

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
 Data File : VX041344.D  
 Acq On : 02 May 2024 11:32  
 Operator : JC/MD  
 Sample : P2264-14  
 Misc : 5.0mL/MSVOA\_X/WATER  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 T2-MW-3-8.0-042424

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
 Title : SW846 8260

Signal : TIC: VX041344.D\data.ms

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 1.581       | 68            | 81          | 82           | rBV5     | 7209           | 20382         | 1.65%           | 0.156%        |
| 2         | 2.770       | 268           | 276         | 279          | rBV      | 25230          | 52436         | 4.25%           | 0.402%        |
| 3         | 2.819       | 279           | 284         | 305          | rVB      | 60423          | 147031        | 11.93%          | 1.128%        |
| 4         | 3.093       | 320           | 329         | 342          | rBV      | 65181          | 143086        | 11.61%          | 1.098%        |
| 5         | 3.721       | 420           | 432         | 441          | rBV2     | 14174          | 35999         | 2.92%           | 0.276%        |
| 6         | 3.831       | 441           | 450         | 466          | rVB3     | 11584          | 30300         | 2.46%           | 0.232%        |
| 7         | 4.093       | 487           | 493         | 501          | rVB4     | 6960           | 16536         | 1.34%           | 0.127%        |
| 8         | 4.209       | 501           | 512         | 517          | rVV2     | 9048           | 25744         | 2.09%           | 0.197%        |
| 9         | 4.294       | 517           | 526         | 538          | rVV      | 59523          | 169036        | 13.72%          | 1.297%        |
| 10        | 4.422       | 538           | 547         | 558          | rVB3     | 11174          | 33802         | 2.74%           | 0.259%        |
| 11        | 5.129       | 646           | 663         | 664          | rBV3     | 8145           | 23297         | 1.89%           | 0.179%        |
| 12        | 5.178       | 666           | 671         | 684          | rVB6     | 13010          | 42947         | 3.48%           | 0.329%        |
| 13        | 5.385       | 693           | 705         | 712          | rBV2     | 84750          | 248237        | 20.14%          | 1.904%        |
| 14        | 5.471       | 712           | 719         | 724          | rVV4     | 19929          | 58476         | 4.75%           | 0.449%        |
| 15        | 5.550       | 724           | 732         | 740          | rVV      | 130545         | 353680        | 28.70%          | 2.713%        |
| 16        | 5.818       | 765           | 776         | 790          | rBV3     | 29944          | 94686         | 7.68%           | 0.726%        |
| 17        | 5.952       | 790           | 798         | 811          | rBV      | 81540          | 218004        | 17.69%          | 1.672%        |
| 18        | 6.153       | 823           | 831         | 840          | rBV4     | 19529          | 68314         | 5.54%           | 0.524%        |
| 19        | 6.251       | 840           | 847         | 856          | rVV3     | 29551          | 86571         | 7.02%           | 0.664%        |
| 20        | 6.355       | 856           | 864         | 877          | rVB3     | 31363          | 89825         | 7.29%           | 0.689%        |
| 21        | 6.598       | 897           | 904         | 914          | rVB2     | 8274           | 20804         | 1.69%           | 0.160%        |
| 22        | 6.763       | 921           | 931         | 940          | rBV      | 231585         | 649845        | 52.73%          | 4.985%        |
| 23        | 6.848       | 941           | 945         | 956          | rVB3     | 56683          | 140918        | 11.43%          | 1.081%        |
| 24        | 6.970       | 958           | 965         | 972          | rBV2     | 13743          | 32107         | 2.61%           | 0.246%        |
| 25        | 7.062       | 973           | 980         | 989          | rBV3     | 9208           | 22286         | 1.81%           | 0.171%        |
| 26        | 7.171       | 989           | 998         | 1006         | rBV      | 73206          | 167106        | 13.56%          | 1.282%        |
| 27        | 7.373       | 1019          | 1031        | 1044         | rVB4     | 75036          | 215490        | 17.49%          | 1.653%        |
| 28        | 7.495       | 1044          | 1051        | 1056         | rBV4     | 12603          | 27817         | 2.26%           | 0.213%        |
| 29        | 7.556       | 1056          | 1061        | 1071         | rVB2     | 12806          | 28364         | 2.30%           | 0.218%        |
| 30        | 7.653       | 1071          | 1077        | 1084         | rBV2     | 15532          | 28753         | 2.33%           | 0.221%        |
| 31        | 7.769       | 1089          | 1096        | 1103         | rVB2     | 17167          | 36082         | 2.93%           | 0.277%        |
| 32        | 7.879       | 1103          | 1114        | 1122         | rBV4     | 23334          | 82278         | 6.68%           | 0.631%        |
| 33        | 7.982       | 1122          | 1131        | 1142         | rVB4     | 23308          | 77911         | 6.32%           | 0.598%        |
| 34        | 8.104       | 1142          | 1151        | 1153         | rBV      | 52417          | 96459         | 7.83%           | 0.740%        |
| 35        | 8.141       | 1153          | 1157        | 1162         | rVV      | 123852         | 218158        | 17.70%          | 1.673%        |
| 36        | 8.196       | 1162          | 1166        | 1176         | rVB3     | 23322          | 51133         | 4.15%           | 0.392%        |

