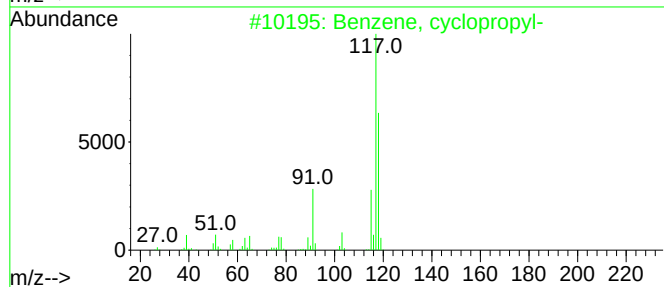
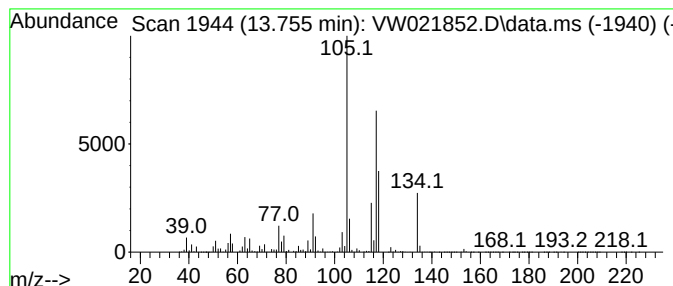
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Quant Title : SFAM01.0

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

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Peak Number 45 Benzene, cyclopropyl- Concentration Rank 9

| R.T.   | EstConc     | Area     | Relative to ISTD       | R.T.   |
|--------|-------------|----------|------------------------|--------|
| 13.755 | 114.54 ug/L | 32600600 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, cyclopropyl-               | 118 | C9H10   | 000873-49-4 | 55   |
| 2    |    |   | Benzene, 1,1'-(1-ethenyl-1,3-pro... | 222 | C17H18  | 061141-97-7 | 52   |
| 3    |    |   | Benzene, 2-propenyl-                | 118 | C9H10   | 000300-57-2 | 50   |
| 4    |    |   | Benzene, 1-ethenyl-2-methyl-        | 118 | C9H10   | 000611-15-4 | 49   |
| 5    |    |   | Benzene, 1-ethenyl-3-methyl-        | 118 | C9H10   | 000100-80-1 | 46   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

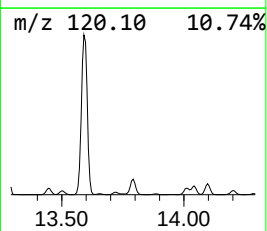
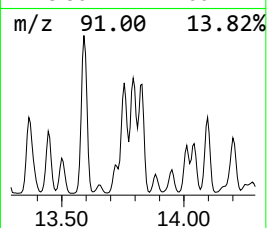
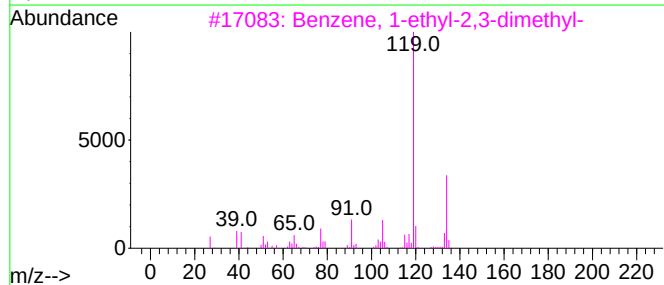
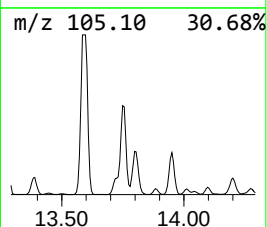
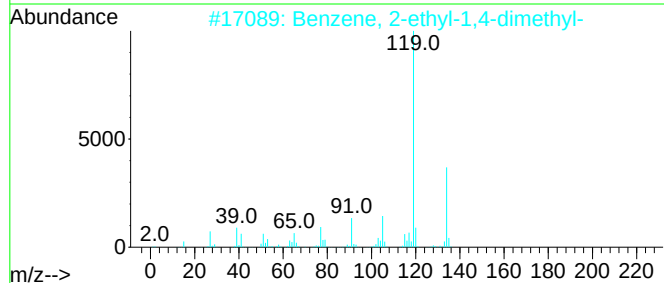
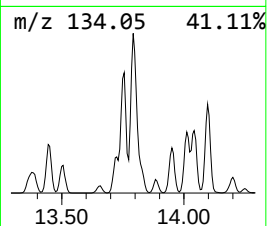
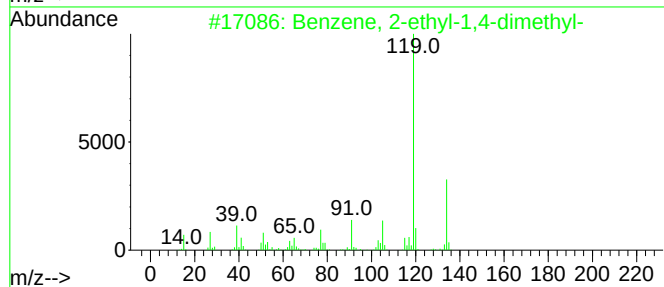
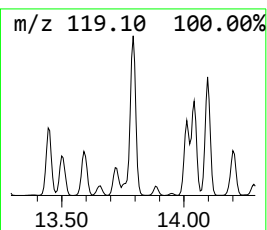
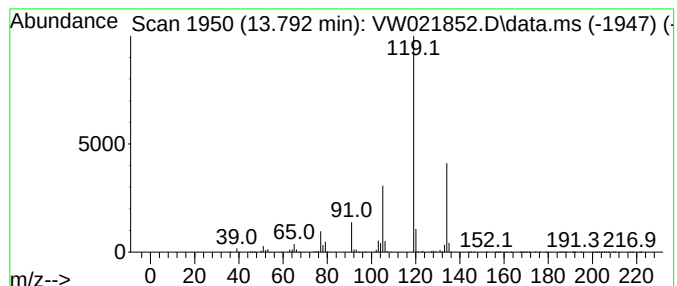
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 46 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 13

| R.T.   | EstConc    | Area     | Relative to ISTD       | R.T.   |
|--------|------------|----------|------------------------|--------|
| 13.792 | 94.65 ug/L | 26939300 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 91   |
| 2    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 91   |
| 3    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 91   |
| 4    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 87   |
| 5    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 87   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

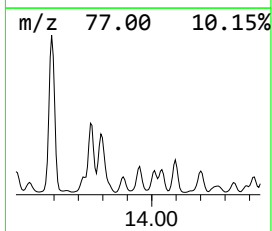
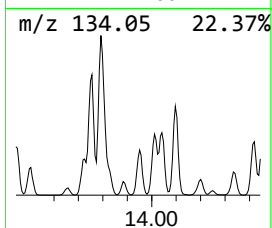
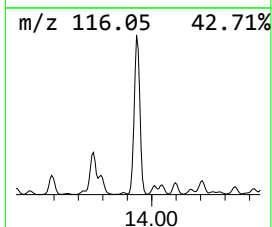
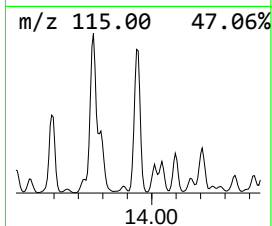
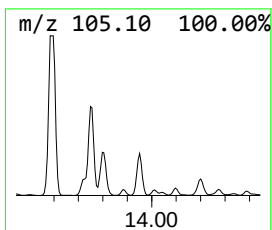
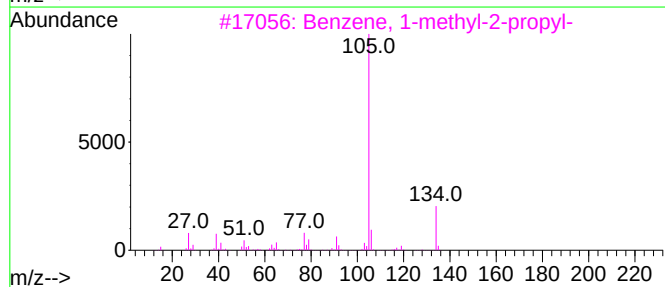
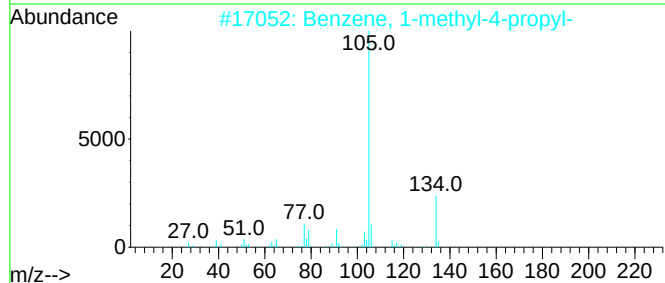
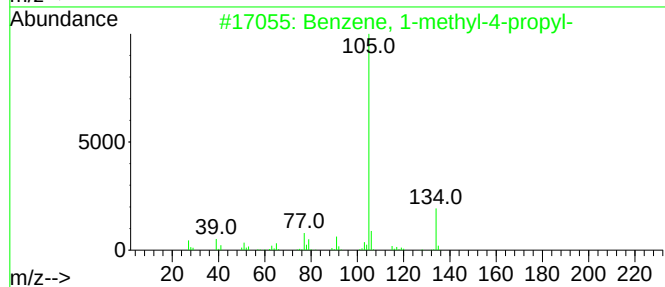
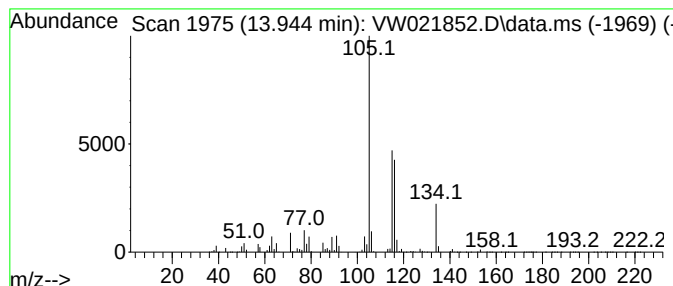
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 47 unknown-03 Concentration Rank 16

| R.T.   | EstConc    | Area     | Relative to ISTD       | R.T.   |
|--------|------------|----------|------------------------|--------|
| 13.944 | 55.80 ug/L | 15880900 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 42   |
| 2    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 42   |
| 3    |    |   | Benzene, 1-methyl-2-propyl- | 134 | C10H14  | 001074-17-5 | 42   |
| 4    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 42   |
| 5    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 38   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

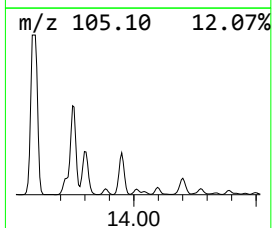
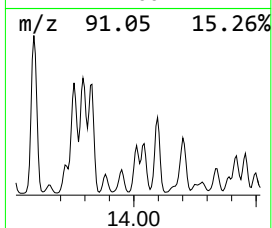
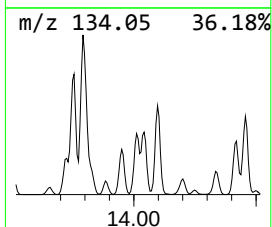
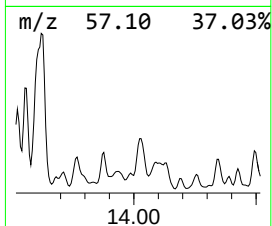
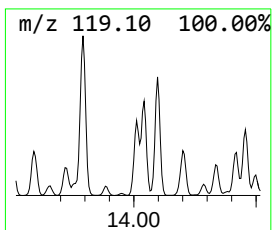
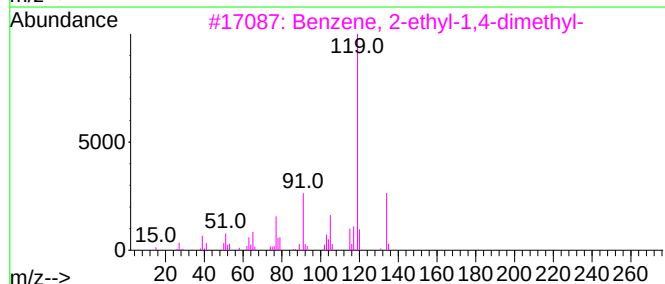
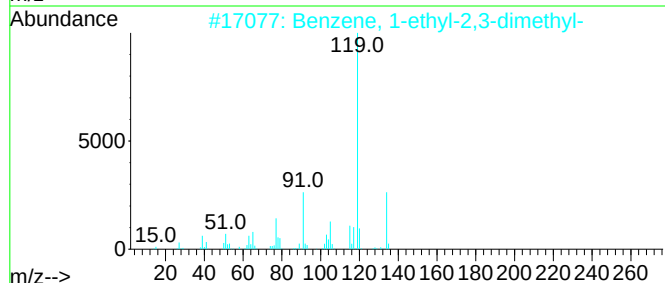
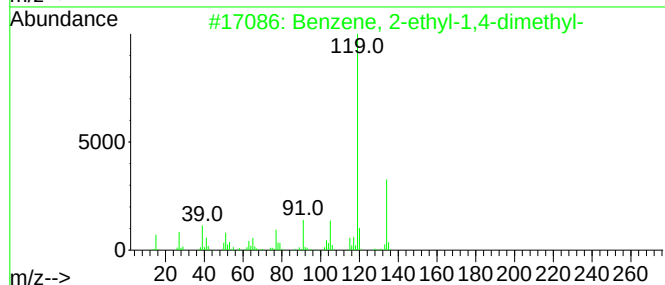
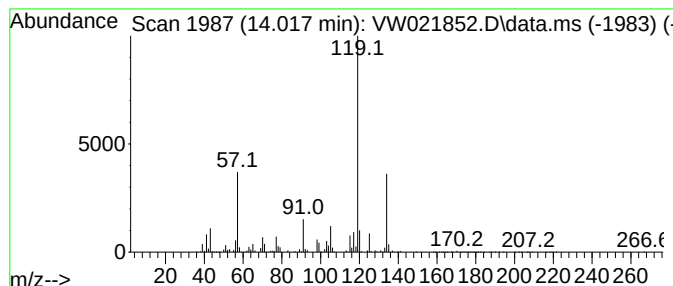
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Quant Title : SFAM01.0

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

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Peak Number 48 Benzene, 1-methyl-3-(1-meth... Concentration Rank 29

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 14.018 | 33.01 ug/L | 9394320 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 95   |
| 2    |    |   | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 95   |
| 3    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 95   |
| 4    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 94   |
| 5    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 94   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