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 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 T2-MW-3-8.0-042424

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
 Title : SW846 8260

|    |        |      |      |      |      |        |         |         |        |
|----|--------|------|------|------|------|--------|---------|---------|--------|
| 37 | 8.305  | 1176 | 1184 | 1192 | rBV3 | 24229  | 48876   | 3.97%   | 0.375% |
| 38 | 8.433  | 1200 | 1205 | 1209 | rBV3 | 10764  | 18557   | 1.51%   | 0.142% |
| 39 | 8.501  | 1209 | 1216 | 1226 | rVB  | 87208  | 152880  | 12.41%  | 1.173% |
| 40 | 8.598  | 1226 | 1232 | 1235 | rBV4 | 17339  | 33867   | 2.75%   | 0.260% |
| 41 | 8.647  | 1235 | 1240 | 1249 | rVB  | 469077 | 730455  | 59.27%  | 5.603% |
| 42 | 8.818  | 1264 | 1268 | 1271 | rBV3 | 8989   | 15502   | 1.26%   | 0.119% |
| 43 | 8.921  | 1280 | 1285 | 1290 | rBV  | 28640  | 46050   | 3.74%   | 0.353% |
| 44 | 8.970  | 1290 | 1293 | 1298 | rVB2 | 12517  | 18181   | 1.48%   | 0.139% |
| 45 | 9.098  | 1310 | 1314 | 1319 | rVV2 | 20679  | 32872   | 2.67%   | 0.252% |
| 46 | 9.159  | 1319 | 1324 | 1334 | rVB  | 104546 | 171471  | 13.91%  | 1.315% |
| 47 | 9.330  | 1345 | 1352 | 1356 | rBV2 | 8637   | 14726   | 1.19%   | 0.113% |
| 48 | 9.378  | 1356 | 1360 | 1364 | rVV3 | 12108  | 17857   | 1.45%   | 0.137% |
| 49 | 9.427  | 1364 | 1368 | 1370 | rVV  | 25568  | 33618   | 2.73%   | 0.258% |
| 50 | 9.458  | 1370 | 1373 | 1376 | rVV  | 37900  | 54486   | 4.42%   | 0.418% |
| 51 | 9.500  | 1376 | 1380 | 1386 | rVB4 | 21713  | 43885   | 3.56%   | 0.337% |
| 52 | 9.604  | 1393 | 1397 | 1402 | rVV  | 18971  | 25622   | 2.08%   | 0.197% |
| 53 | 9.659  | 1402 | 1406 | 1410 | rVB  | 14665  | 21736   | 1.76%   | 0.167% |
| 54 | 9.762  | 1416 | 1423 | 1431 | rBV3 | 33226  | 63717   | 5.17%   | 0.489% |
| 55 | 9.921  | 1443 | 1449 | 1454 | rBV  | 9664   | 17040   | 1.38%   | 0.131% |
| 56 | 10.055 | 1466 | 1471 | 1476 | rBV  | 539042 | 727063  | 59.00%  | 5.577% |
| 57 | 10.153 | 1485 | 1487 | 1491 | rVB  | 18713  | 22504   | 1.83%   | 0.173% |
| 58 | 10.220 | 1491 | 1498 | 1502 | rVV3 | 9739   | 23957   | 1.94%   | 0.184% |
| 59 | 10.268 | 1502 | 1506 | 1509 | rVV2 | 13933  | 23504   | 1.91%   | 0.180% |
| 60 | 10.305 | 1509 | 1512 | 1517 | rVB2 | 20962  | 31297   | 2.54%   | 0.240% |
| 61 | 10.464 | 1535 | 1538 | 1546 | rVB4 | 7734   | 15116   | 1.23%   | 0.116% |
| 62 | 10.586 | 1546 | 1558 | 1564 | rBV  | 16884  | 34852   | 2.83%   | 0.267% |
| 63 | 10.775 | 1581 | 1589 | 1592 | rBV6 | 8536   | 17550   | 1.42%   | 0.135% |
| 64 | 10.878 | 1598 | 1606 | 1614 | rVB7 | 10730  | 29137   | 2.36%   | 0.224% |
| 65 | 10.963 | 1614 | 1620 | 1627 | rBV  | 357902 | 474751  | 38.52%  | 3.642% |
| 66 | 11.079 | 1634 | 1639 | 1651 | rVB  | 469188 | 616679  | 50.04%  | 4.730% |
| 67 | 11.305 | 1670 | 1676 | 1685 | rBV  | 495949 | 616308  | 50.01%  | 4.727% |
| 68 | 11.610 | 1722 | 1726 | 1735 | rVB  | 39586  | 56247   | 4.56%   | 0.431% |
| 69 | 11.866 | 1763 | 1768 | 1770 | rBV  | 25200  | 34254   | 2.78%   | 0.263% |
| 70 | 11.890 | 1770 | 1772 | 1782 | rVB  | 25186  | 35077   | 2.85%   | 0.269% |
| 71 | 12.018 | 1788 | 1793 | 1801 | rBV  | 535512 | 704201  | 57.14%  | 5.402% |
| 72 | 12.244 | 1824 | 1830 | 1837 | rBV2 | 963131 | 1232358 | 100.00% | 9.453% |
| 73 | 12.311 | 1837 | 1841 | 1843 | rBV  | 32860  | 42380   | 3.44%   | 0.325% |
| 74 | 12.402 | 1850 | 1856 | 1862 | rBV  | 50939  | 66272   | 5.38%   | 0.508% |
| 75 | 12.463 | 1862 | 1866 | 1871 | rVB  | 11659  | 15262   | 1.24%   | 0.117% |
| 76 | 12.610 | 1883 | 1890 | 1893 | rBV  | 253716 | 310847  | 25.22%  | 2.384% |

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Instrument :  
 MSVOA\_X  
 ClientSampleId :  
 T2-MW-3-8.0-042424

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
 Title : SW846 8260

|     |        |      |      |      |      |        |        |        |        |
|-----|--------|------|------|------|------|--------|--------|--------|--------|
| 77  | 12.646 | 1893 | 1896 | 1900 | rVV  | 99755  | 125397 | 10.18% | 0.962% |
| 78  | 12.695 | 1900 | 1904 | 1912 | rVV  | 354092 | 486215 | 39.45% | 3.730% |
| 79  | 12.841 | 1923 | 1928 | 1933 | rVB  | 23602  | 33841  | 2.75%  | 0.260% |
| 80  | 12.920 | 1933 | 1941 | 1946 | rBV  | 279902 | 348205 | 28.26% | 2.671% |
| 81  | 13.030 | 1954 | 1959 | 1966 | rVB3 | 32512  | 52988  | 4.30%  | 0.406% |
| 82  | 13.134 | 1966 | 1976 | 1978 | rBV2 | 45152  | 71696  | 5.82%  | 0.550% |
| 83  | 13.170 | 1978 | 1982 | 1989 | rVV  | 195302 | 274640 | 22.29% | 2.107% |
| 84  | 13.237 | 1989 | 1993 | 1994 | rVV  | 16458  | 22390  | 1.82%  | 0.172% |
| 85  | 13.292 | 1994 | 2002 | 2007 | rVV2 | 329913 | 487835 | 39.59% | 3.742% |
| 86  | 13.341 | 2007 | 2010 | 2013 | rVV4 | 25925  | 41719  | 3.39%  | 0.320% |
| 87  | 13.372 | 2013 | 2015 | 2020 | rVV2 | 20758  | 29125  | 2.36%  | 0.223% |
| 88  | 13.426 | 2020 | 2024 | 2030 | rVB  | 26477  | 37115  | 3.01%  | 0.285% |
| 89  | 13.506 | 2033 | 2037 | 2040 | rBV2 | 17465  | 28750  | 2.33%  | 0.221% |
| 90  | 13.542 | 2040 | 2043 | 2046 | rVV  | 54797  | 76109  | 6.18%  | 0.584% |
| 91  | 13.579 | 2046 | 2049 | 2053 | rVV3 | 51097  | 74287  | 6.03%  | 0.570% |
| 92  | 13.640 | 2056 | 2059 | 2060 | rVV  | 31671  | 36462  | 2.96%  | 0.280% |
| 93  | 13.664 | 2060 | 2063 | 2074 | rVV2 | 47253  | 75960  | 6.16%  | 0.583% |
| 94  | 13.780 | 2078 | 2082 | 2091 | rVB3 | 12484  | 28424  | 2.31%  | 0.218% |
| 95  | 13.926 | 2099 | 2106 | 2110 | rBV2 | 16391  | 27822  | 2.26%  | 0.213% |
| 96  | 14.073 | 2126 | 2130 | 2135 | rBV  | 15029  | 18569  | 1.51%  | 0.142% |
| 97  | 14.213 | 2149 | 2153 | 2159 | rBV  | 17274  | 27325  | 2.22%  | 0.210% |
| 98  | 14.317 | 2166 | 2170 | 2176 | rBV  | 20978  | 31472  | 2.55%  | 0.241% |
| 99  | 14.371 | 2176 | 2179 | 2184 | rVB3 | 11522  | 14707  | 1.19%  | 0.113% |
| 100 | 14.780 | 2241 | 2246 | 2255 | rVB2 | 28581  | 40718  | 3.30%  | 0.312% |