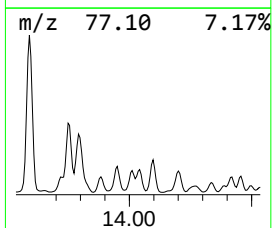
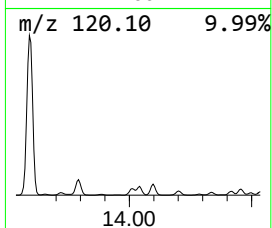
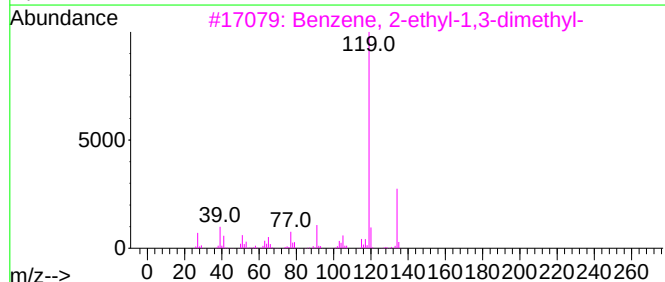
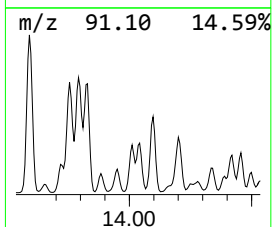
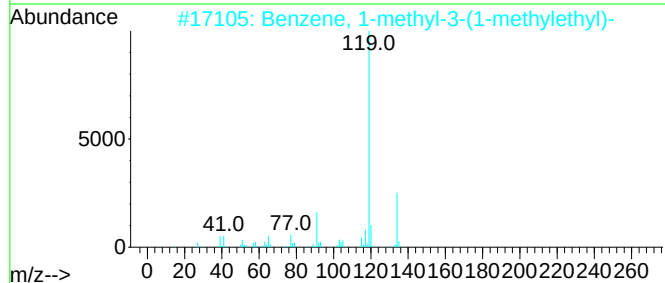
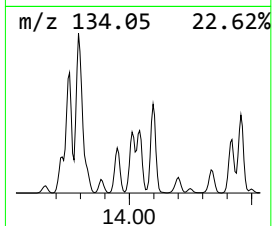
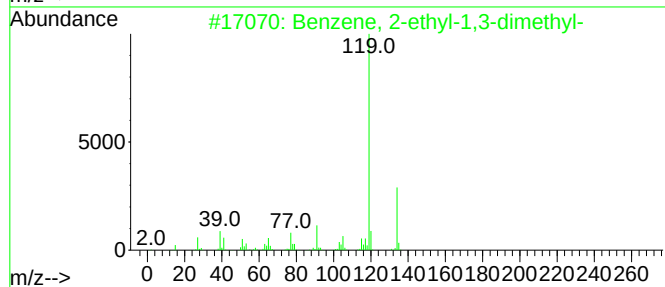
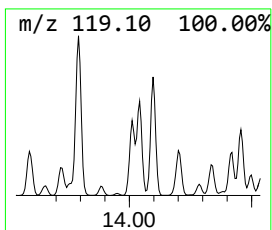
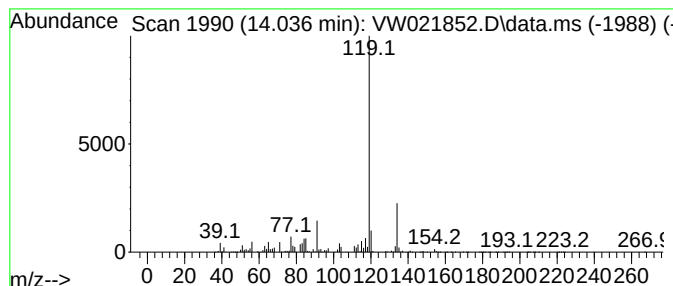
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 49 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 22

| R.T.   | EstConc    | Area     | Relative to ISTD       | R.T.   |
|--------|------------|----------|------------------------|--------|
| 14.036 | 43.32 ug/L | 12328000 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,3-dimethyl-       | 134 | C10H14  | 002870-04-4 | 87   |
| 2    |    |   | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 87   |
| 3    |    |   | Benzene, 2-ethyl-1,3-dimethyl-       | 134 | C10H14  | 002870-04-4 | 87   |
| 4    |    |   | Benzene, 4-ethyl-1,2-dimethyl-       | 134 | C10H14  | 000934-80-5 | 87   |
| 5    |    |   | Benzene, 1,2,3,5-tetramethyl-        | 134 | C10H14  | 000527-53-7 | 87   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

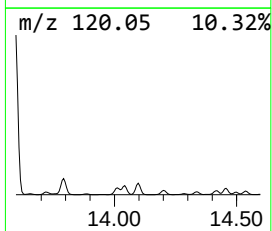
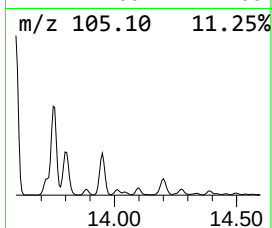
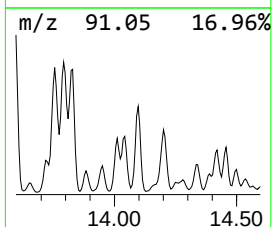
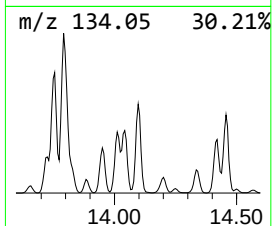
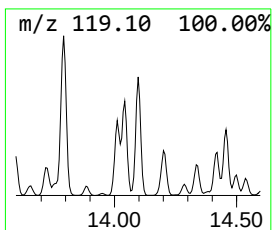
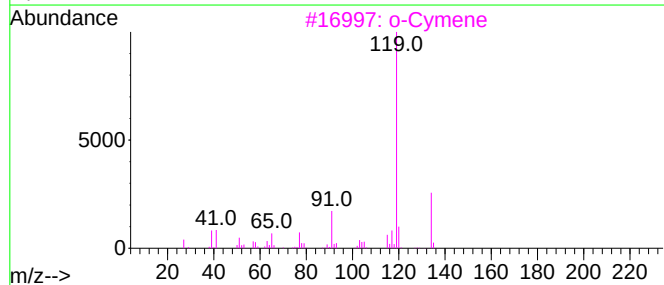
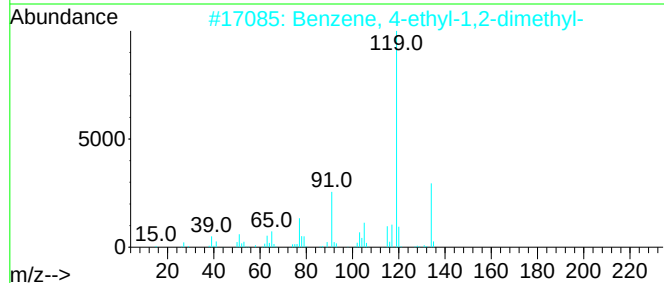
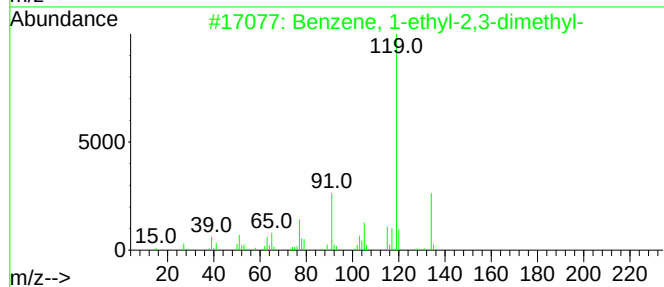
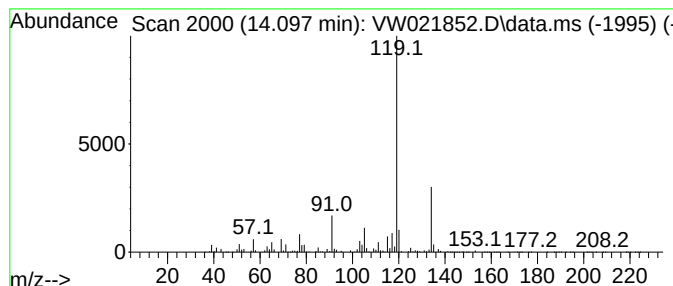
Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

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Peak Number 50 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 19

| R.T.      | EstConc                        | Area     | Relative to ISTD       |         |             | R.T.   |
|-----------|--------------------------------|----------|------------------------|---------|-------------|--------|
| 14.097    | 49.68 ug/L                     | 14138800 | 1,4-Dichlorobenzene-d4 |         |             | 13.560 |
| Hit# of 5 | Tentative ID                   |          | MW                     | MolForm | CAS#        | Qual   |
| 1         | Benzene, 1-ethyl-2,3-dimethyl- |          | 134                    | C10H14  | 000933-98-2 | 96     |
| 2         | Benzene, 4-ethyl-1,2-dimethyl- |          | 134                    | C10H14  | 000934-80-5 | 96     |
| 3         | o-Cymene                       |          | 134                    | C10H14  | 000527-84-4 | 95     |
| 4         | Benzene, 2-ethyl-1,4-dimethyl- |          | 134                    | C10H14  | 001758-88-9 | 94     |
| 5         | Benzene, 4-ethyl-1,2-dimethyl- |          | 134                    | C10H14  | 000934-80-5 | 94     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

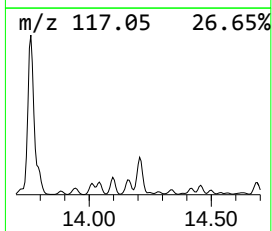
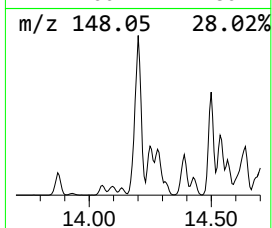
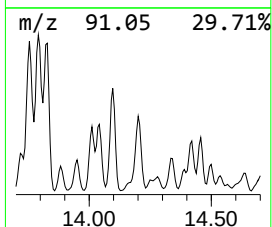
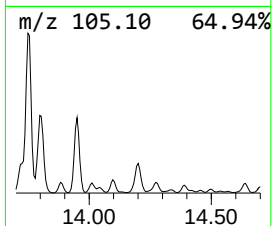
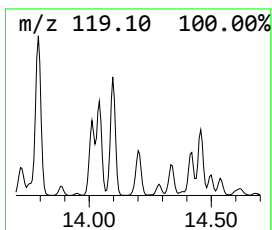
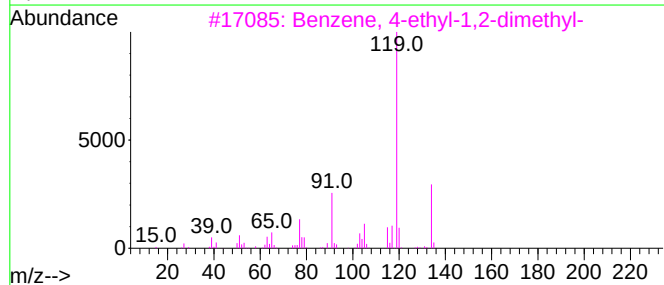
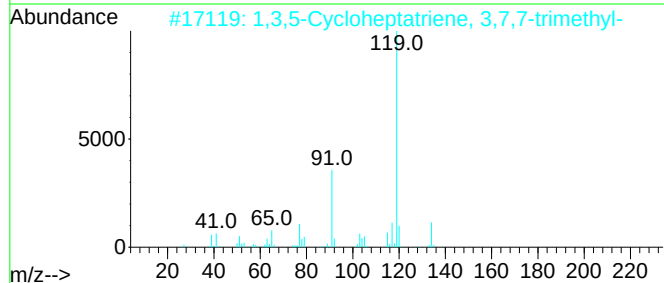
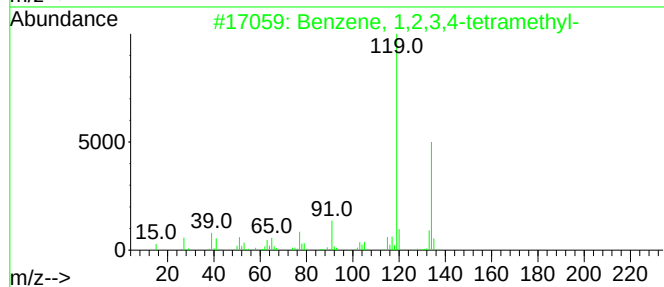
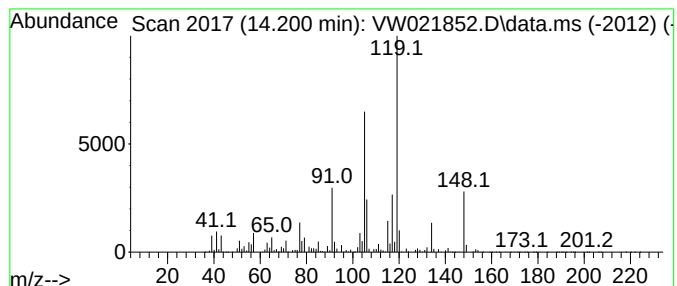
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 51 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 25

| R.T.   | EstConc    | Area     | Relative to ISTD       | R.T.   |
|--------|------------|----------|------------------------|--------|
| 14.200 | 36.34 ug/L | 10343300 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 64   |
| 2    |    |   | 1,3,5-Cycloheptatriene, 3,7,7-tr... | 134 | C10H14  | 003479-89-8 | 64   |
| 3    |    |   | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 64   |
| 4    |    |   | p-Cymene                            | 134 | C10H14  | 000099-87-6 | 60   |
| 5    |    |   | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 60   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

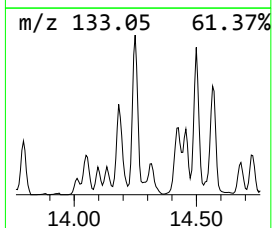
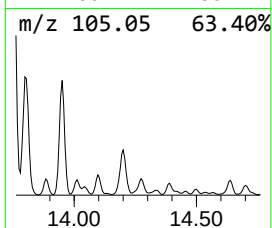
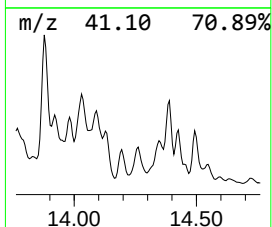
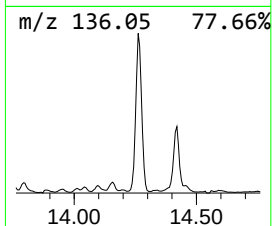
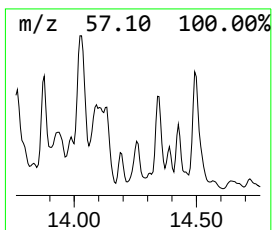
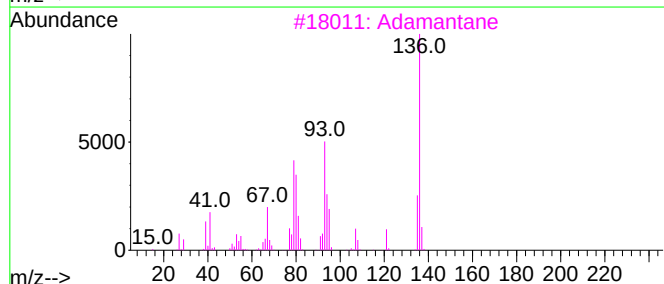
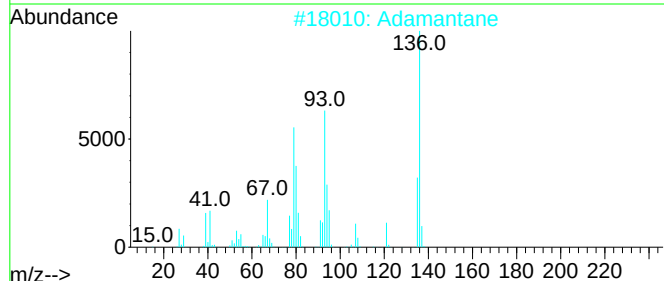
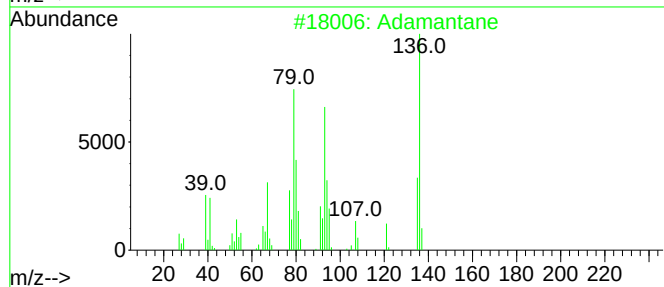
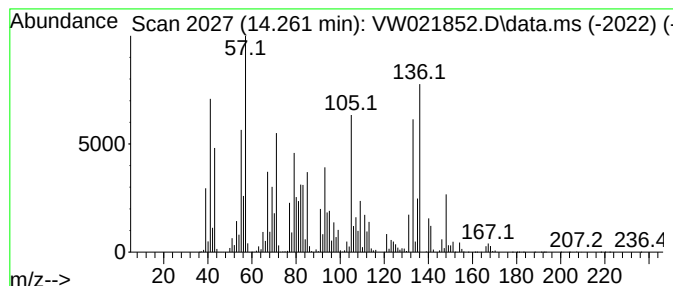
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Quant Title : SFAM01.0

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

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Peak Number 52 Adamantane Concentration Rank 45

| R.T.      | EstConc      | Area    | Relative to ISTD       |         | R.T.        |      |
|-----------|--------------|---------|------------------------|---------|-------------|------|
| 14.261    | 22.55 ug/L   | 6417840 | 1,4-Dichlorobenzene-d4 |         | 13.560      |      |
| Hit# of 5 | Tentative ID |         | MW                     | MolForm | CAS#        | Qual |
| 1         | Adamantane   |         | 136                    | C10H16  | 000281-23-2 | 92   |
| 2         | Adamantane   |         | 136                    | C10H16  | 000281-23-2 | 64   |
| 3         | Adamantane   |         | 136                    | C10H16  | 000281-23-2 | 60   |
| 4         | Adamantane   |         | 136                    | C10H16  | 000281-23-2 | 60   |
| 5         | Adamantane   |         | 136                    | C10H16  | 000281-23-2 | 49   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

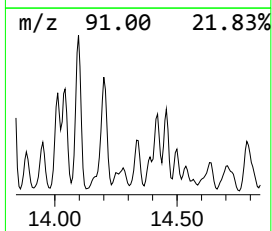
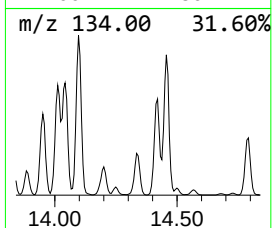
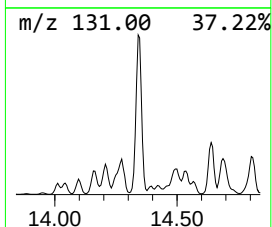
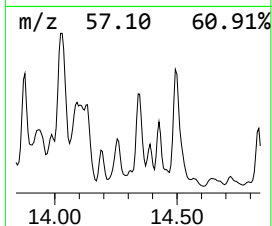
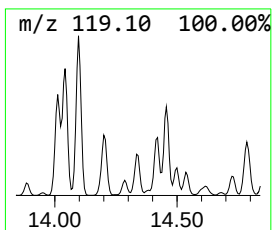
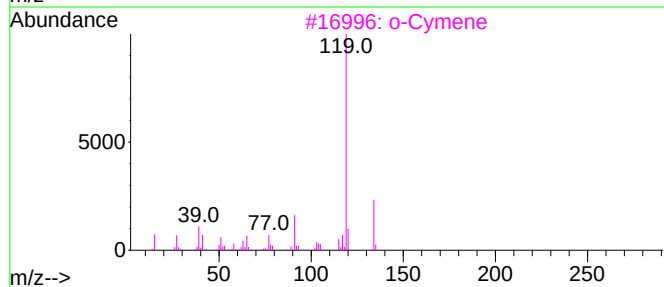
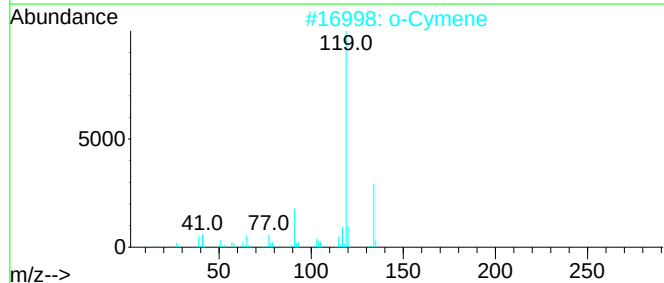
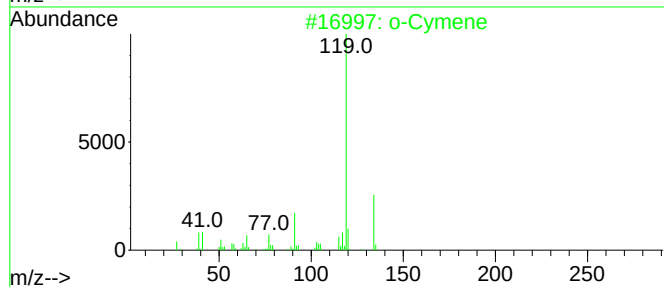
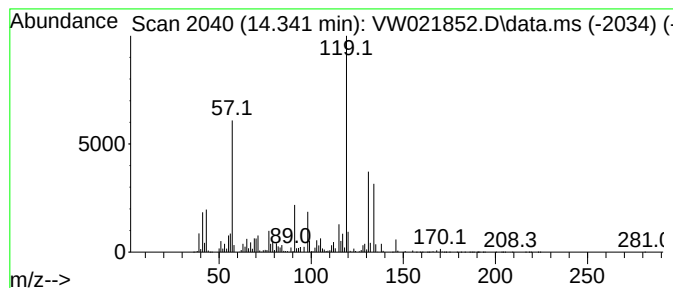
Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

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Peak Number 53 o-Cymene Concentration Rank 37

| R.T.      | EstConc                        | Area    | Relative to ISTD       |         |             | R.T.   |
|-----------|--------------------------------|---------|------------------------|---------|-------------|--------|
| 14.341    | 29.86 ug/L                     | 8498300 | 1,4-Dichlorobenzene-d4 |         |             | 13.560 |
| Hit# of 5 | Tentative ID                   |         | MW                     | MolForm | CAS#        | Qual   |
| 1         | o-Cymene                       |         | 134                    | C10H14  | 000527-84-4 | 92     |
| 2         | o-Cymene                       |         | 134                    | C10H14  | 000527-84-4 | 92     |
| 3         | o-Cymene                       |         | 134                    | C10H14  | 000527-84-4 | 92     |
| 4         | Benzene, 1-ethyl-2,4-dimethyl- |         | 134                    | C10H14  | 000874-41-9 | 86     |
| 5         | Benzene, 1,2,4,5-tetramethyl-  |         | 134                    | C10H14  | 000095-93-2 | 70     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

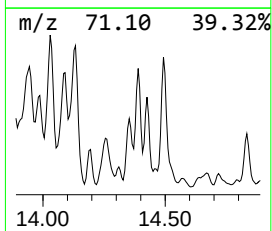
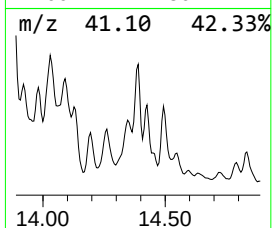
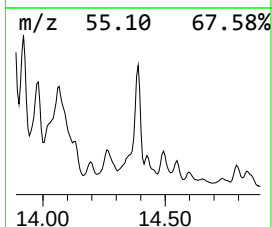
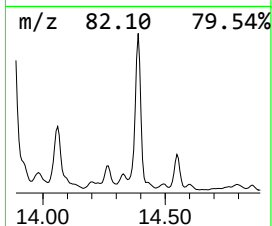
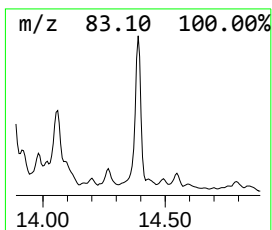
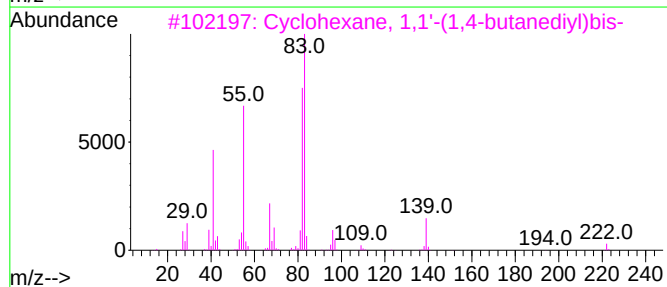
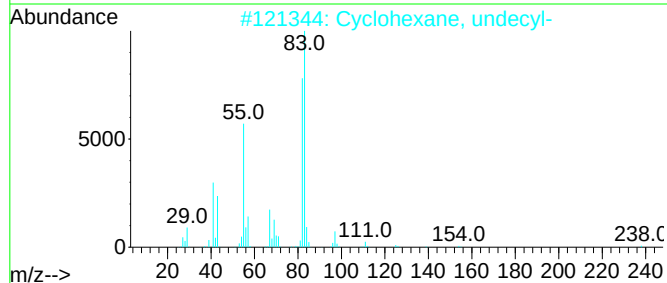
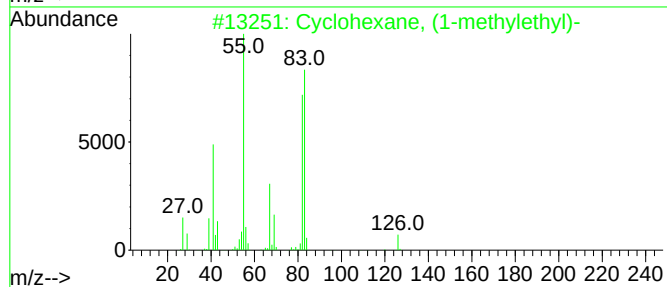
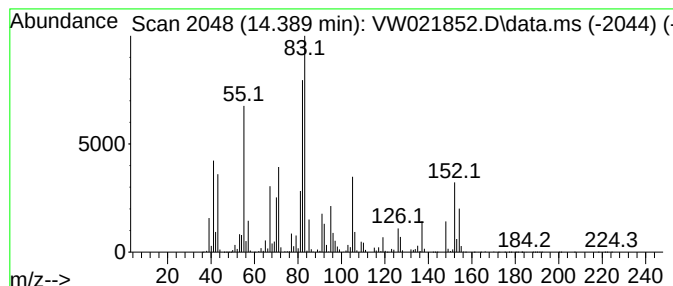
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Quant Title : SFAM01.0

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

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Peak Number 54 (DEL) Alkane: Cyclic14.389 Concentration Rank 26

| R.T.   | EstConc    | Area     | Relative to ISTD       | R.T.   |
|--------|------------|----------|------------------------|--------|
| 14.389 | 36.18 ug/L | 10297500 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, (1-methylethyl)-       | 126 | C9H18   | 000696-29-7 | 46   |
| 2    |    |   | Cyclohexane, undecyl-               | 238 | C17H34  | 054105-66-7 | 43   |
| 3    |    |   | Cyclohexane, 1,1'-(1,4-butanediyl)- | 222 | C16H30  | 006165-44-2 | 43   |
| 4    |    |   | Cyclohexane, (1-methylethyl)-       | 126 | C9H18   | 000696-29-7 | 43   |
| 5    |    |   | Cyclohexane, 1,1'-(1,4-butanediyl)- | 222 | C16H30  | 006165-44-2 | 43   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

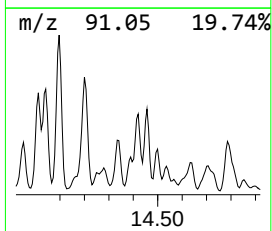
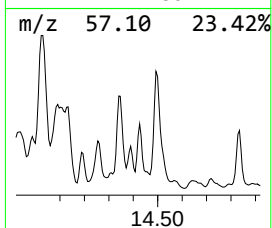
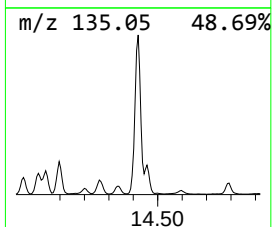
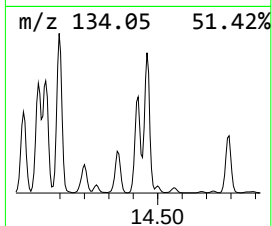
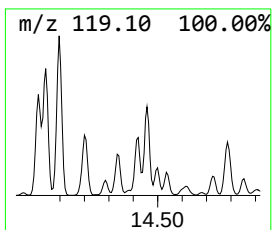
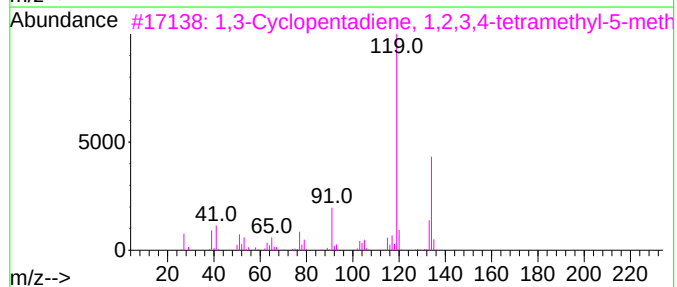
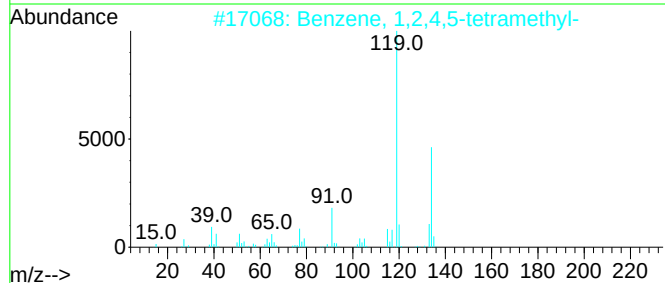
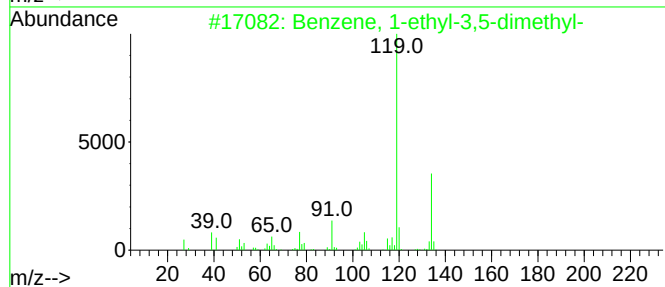
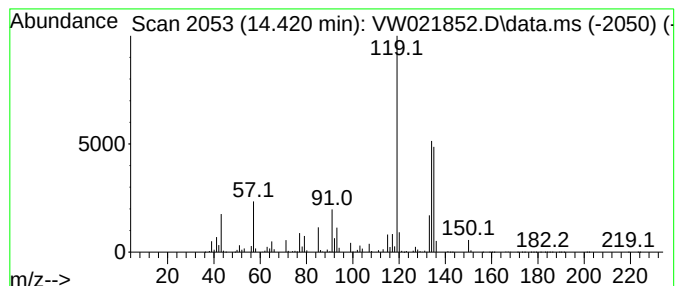
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Quant Title : SFAM01.0

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

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Peak Number 55 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 28

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 14.420 | 34.35 ug/L | 9775250 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 70   |
| 2    |    |   | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 70   |
| 3    |    |   | 1,3-Cyclopentadiene, 1,2,3,4-tet... | 134 | C10H14  | 076089-59-3 | 70   |
| 4    |    |   | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 70   |
| 5    |    |   | Benzene, 1,2,3,5-tetramethyl-       | 134 | C10H14  | 000527-53-7 | 70   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