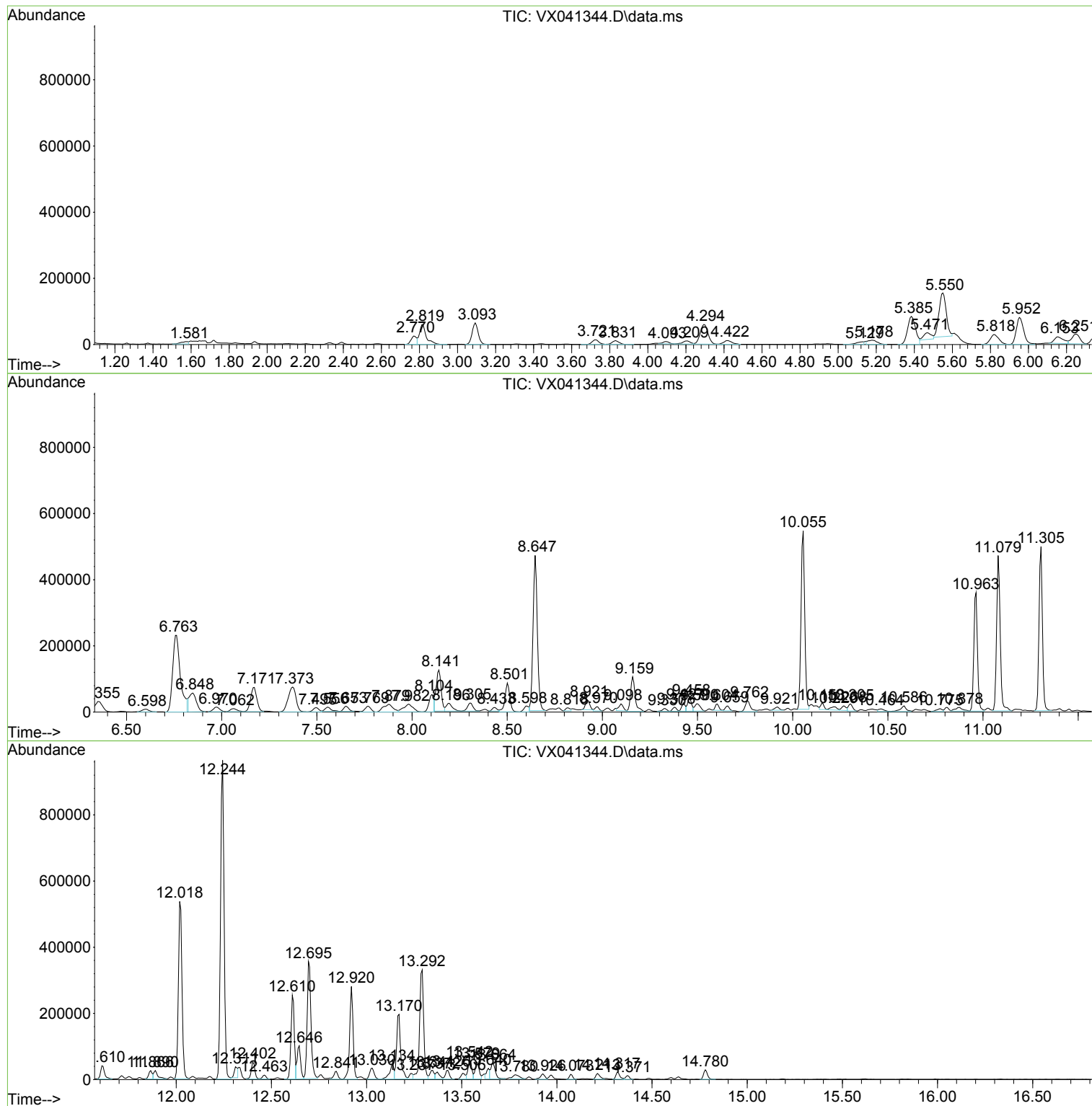
Sum of corrected areas: 13036685

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Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
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Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
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Instrument :  
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ClientSampleId :  
T2-MW-3-8.0-042424

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

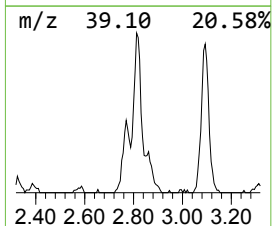
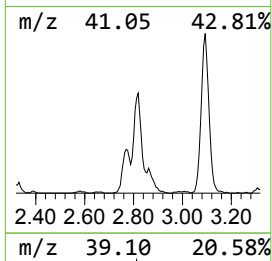
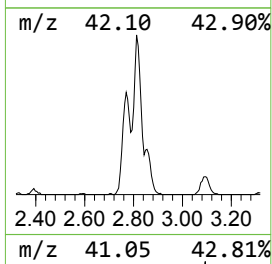
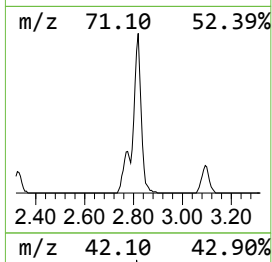
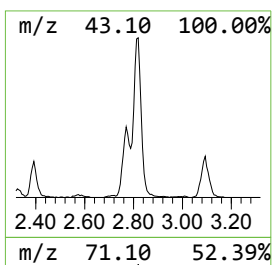
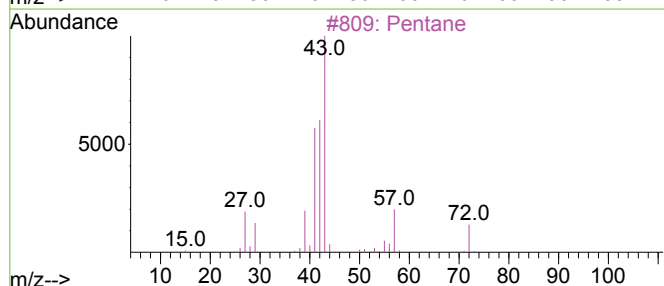
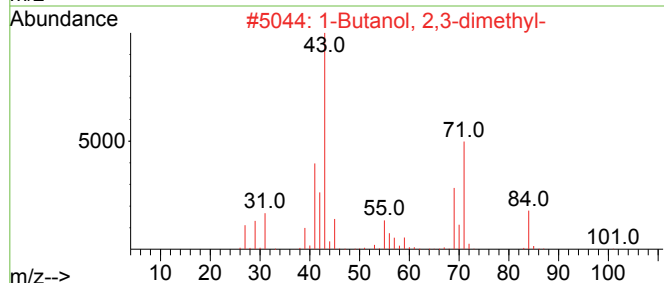
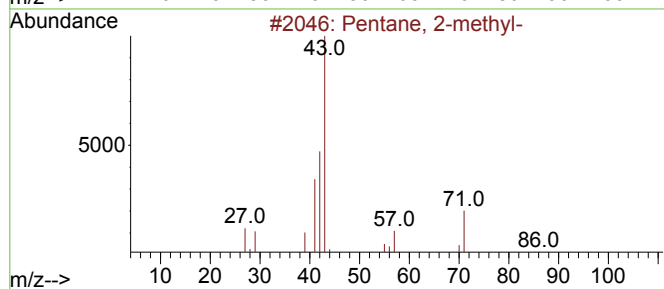
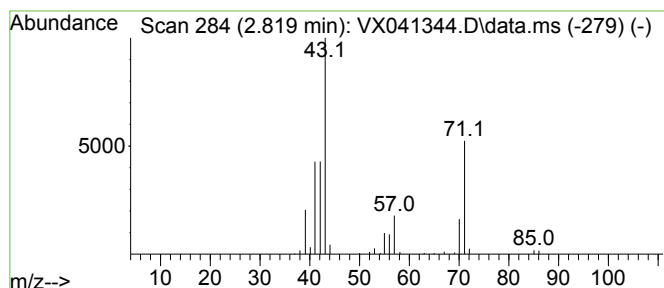
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Pentane, 2-methyl- Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD   | R.T.  |
|-------|------------|--------|--------------------|-------|
| 2.819 | 20.79 ug/l | 147031 | Pentafluorobenzene | 5.550 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentane, 2-methyl-       | 86  | C6H14   | 000107-83-5 | 74   |
| 2    |    |   | 1-Butanol, 2,3-dimethyl- | 102 | C6H14O  | 019550-30-2 | 50   |
| 3    |    |   | Pentane                  | 72  | C5H12   | 000109-66-0 | 43   |
| 4    |    |   | Hexane, 2,3-dimethyl-    | 114 | C8H18   | 000584-94-1 | 40   |
| 5    |    |   | Pentane, 3,3-dimethyl-   | 100 | C7H16   | 000562-49-2 | 38   |



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MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