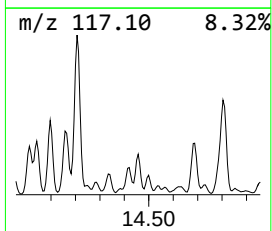
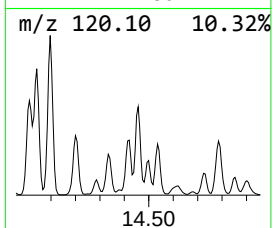
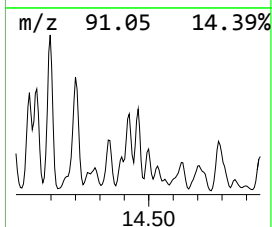
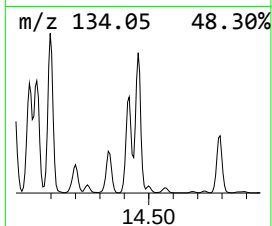
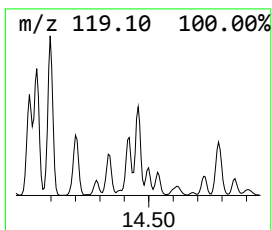
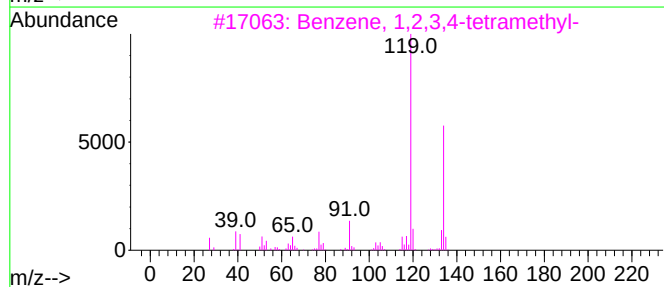
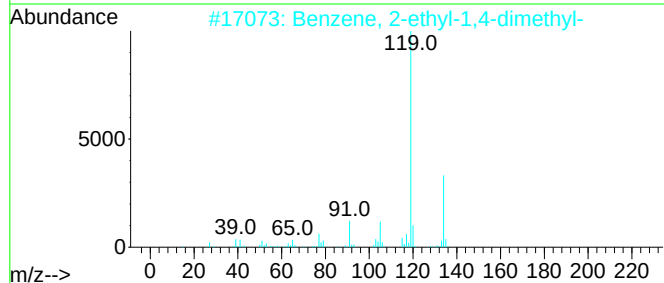
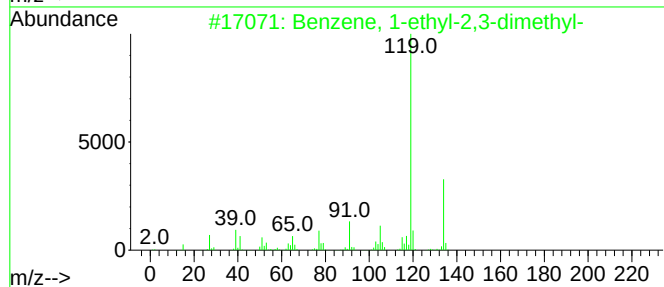
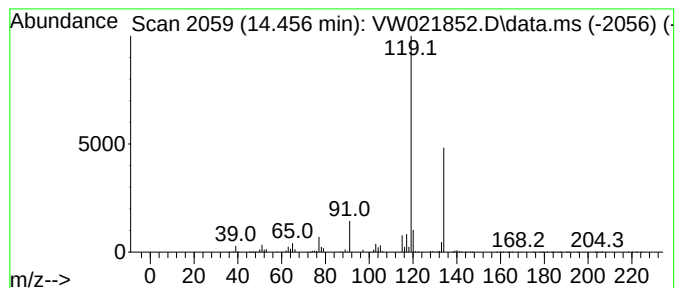
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Quant Title : SFAM01.0

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

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Peak Number 56 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 43

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 14.456 | 23.64 ug/L | 6729340 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 91   |
| 2    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 91   |
| 3    |    |   | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 91   |
| 4    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 91   |
| 5    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

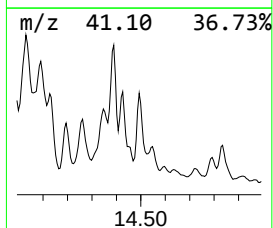
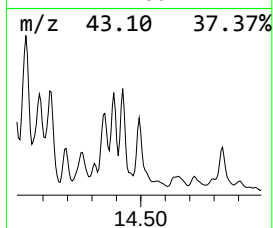
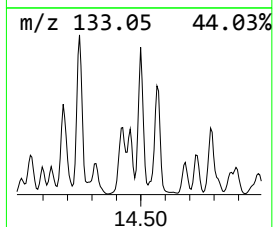
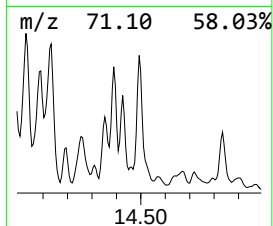
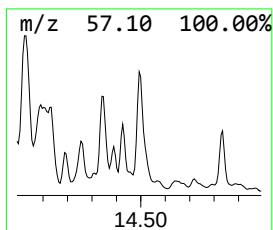
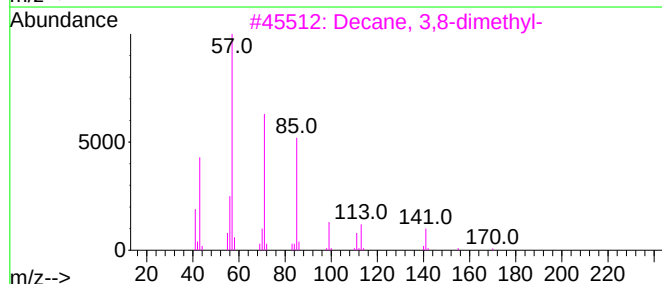
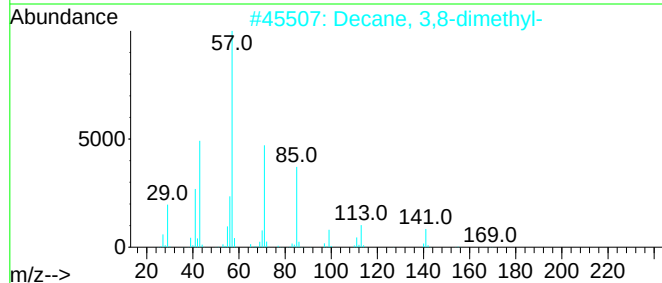
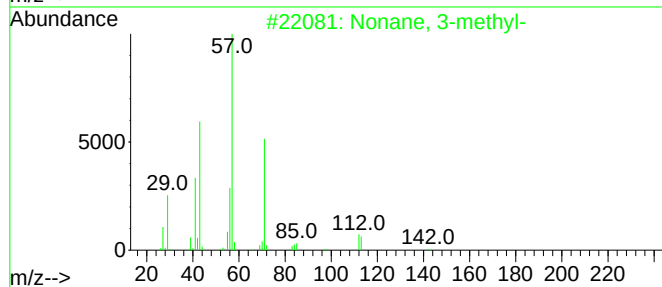
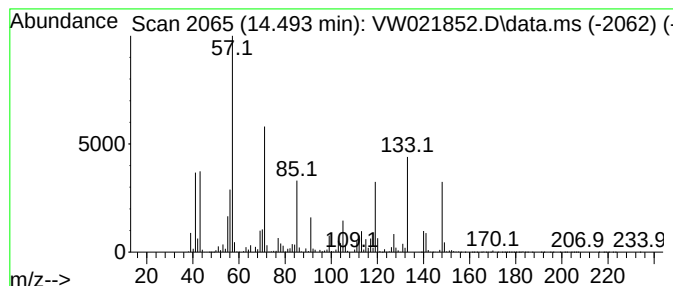
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Quant Title : SFAM01.0

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

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Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 41

| R.T.      | EstConc               | Area    | Relative to ISTD       | R.T.        |      |
|-----------|-----------------------|---------|------------------------|-------------|------|
| 14.493    | 25.07 ug/L            | 7135440 | 1,4-Dichlorobenzene-d4 | 13.560      |      |
| Hit# of 5 | Tentative ID          | MW      | MolForm                | CAS#        | Qual |
| 1         | Nonane, 3-methyl-     | 142     | C10H22                 | 005911-04-6 | 46   |
| 2         | Decane, 3,8-dimethyl- | 170     | C12H26                 | 017312-55-9 | 43   |
| 3         | Decane, 3,8-dimethyl- | 170     | C12H26                 | 017312-55-9 | 38   |
| 4         | Undecane, 3-methyl-   | 170     | C12H26                 | 001002-43-3 | 38   |
| 5         | Octane, 2,6-dimethyl- | 142     | C10H22                 | 002051-30-1 | 38   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

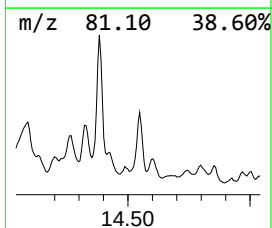
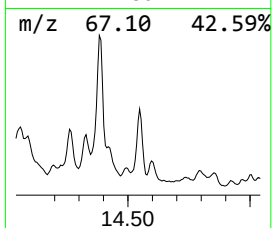
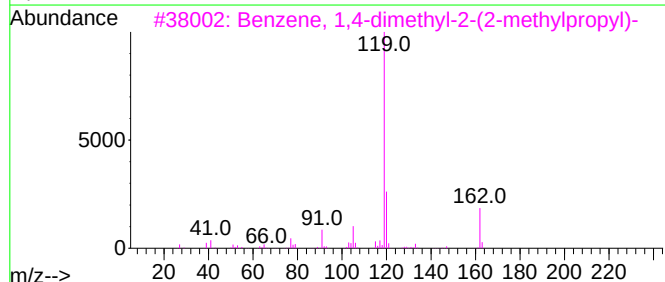
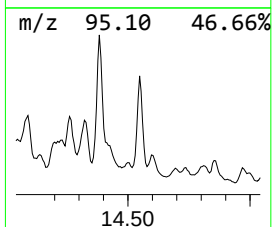
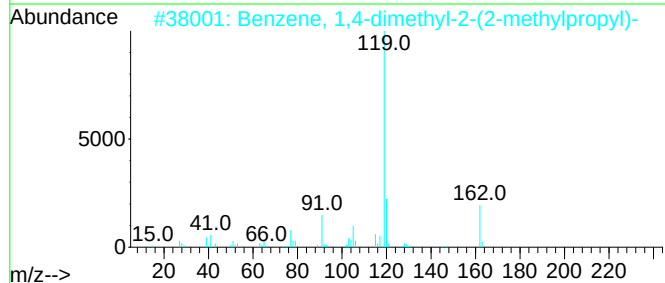
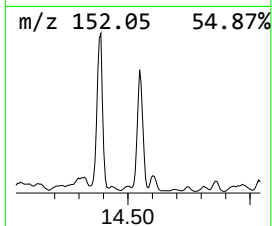
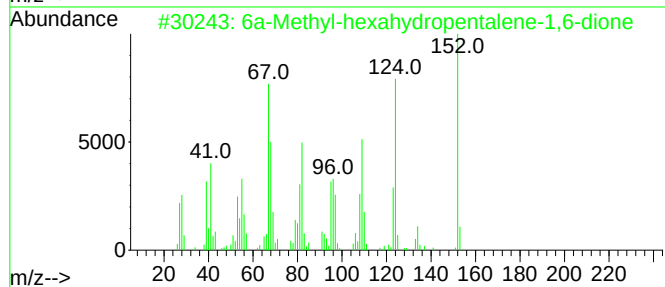
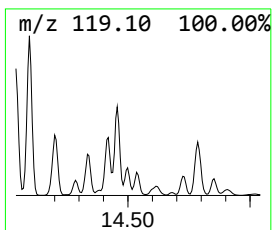
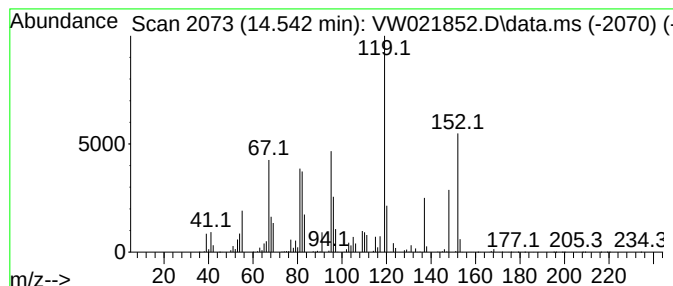
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 58 unknown-04 Concentration Rank 50

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 14.542 | 14.20 ug/L | 4041160 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|------|-------------------------------------|-----|---------|--------------|------|
| 1    |      | 6a-Methyl-hexahydropentalene-1,6... | 152 | C9H12O2 | 092485-86-4  | 27   |
| 2    |      | Benzene, 1,4-dimethyl-2-(2-methy... | 162 | C12H18  | 055669-88-0  | 20   |
| 3    |      | Benzene, 1,4-dimethyl-2-(2-methy... | 162 | C12H18  | 055669-88-0  | 14   |
| 4    |      | Benzene, 1,4-dimethyl-2-(2-methy... | 162 | C12H18  | 055669-88-0  | 14   |
| 5    |      | Methanimidamide, N,N-dimethyl-N'... | 152 | C7H12N4 | 1000327-41-1 | 14   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

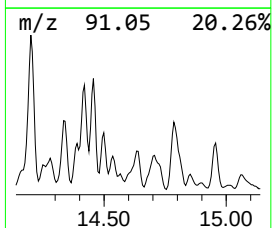
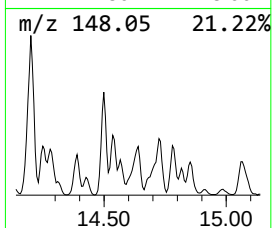
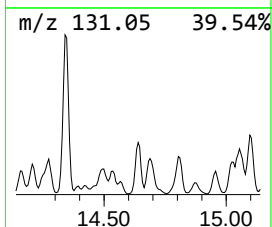
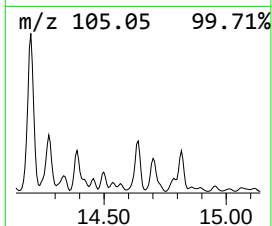
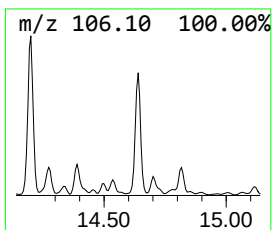
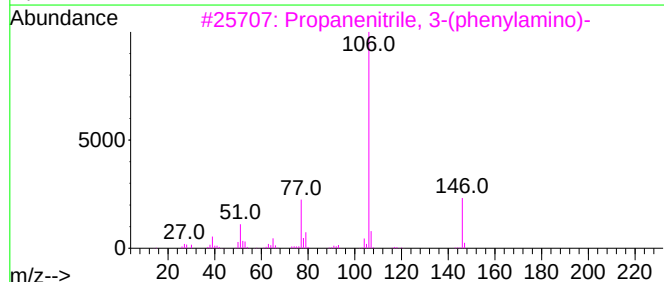
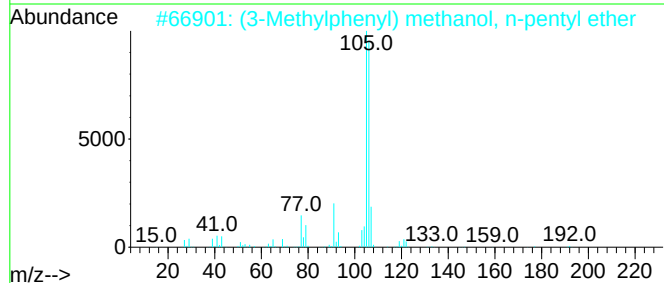
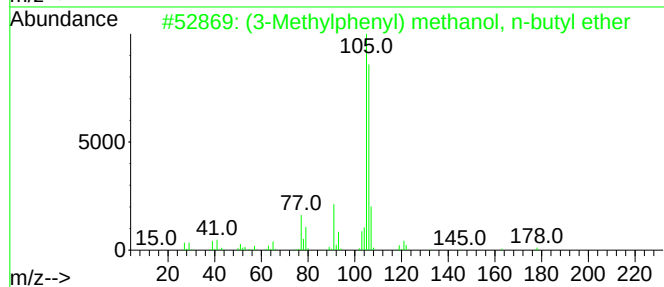
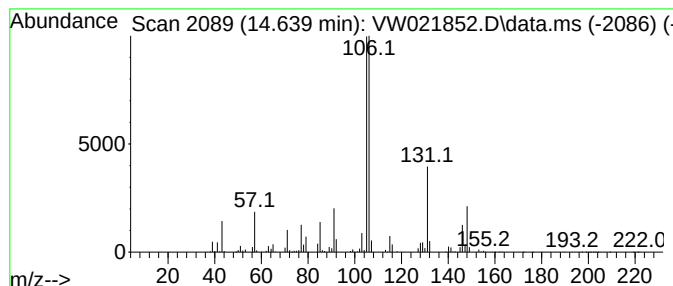
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 59 unknown-05 Concentration Rank 64

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 14.639 | 6.11 ug/L | 1740220 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | (3-Methylphenyl) methanol, n-but... | 178 | C12H18O | 1000374-44-2 | 38   |
| 2    |    |   | (3-Methylphenyl) methanol, n-pen... | 192 | C13H20O | 1000374-44-0 | 38   |
| 3    |    |   | Propanenitrile, 3-(phenylamino)-    | 146 | C9H10N2 | 001075-76-9  | 35   |
| 4    |    |   | (3-Methylphenyl) methanol, 3-met... | 192 | C13H20O | 1000374-58-6 | 32   |
| 5    |    |   | (3-Methylphenyl) methanol, neope... | 192 | C13H20O | 1000374-44-1 | 32   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

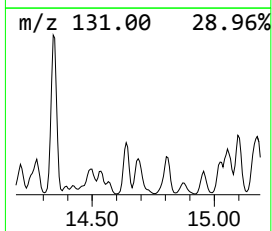
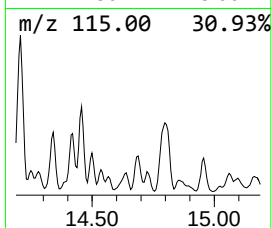
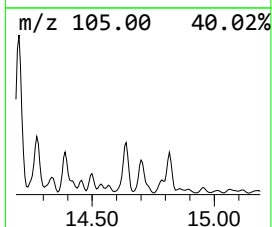
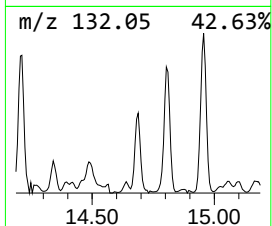
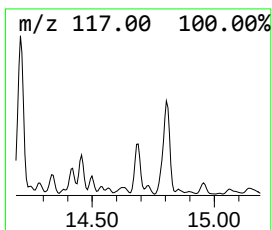
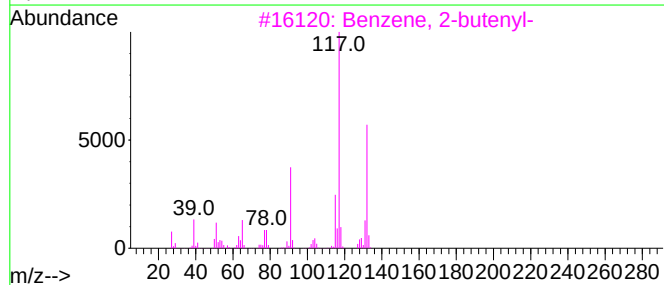
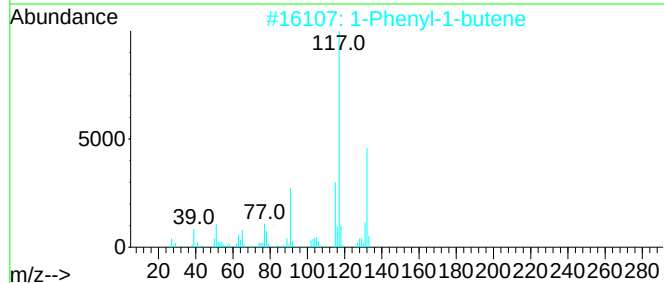
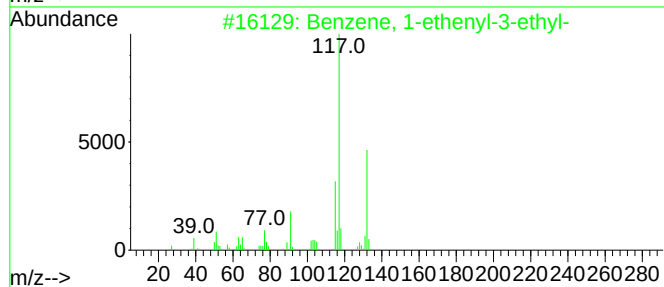
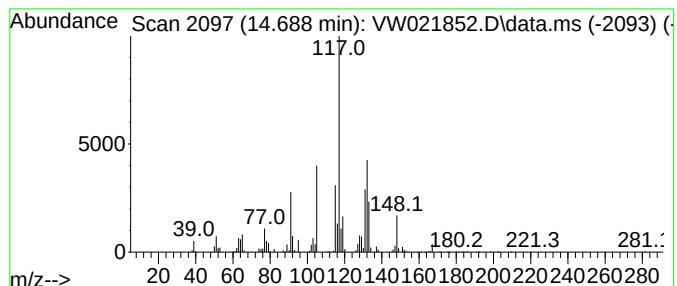
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Quant Title : SFAM01.0

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

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Peak Number 60 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 68

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 14.688 | 4.82 ug/L | 1373180 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethenyl-3-ethyl-         | 132 | C10H12  | 007525-62-4 | 76   |
| 2    |    |   | 1-Phenyl-1-butene                   | 132 | C10H12  | 000824-90-8 | 76   |
| 3    |    |   | Benzene, 2-butenyl-                 | 132 | C10H12  | 001560-06-1 | 76   |
| 4    |    |   | Benzene, 2-ethenyl-1,3-dimethyl-    | 132 | C10H12  | 002039-90-9 | 76   |
| 5    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000767-99-7 | 76   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

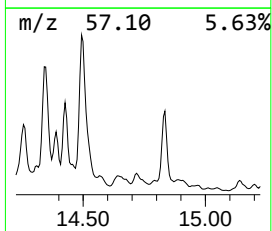
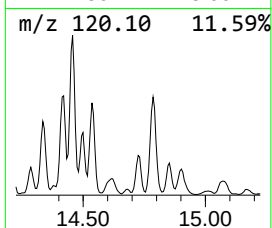
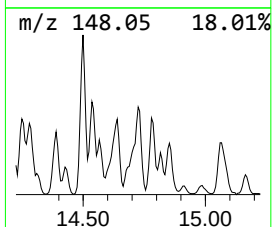
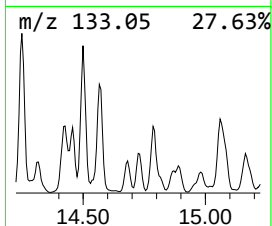
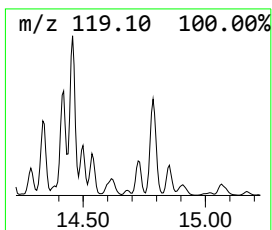
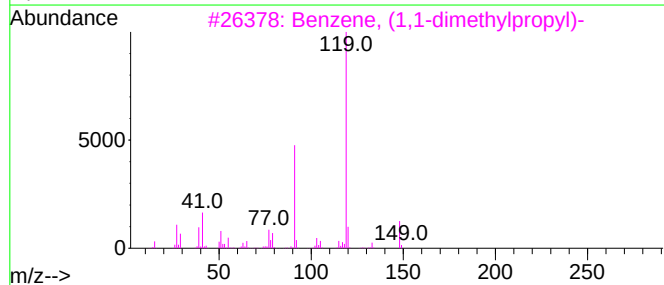
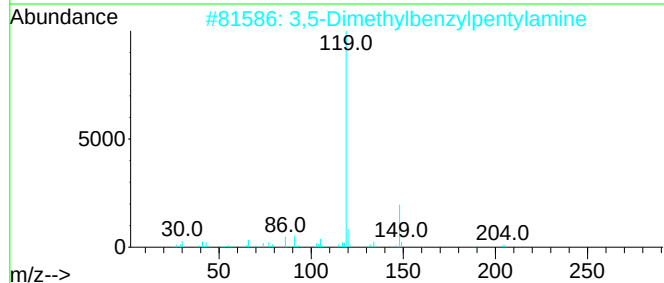
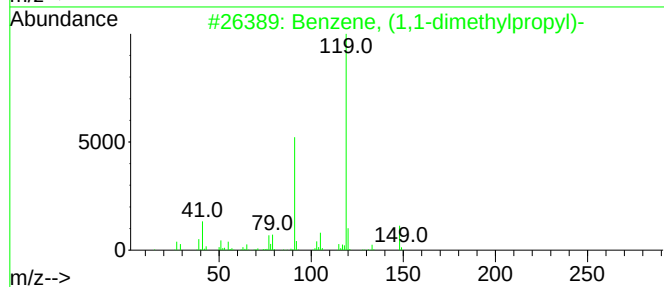
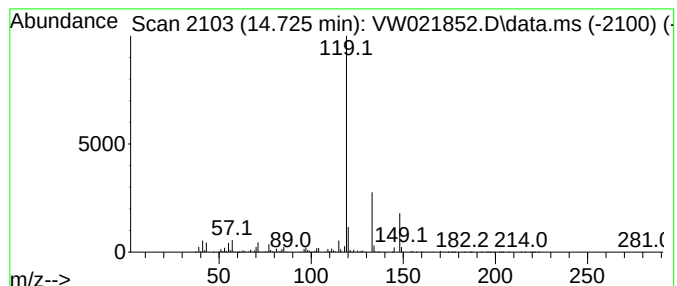
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Quant Title : SFAM01.0

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

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Peak Number 61 Benzene, (1,1-dimethylpropyl)- Concentration Rank 60

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 14.725 | 7.03 ug/L | 2000090 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8  | 64   |
| 2    |    |   | 3,5-Dimethylbenzylpentylamine       | 205 | C14H23N | 1000491-60-8 | 64   |
| 3    |    |   | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8  | 64   |
| 4    |    |   | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16  | 001595-16-0  | 64   |
| 5    |    |   | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8  | 59   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

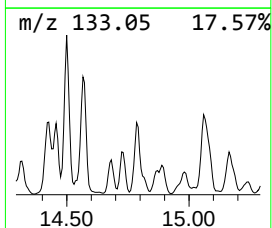
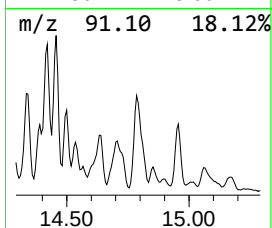
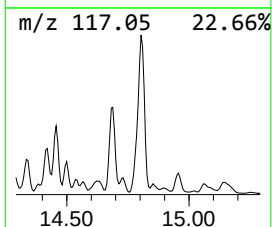
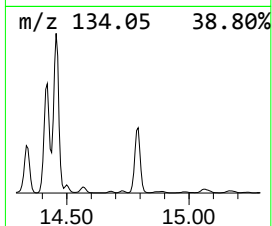
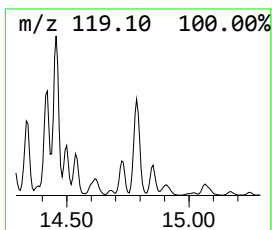
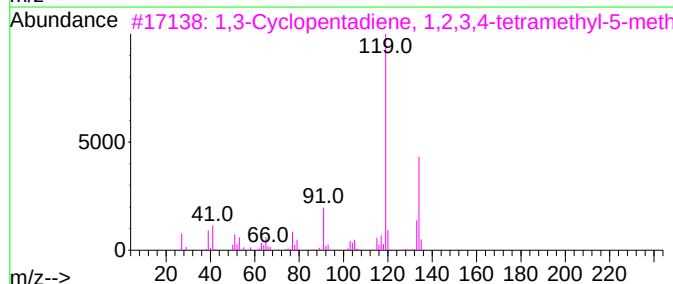
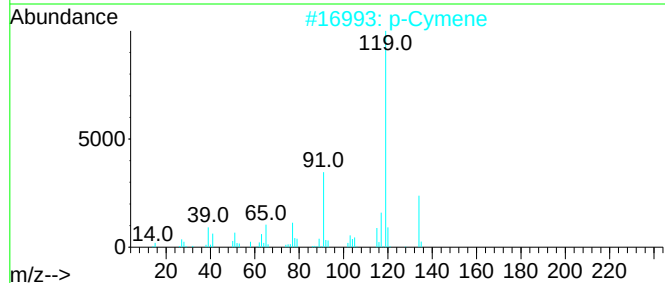
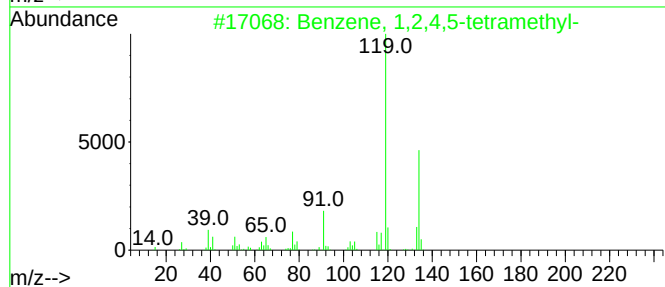
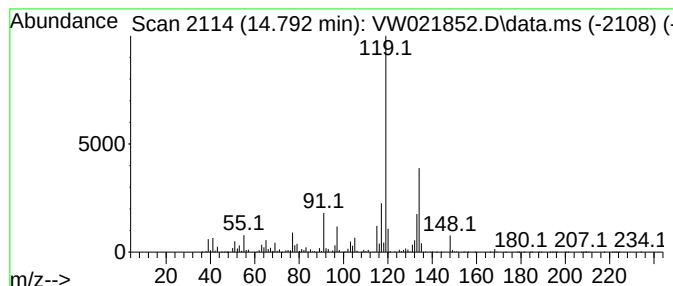
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Quant Title : SFAM01.0

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

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Peak Number 62 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 38

| R.T.      | EstConc                             | Area    | Relative to ISTD       |         |             | R.T.   |
|-----------|-------------------------------------|---------|------------------------|---------|-------------|--------|
| 14.792    | 29.03 ug/L                          | 8262630 | 1,4-Dichlorobenzene-d4 |         |             | 13.560 |
| Hit# of 5 | Tentative ID                        |         | MW                     | MolForm | CAS#        | Qual   |
| 1         | Benzene, 1,2,4,5-tetramethyl-       |         | 134                    | C10H14  | 000095-93-2 | 94     |
| 2         | p-Cymene                            |         | 134                    | C10H14  | 000099-87-6 | 81     |
| 3         | 1,3-Cyclopentadiene, 1,2,3,4-tet... |         | 134                    | C10H14  | 076089-59-3 | 81     |
| 4         | Benzene, 1,2,3,4-tetramethyl-       |         | 134                    | C10H14  | 000488-23-3 | 76     |
| 5         | o-Cymene                            |         | 134                    | C10H14  | 000527-84-4 | 76     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