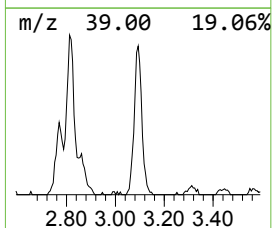
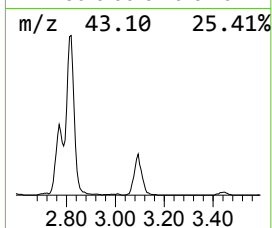
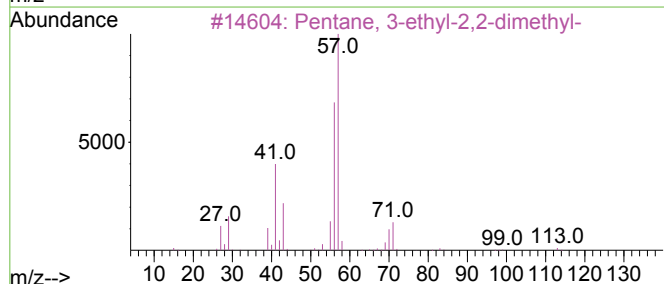
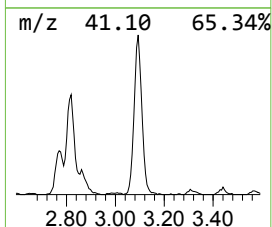
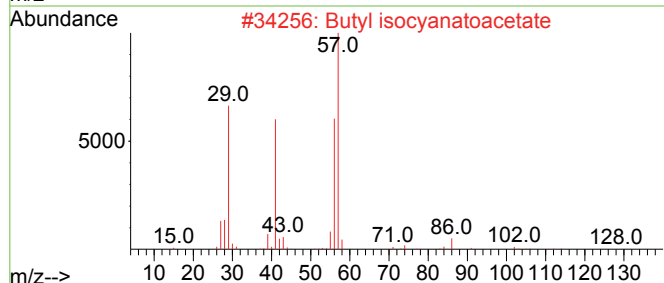
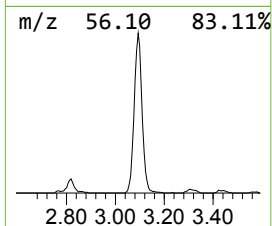
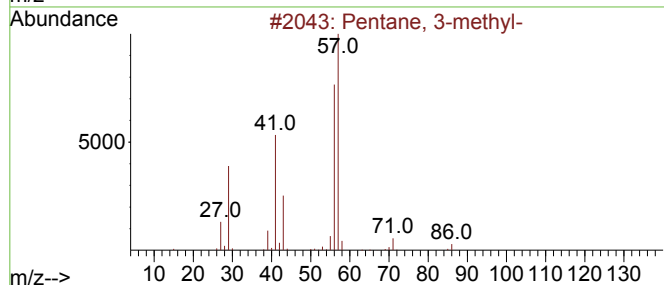
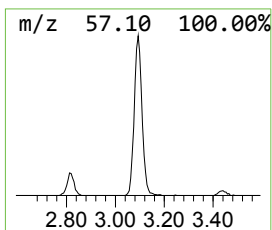
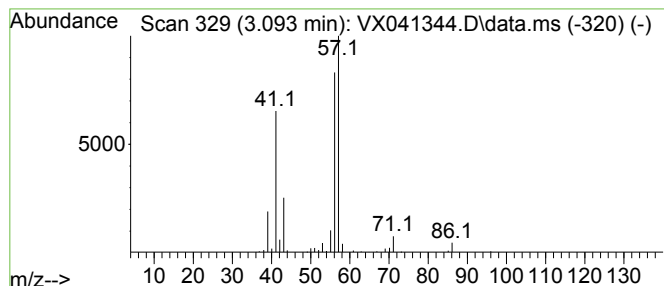
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

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

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Peak Number 2 Pentane, 3-methyl- Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD   | R.T.  |
|-------|------------|--------|--------------------|-------|
| 3.093 | 20.23 ug/l | 143086 | Pentafluorobenzene | 5.550 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Pentane, 3-methyl-                 | 86  | C6H14    | 000096-14-0 | 91   |
| 2    |    |   | Butyl isocyanatoacetate            | 157 | C7H11NO3 | 017046-22-9 | 64   |
| 3    |    |   | Pentane, 3-ethyl-2,2-dimethyl-     | 128 | C9H20    | 016747-32-3 | 64   |
| 4    |    |   | Oxirane, (1-methylethyl)-          | 86  | C5H10O   | 001438-14-8 | 59   |
| 5    |    |   | 3,4-Dimethyldihydrofuran-2,5-dione | 128 | C6H8O3   | 007475-92-5 | 53   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

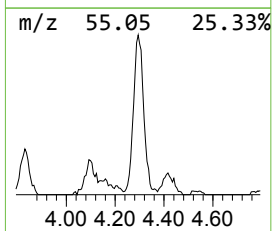
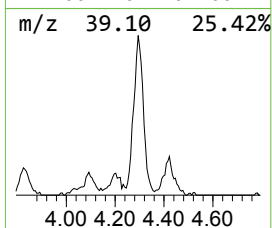
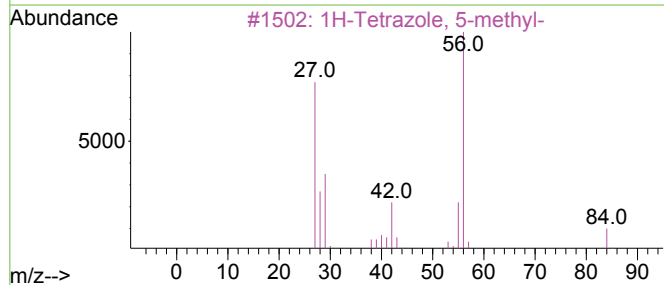
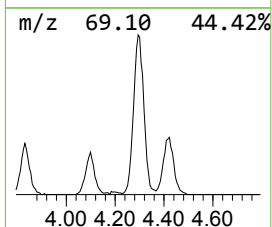
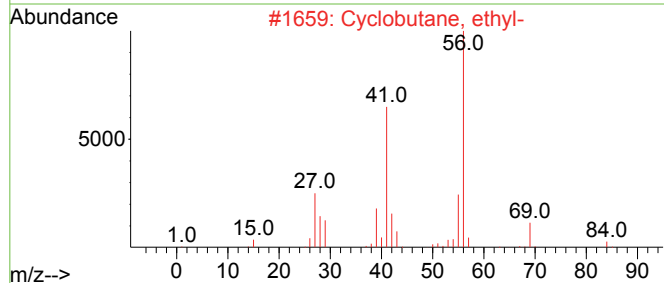
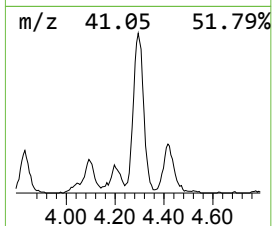
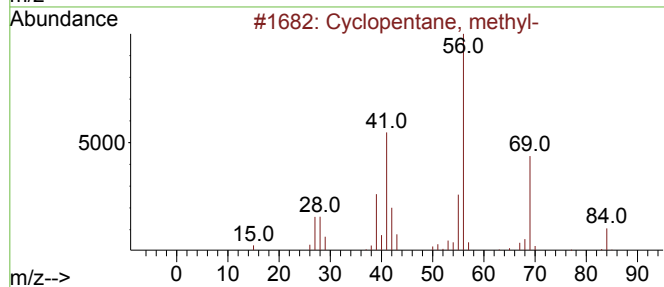
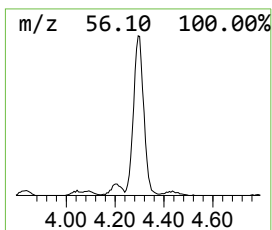
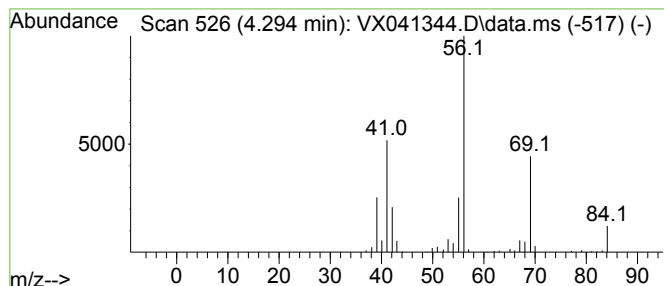
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Quant Title : SW846 8260

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

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Peak Number 3 Cyclopentane, methyl- Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD   | R.T.  |
|-------|------------|--------|--------------------|-------|
| 4.294 | 23.90 ug/l | 169036 | Pentafluorobenzene | 5.550 |

| Hit# | of | 5 | Tentative ID                        | MW | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|----|---------|-------------|------|
| 1    |    |   | Cyclopentane, methyl-               | 84 | C6H12   | 000096-37-7 | 91   |
| 2    |    |   | Cyclobutane, ethyl-                 | 84 | C6H12   | 004806-61-5 | 72   |
| 3    |    |   | 1H-Tetrazole, 5-methyl-             | 84 | C2H4N4  | 004076-36-2 | 64   |
| 4    |    |   | Cyclobutane                         | 56 | C4H8    | 000287-23-0 | 50   |
| 5    |    |   | Cyclopropane, 1-ethyl-2-methyl-,... | 84 | C6H12   | 019781-68-1 | 38   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

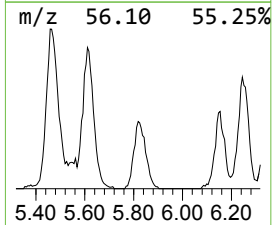
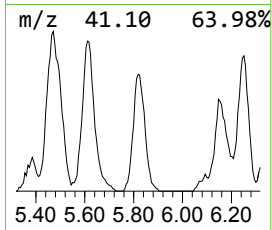
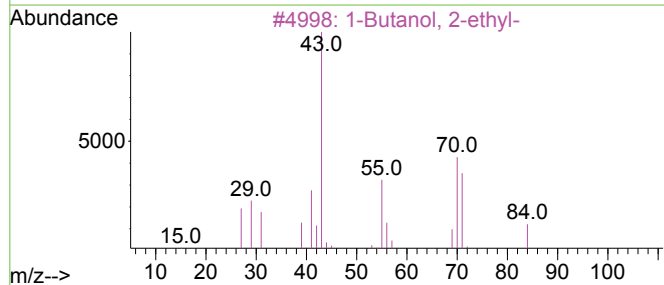
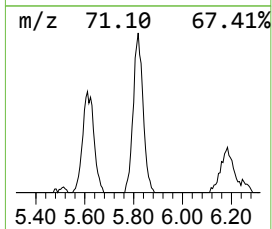
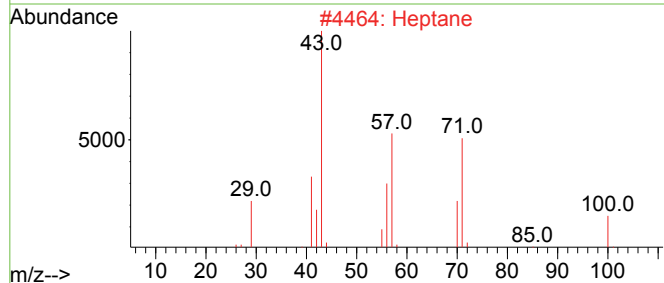
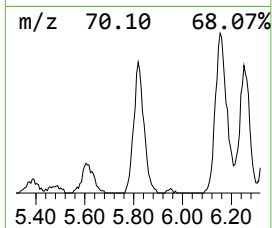
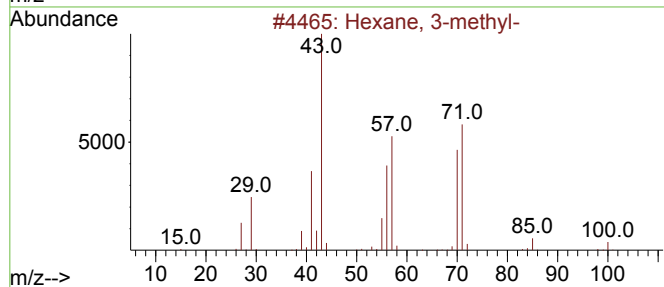
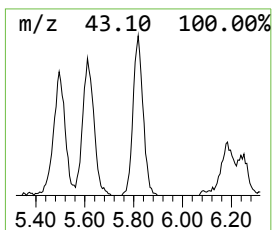
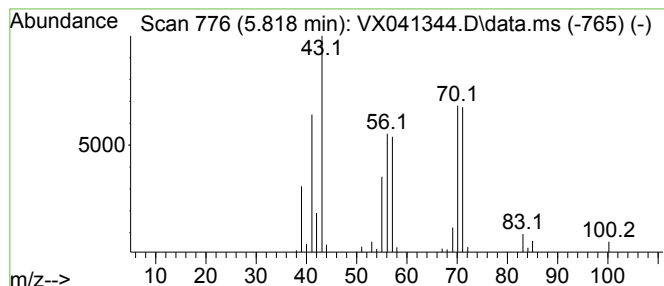
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 Hexane, 3-methyl- Concentration Rank 11