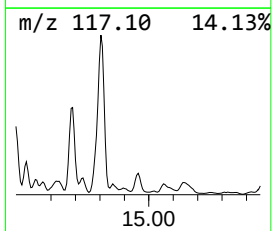
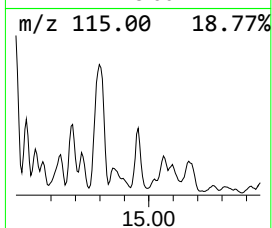
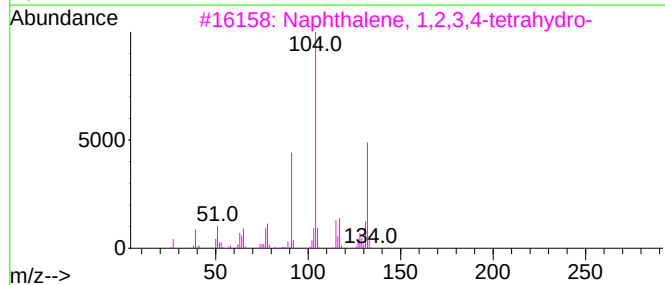
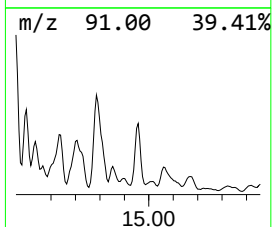
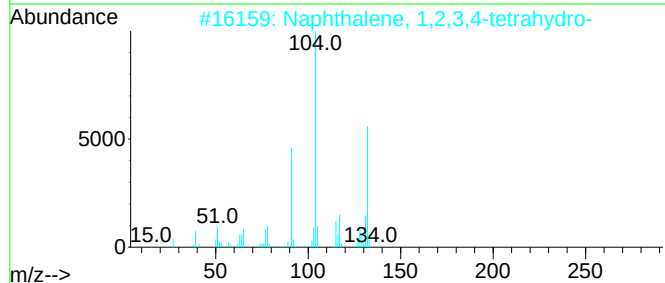
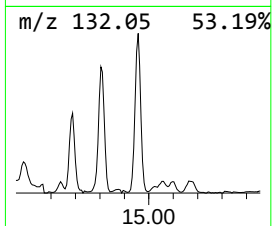
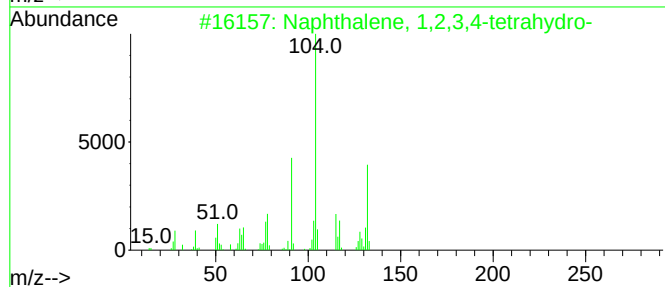
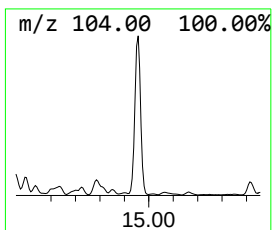
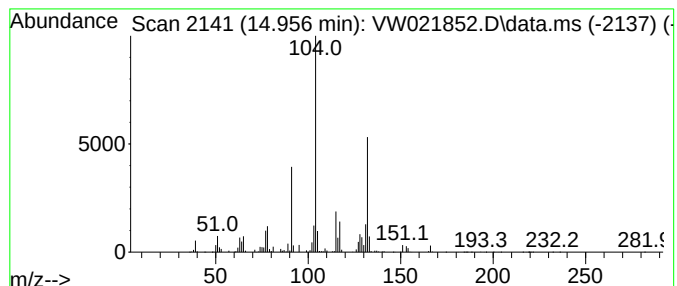
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 63 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 61

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 14.956 | 6.72 ug/L | 1911190 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 96   |
| 2    |    |   | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 96   |
| 3    |    |   | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 96   |
| 4    |    |   | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 94   |
| 5    |    |   | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 93   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

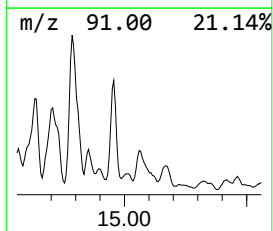
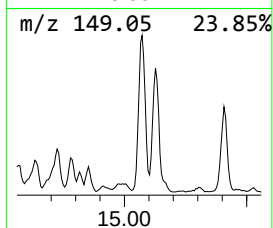
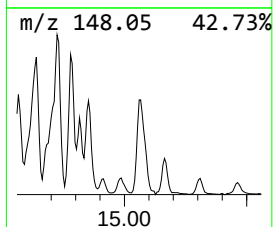
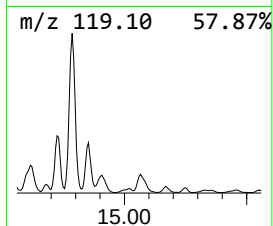
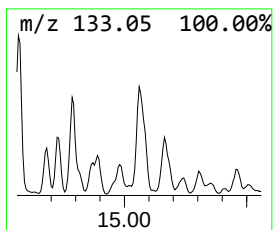
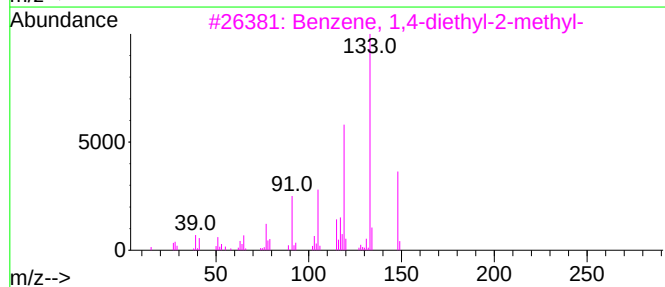
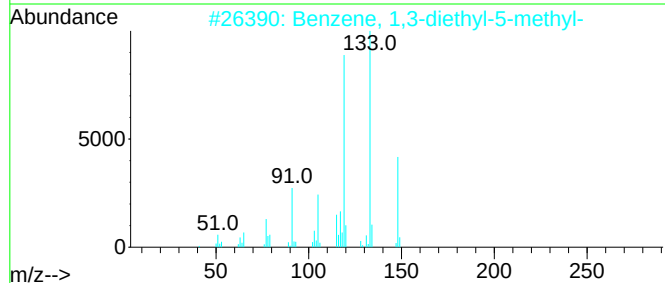
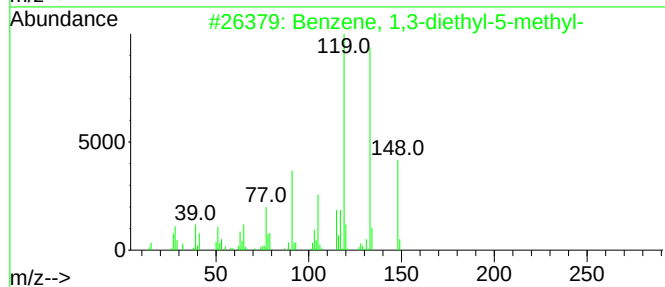
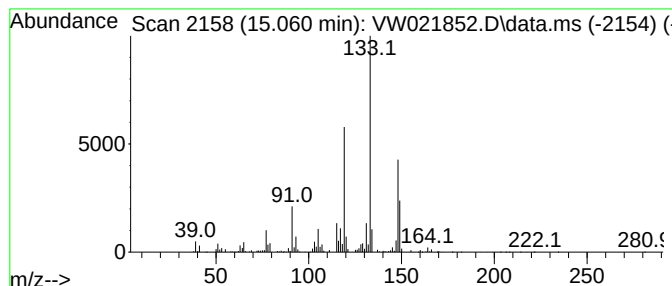
Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

\*\*\*\*\*  
Peak Number 65 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 67

| R.T.      | EstConc                             | Area    | Relative to ISTD       |         |             | R.T.   |
|-----------|-------------------------------------|---------|------------------------|---------|-------------|--------|
| 15.060    | 5.04 ug/L                           | 1433970 | 1,4-Dichlorobenzene-d4 |         |             | 13.560 |
| Hit# of 5 | Tentative ID                        |         | MW                     | MolForm | CAS#        | Qual   |
| 1         | Benzene, 1,3-diethyl-5-methyl-      |         | 148                    | C11H16  | 002050-24-0 | 91     |
| 2         | Benzene, 1,3-diethyl-5-methyl-      |         | 148                    | C11H16  | 002050-24-0 | 87     |
| 3         | Benzene, 1,4-diethyl-2-methyl-      |         | 148                    | C11H16  | 013632-94-5 | 87     |
| 4         | Benzene, pentamethyl-               |         | 148                    | C11H16  | 000700-12-9 | 70     |
| 5         | Benzene, 1-ethyl-4-(1-methylethyl)- |         | 148                    | C11H16  | 004218-48-8 | 64     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

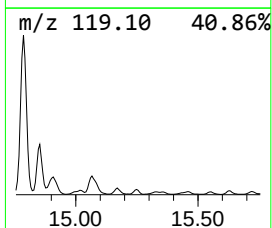
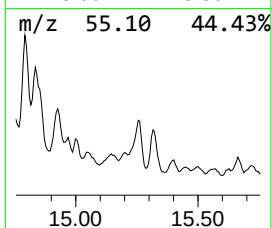
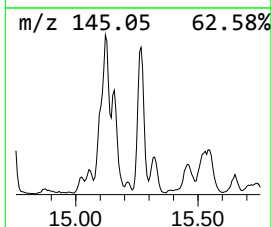
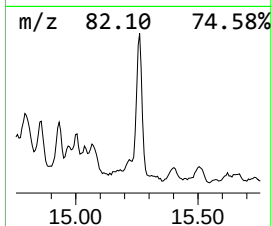
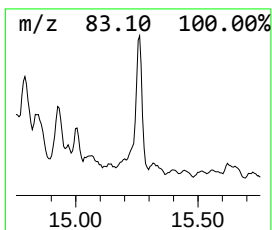
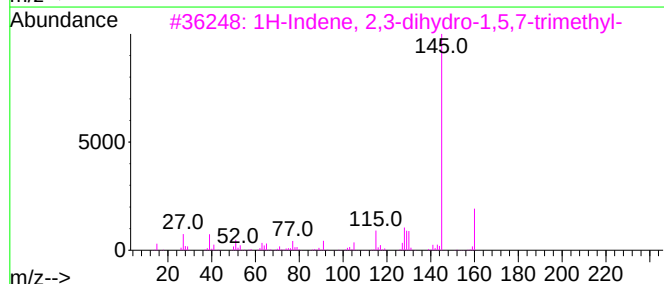
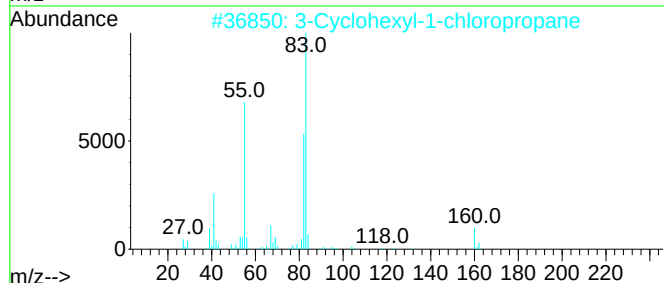
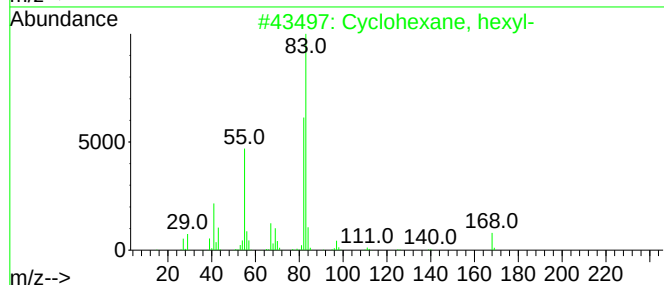
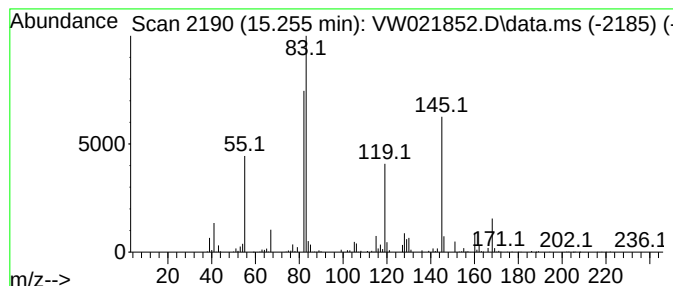
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Quant Title : SFAM01.0

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

\*\*\*\*\*  
Peak Number 67 (DEL) Alkane: Cyclic15.255 Concentration Rank 70

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 15.255 | 3.99 ug/L | 1134320 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, hexyl-                 | 168 | C12H24  | 004292-75-5 | 38   |
| 2    |    |   | 3-Cyclohexyl-1-chloropropane        | 160 | C9H17Cl | 001124-62-5 | 35   |
| 3    |    |   | 1H-Indene, 2,3-dihydro-1,5,7-tri... | 160 | C12H16  | 054340-88-4 | 25   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-1,1,4-tri... | 160 | C12H16  | 016204-72-1 | 25   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-1,1,5-tri... | 160 | C12H16  | 040650-41-7 | 25   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

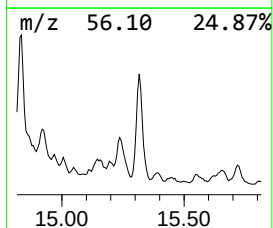
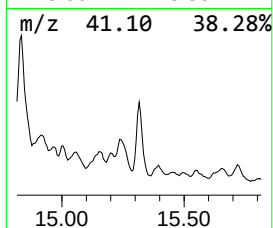
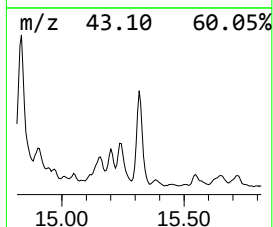
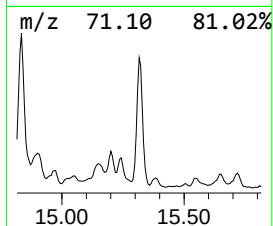
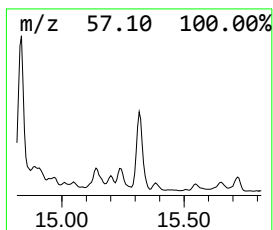
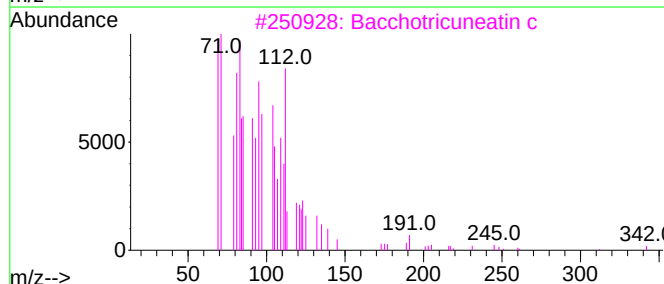
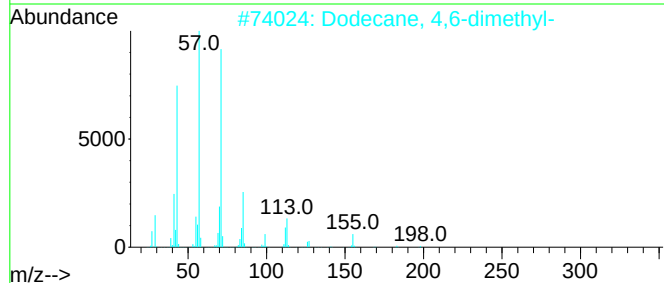
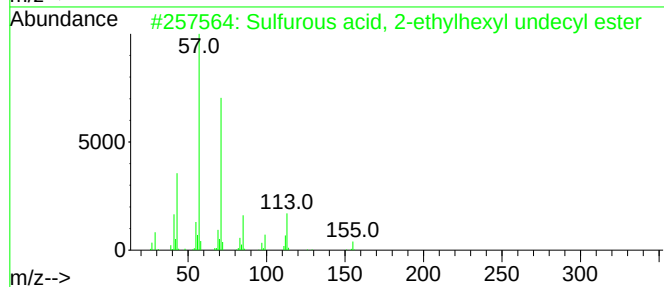
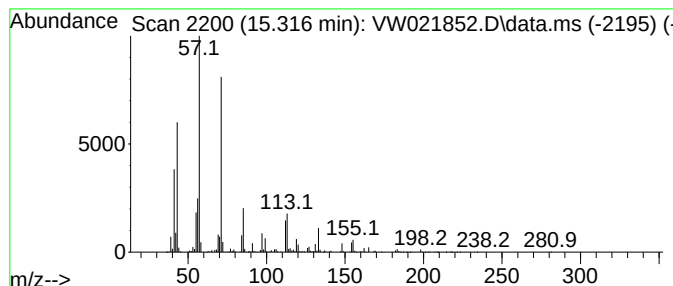
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 68 Sulfurous acid, 2-ethylhexy... Concentration Rank 66

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 15.316 | 5.48 ug/L | 1559320 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, 2-ethylhexyl und... | 348 | C19H40O3S | 1000309-19-4 | 72   |
| 2    |    |   | Dodecane, 4,6-dimethyl-             | 198 | C14H30    | 061141-72-8  | 64   |
| 3    |    |   | Bacchotricuneatin c                 | 342 | C20H22O5  | 066563-30-2  | 64   |
| 4    |    |   | Tridecane, 7-methyl-                | 198 | C14H30    | 026730-14-3  | 62   |
| 5    |    |   | Dodecane, 2,6,11-trimethyl-         | 212 | C15H32    | 031295-56-4  | 59   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

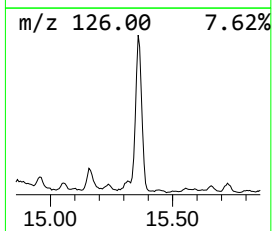
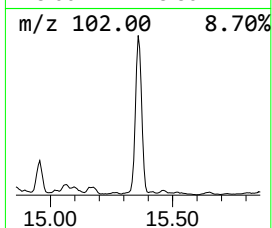
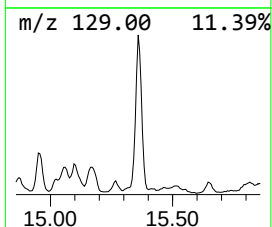
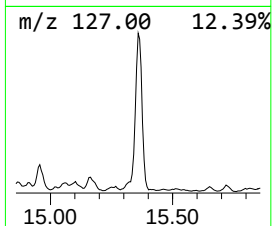
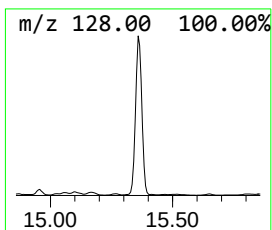
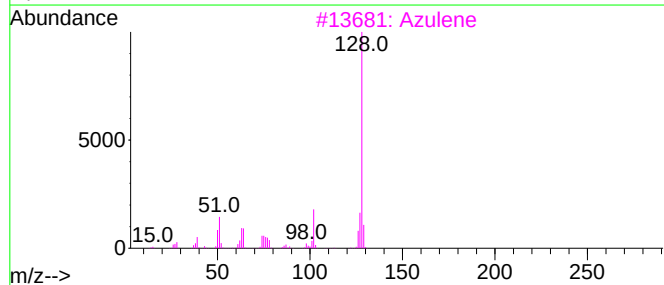
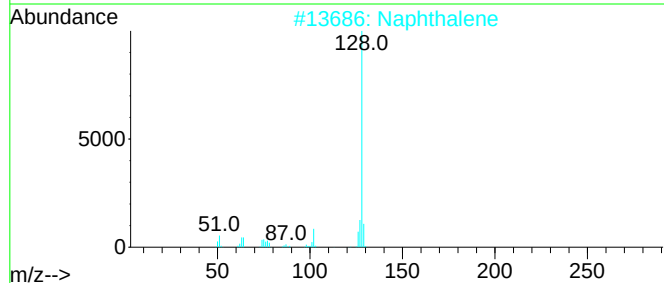
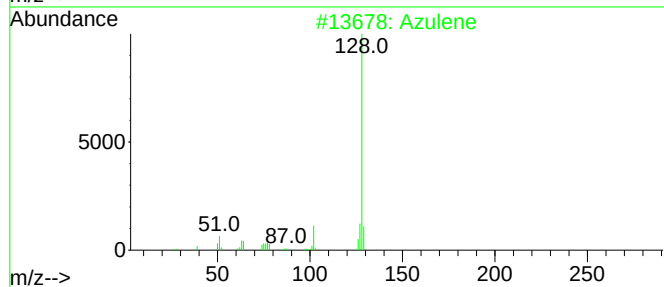
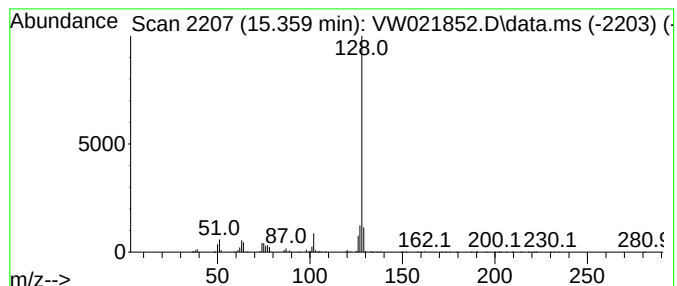
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Quant Title : SFAM01.0

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

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Peak Number 69 Azulene Concentration Rank 56

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 15.359 | 8.36 ug/L | 2380700 | 1,4-Dichlorobenzene-d4 | 13.560 |

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

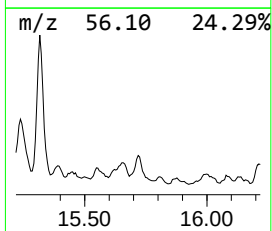
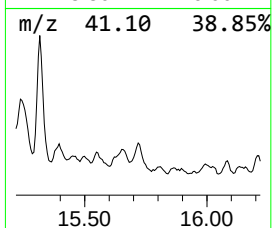
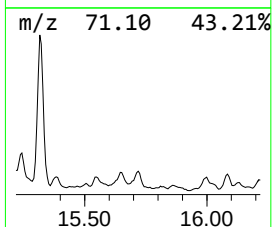
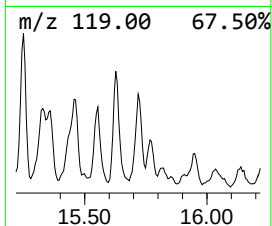
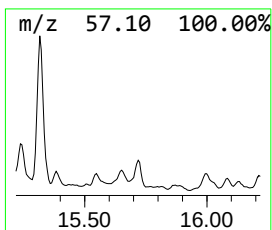
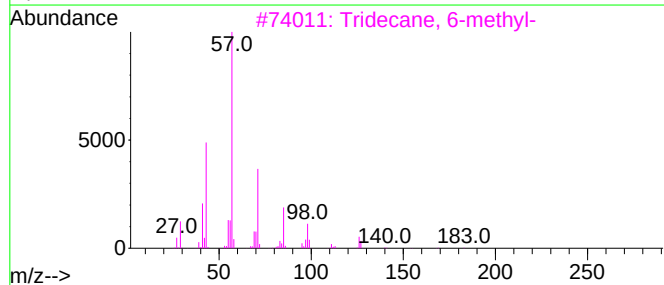
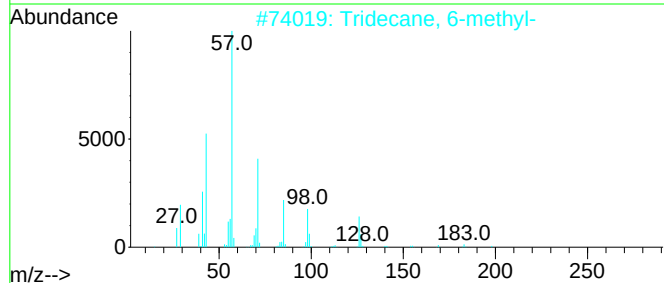
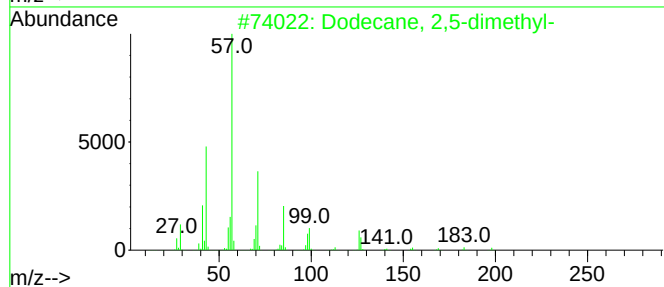
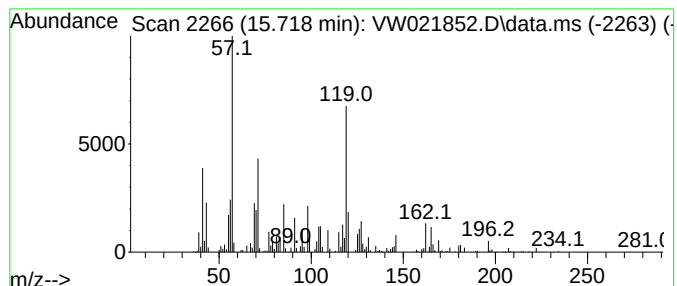
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Quant Title : SFAM01.0

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

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Peak Number 71 (DEL) Alkane: Straight-Chai... Concentration Rank 73

| R.T.      | EstConc                             | Area   | Relative to ISTD       |         | R.T.        |      |
|-----------|-------------------------------------|--------|------------------------|---------|-------------|------|
| 15.718    | 1.68 ug/L                           | 476765 | 1,4-Dichlorobenzene-d4 |         | 13.560      |      |
| Hit# of 5 | Tentative ID                        |        | MW                     | MolForm | CAS#        | Qual |
| 1         | Dodecane, 2,5-dimethyl-             |        | 198                    | C14H30  | 056292-65-0 | 38   |
| 2         | Tridecane, 6-methyl-                |        | 198                    | C14H30  | 013287-21-3 | 38   |
| 3         | Tridecane, 6-methyl-                |        | 198                    | C14H30  | 013287-21-3 | 35   |
| 4         | Benzene, 1,4-dimethyl-2-(2-methy... |        | 162                    | C12H18  | 055669-88-0 | 25   |
| 5         | 3-Ethyl-3-methylheptane             |        | 142                    | C10H22  | 017302-01-1 | 22   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

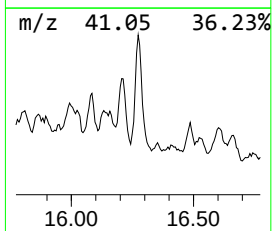
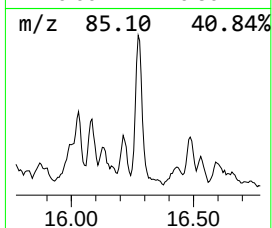
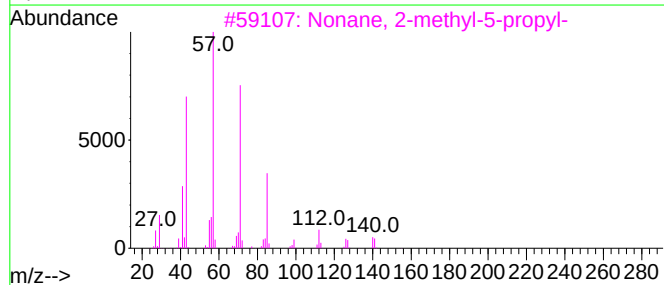
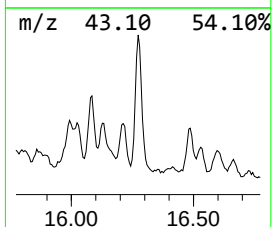
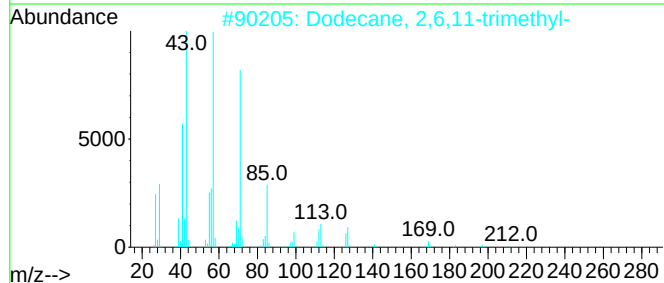
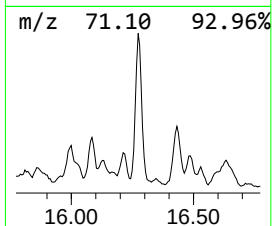
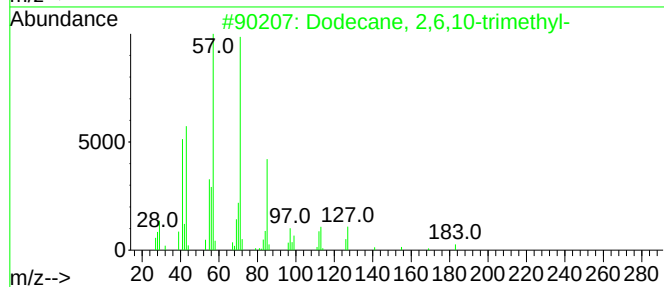
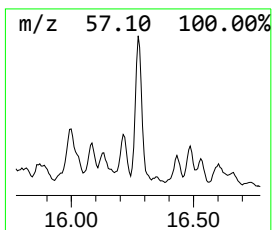
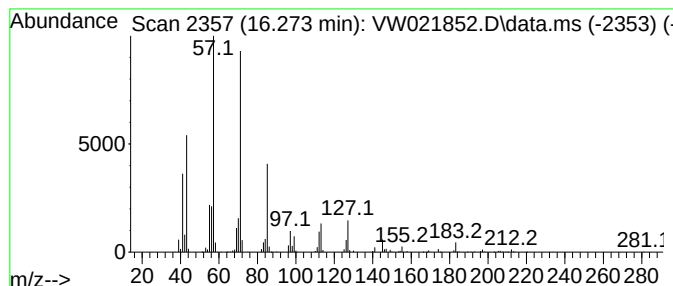
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 72 (DEL) Alkane: Straight-Chai... Concentration Rank 72

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 16.273 | 1.93 ug/L | 550503 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3 | 90   |
| 2    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 86   |
| 3    |    |   | Nonane, 2-methyl-5-propyl-  | 184 | C13H28  | 031081-17-1 | 72   |
| 4    |    |   | Tetradecane, 4-methyl-      | 212 | C15H32  | 025117-24-2 | 72   |
| 5    |    |   | Dodecane, 4,6-dimethyl-     | 198 | C14H30  | 061141-72-8 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

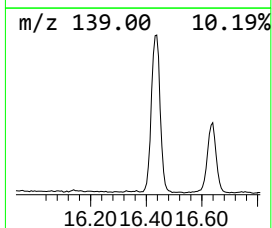
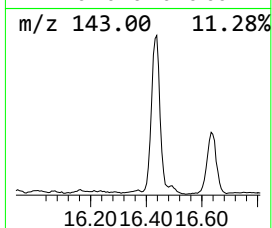
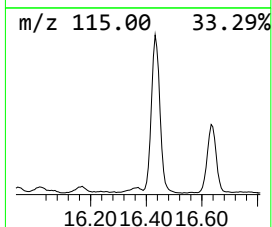
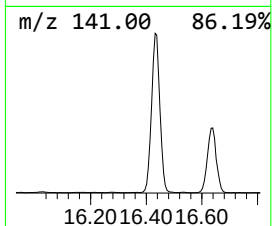
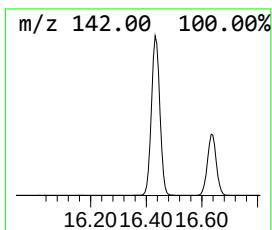
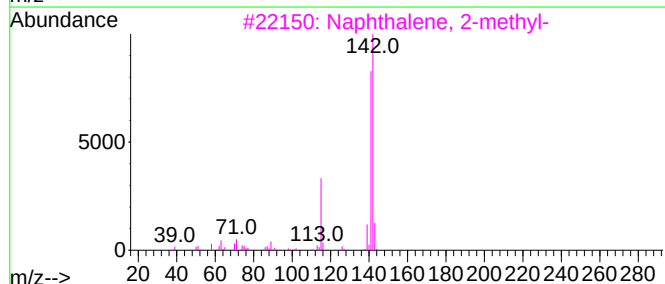
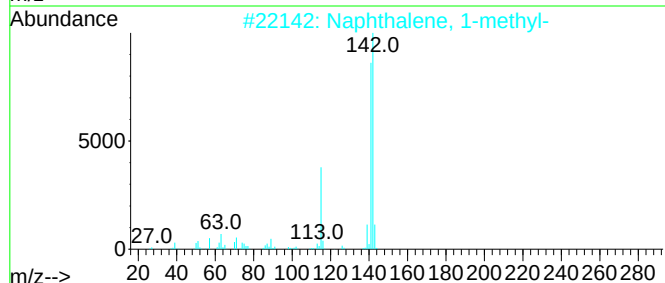
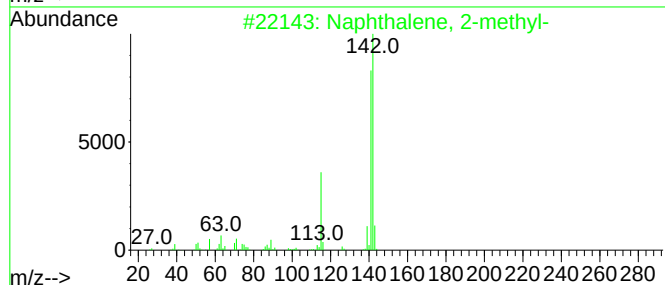
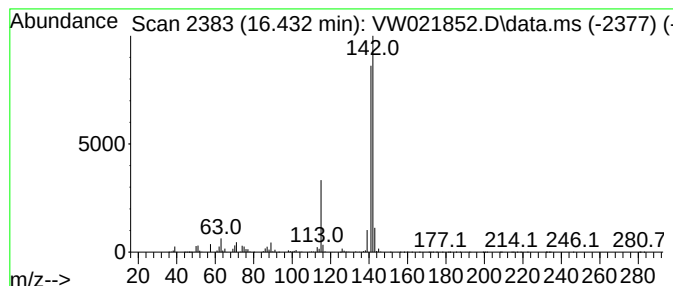
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

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Peak Number 73 Naphthalene, 2-methyl- Concentration Rank 52

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 16.432 | 11.27 ug/L | 3208210 | 1,4-Dichlorobenzene-d4 | 13.560 |

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
Data File : VW021852.D  
Acq On : 11 Jan 2022 23:20  
Operator : SY/VA  
Sample : N1084-10  
Misc : 6.57g/10mL/MSVOA\_W/SOIL  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
BGLN3

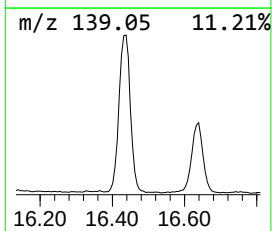
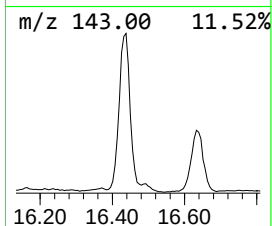
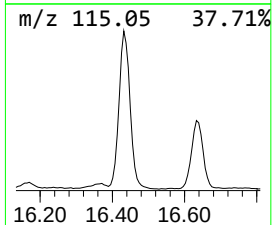
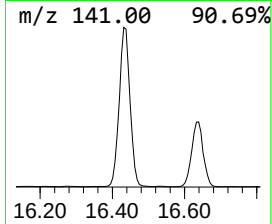
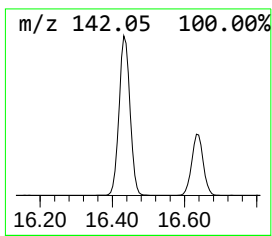
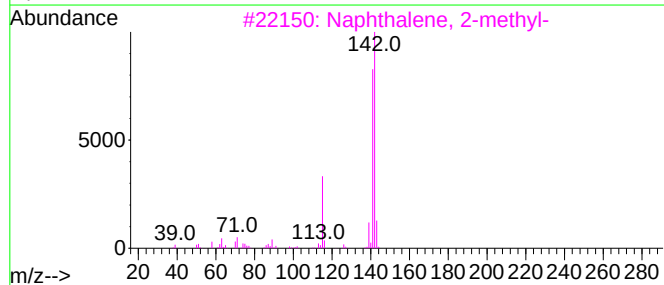
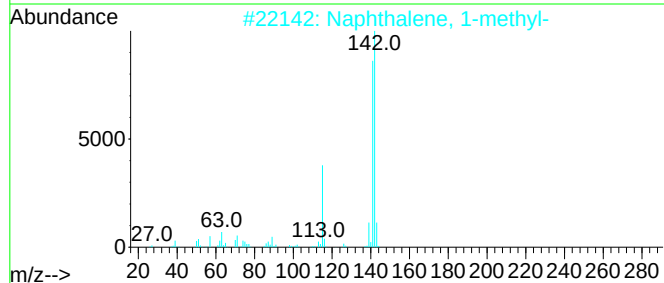
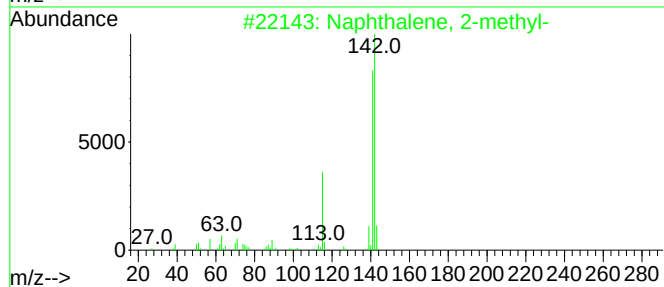
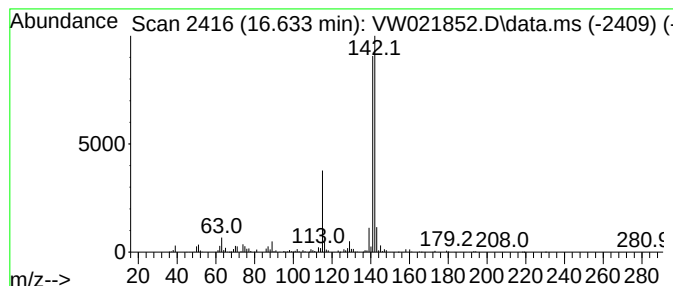
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
Quant Title : SFAM01.0

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

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Peak Number 74 Naphthalene, 1-methyl- Concentration Rank 63

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 16.633 | 6.40 ug/L | 1821160 | 1,4-Dichlorobenzene-d4 | 13.560 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 96   |
| 2    |    |   | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 96   |
| 3    |    |   | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 94   |
| 4    |    |   | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 94   |
| 5    |    |   | Benzocycloheptatriene  | 142 | C11H10  | 000264-09-5 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_W\Data\VW011122\  
 Data File : VW021852.D  
 Acq On : 11 Jan 2022 23:20  
 Operator : SY/VA  
 Sample : N1084-10  
 Misc : 6.57g/10mL/MSVOA\_W/SOIL  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_W  
 ClientSampleId :  
 BGLN3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_W\Method\SFAMWLM011122SMA.M  
 Quant Title : SFAM01.0

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... | 3.026  | 8.3     | ug/L  | 159459   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 3.385  | 9.3     | ug/L  | 179967   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 4.964  | 37.0    | ug/L  | 715506   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 5.440  | 28.2    | ug/L  | 544829   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 5.940  | 55.5    | ug/L  | 1072410  | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: C... | 6.988  | 44.7    | ug/L  | 863296   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 8.159  | 31.1    | ug/L  | 601585   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 8.433  | 9.5     | ug/L  | 184204   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 8.488  | 12.4    | ug/L  | 239438   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 8.695  | 44.7    | ug/L  | 863839   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 9.957  | 19.7    | ug/L  | 379886   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 10.097 | 30.6    | ug/L  | 591973   | 1                     | 8.842  | 483139  | 25.0 |
| (DEL) Alkane: S... | 10.506 | 31.3    | ug/L  | 1034150  | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: C... | 10.671 | 19.3    | ug/L  | 638517   | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: C... | 10.762 | 22.8    | ug/L  | 752599   | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: S... | 10.847 | 10.2    | ug/L  | 337028   | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: S... | 11.042 | 30.1    | ug/L  | 993184   | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: C... | 11.116 | 41.0    | ug/L  | 1354570  | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: C... | 11.183 | 32.7    | ug/L  | 1081180  | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: C... | 11.238 | 32.1    | ug/L  | 1059770  | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: C... | 11.286 | 6.5     | ug/L  | 214810   | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: S... | 11.347 | 31.7    | ug/L  | 1046640  | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: S... | 11.420 | 206.3   | ug/L  | 6818500  | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: S... | 11.524 | 112.4   | ug/L  | 3716140  | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: C... | 11.939 | 60.7    | ug/L  | 2004810  | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: C... | 11.981 | 8.2     | ug/L  | 271132   | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: S... | 12.012 | 7.5     | ug/L  | 246736   | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: S... | 12.067 | 112.3   | ug/L  | 3713440  | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: S... | 12.256 | 277.9   | ug/L  | 9187030  | 2                     | 11.634 | 826306  | 25.0 |
| unknown-01         | 12.341 | 275.1   | ug/L  | 9090940  | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: C... | 12.384 | 395.2   | ug/L  | 13062200 | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: S... | 12.548 | 346.4   | ug/L  | 11448100 | 2                     | 11.634 | 826306  | 25.0 |
| (DEL) Alkane: S... | 12.573 | 391.4   | ug/L  | 12936600 | 2                     | 11.634 | 826306  | 25.0 |
| Benzene, propyl-   | 12.804 | 96.8    | ug/L  | 27538700 | 3                     | 13.560 | 7115280 | 25.0 |
| Benzene, 1-ethy... | 12.871 | 269.8   | ug/L  | 76783300 | 3                     | 13.560 | 7115280 | 25.0 |
| Ethanone, 1-(1-... | 12.999 | 23.7    | ug/L  | 6747850  | 3                     | 13.560 | 7115280 | 25.0 |
| Benzene, 1-ethy... | 13.109 | 165.8   | ug/L  | 47195800 | 3                     | 13.560 | 7115280 | 25.0 |
| (DEL) Alkane: S... | 13.170 | 83.2    | ug/L  | 23688600 | 3                     | 13.560 | 7115280 | 25.0 |
| (DEL) Alkane: S... | 13.353 | 22.4    | ug/L  | 6360010  | 3                     | 13.560 | 7115280 | 25.0 |
| unknown-02         | 13.383 | 5.5     | ug/L  | 1570110  | 3                     | 13.560 | 7115280 | 25.0 |
| (DEL) Alkane: S... | 13.420 | 19.5    | ug/L  | 5547820  | 3                     | 13.560 | 7115280 | 25.0 |
| p-Cymene           | 13.451 | 53.6    | ug/L  | 15243300 | 3                     | 13.560 | 7115280 | 25.0 |
| (DEL) Alkane: S... | 13.493 | 35.4    | ug/L  | 10064200 | 3                     | 13.560 | 7115280 | 25.0 |
| Benzene, 1,3-di... | 13.719 | 28.9    | ug/L  | 8226670  | 3                     | 13.560 | 7115280 | 25.0 |
| Benzene, cyclop... | 13.755 | 114.5   | ug/L  | 32600600 | 3                     | 13.560 | 7115280 | 25.0 |
| Benzene, 2-ethy... | 13.792 | 94.7    | ug/L  | 26939300 | 3                     | 13.560 | 7115280 | 25.0 |
| unknown-03         | 13.944 | 55.8    | ug/L  | 15880900 | 3                     | 13.560 | 7115280 | 25.0 |
| Benzene, 1-meth... | 14.018 | 33.0    | ug/L  | 9394320  | 3                     | 13.560 | 7115280 | 25.0 |
| Benzene, 2-ethy... | 14.036 | 43.3    | ug/L  | 12328000 | 3                     | 13.560 | 7115280 | 25.0 |
| Benzene, 4-ethy... | 14.097 | 49.7    | ug/L  | 14138800 | 3                     | 13.560 | 7115280 | 25.0 |
| Benzene, 1,2,3,... | 14.200 | 36.3    | ug/L  | 10343300 | 3                     | 13.560 | 7115280 | 25.0 |
| Adamantane         | 14.261 | 22.6    | ug/L  | 6417840  | 3                     | 13.560 | 7115280 | 25.0 |

|                    |        |      |      |          |   |        |         |      |
|--------------------|--------|------|------|----------|---|--------|---------|------|
| o-Cymene           | 14.341 | 29.9 | ug/L | 8498300  | 3 | 13.560 | 7115280 | 25.0 |
| (DEL) Alkane: C... | 14.389 | 36.2 | ug/L | 10297500 | 3 | 13.560 | 7115280 | 25.0 |
| Benzene, 1-ethy... | 14.420 | 34.4 | ug/L | 9775250  | 3 | 13.560 | 7115280 | 25.0 |
| Benzene, 1-ethy... | 14.456 | 23.6 | ug/L | 6729340  | 3 | 13.560 | 7115280 | 25.0 |
| (DEL) Alkane: S... | 14.493 | 25.1 | ug/L | 7135440  | 3 | 13.560 | 7115280 | 25.0 |
| unknown-04         | 14.542 | 14.2 | ug/L | 4041160  | 3 | 13.560 | 7115280 | 25.0 |
| unknown-05         | 14.639 | 6.1  | ug/L | 1740220  | 3 | 13.560 | 7115280 | 25.0 |
| Benzene, 1-ethe... | 14.688 | 4.8  | ug/L | 1373180  | 3 | 13.560 | 7115280 | 25.0 |
| Benzene, (1,1-d... | 14.725 | 7.0  | ug/L | 2000090  | 3 | 13.560 | 7115280 | 25.0 |
| Benzene, 1,2,4,... | 14.792 | 29.0 | ug/L | 8262630  | 3 | 13.560 | 7115280 | 25.0 |
| Naphthalene, 1,... | 14.956 | 6.7  | ug/L | 1911190  | 3 | 13.560 | 7115280 | 25.0 |
| Benzene, 1,3-di... | 15.060 | 5.0  | ug/L | 1433970  | 3 | 13.560 | 7115280 | 25.0 |
| (DEL) Alkane: C... | 15.255 | 4.0  | ug/L | 1134320  | 3 | 13.560 | 7115280 | 25.0 |
| Sulfurous acid,... | 15.316 | 5.5  | ug/L | 1559320  | 3 | 13.560 | 7115280 | 25.0 |
| Azulene            | 15.359 | 8.4  | ug/L | 2380700  | 3 | 13.560 | 7115280 | 25.0 |
| (DEL) Alkane: S... | 15.718 | 1.7  | ug/L | 476765   | 3 | 13.560 | 7115280 | 25.0 |
| (DEL) Alkane: S... | 16.273 | 1.9  | ug/L | 550503   | 3 | 13.560 | 7115280 | 25.0 |
| Naphthalene, 2-... | 16.432 | 11.3 | ug/L | 3208210  | 3 | 13.560 | 7115280 | 25.0 |
| Naphthalene, 1-... | 16.633 | 6.4  | ug/L | 1821160  | 3 | 13.560 | 7115280 | 25.0 |

Instrument :

MSVOA\_W

ClientSampleId :

BGLN3