| R.T.  | EstConc    | Area  | Relative to ISTD   | R.T.  |
|-------|------------|-------|--------------------|-------|
| 5.818 | 13.39 ug/l | 94686 | Pentafluorobenzene | 5.550 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Hexane, 3-methyl-     | 100 | C7H16   | 000589-34-4 | 90   |
| 2    |    |   | Heptane               | 100 | C7H16   | 000142-82-5 | 58   |
| 3    |    |   | 1-Butanol, 2-ethyl-   | 102 | C6H14O  | 000097-95-0 | 50   |
| 4    |    |   | Pyrrolidine           | 71  | C4H9N   | 000123-75-1 | 47   |
| 5    |    |   | Hexane, 2,3-dimethyl- | 114 | C8H18   | 000584-94-1 | 47   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

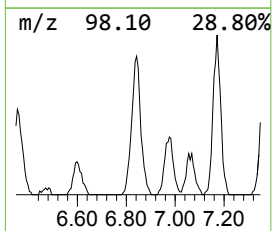
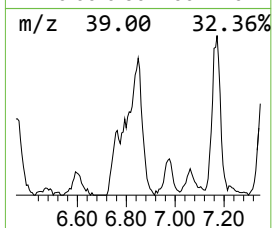
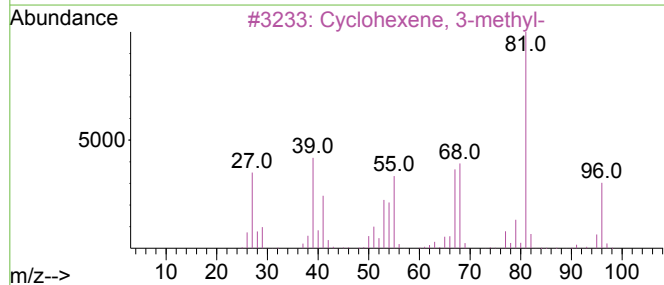
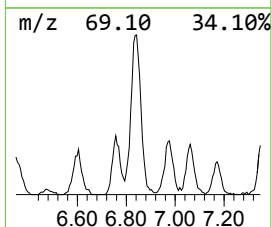
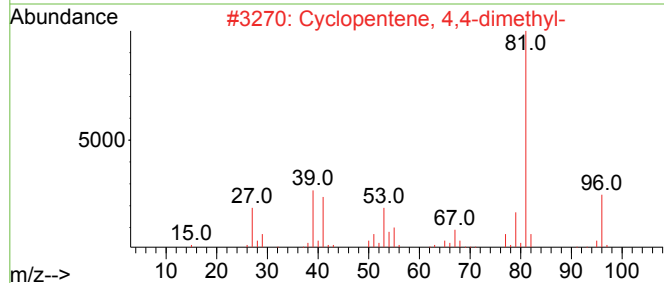
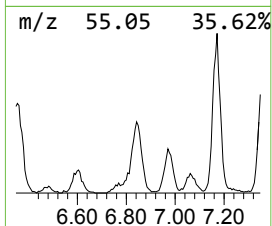
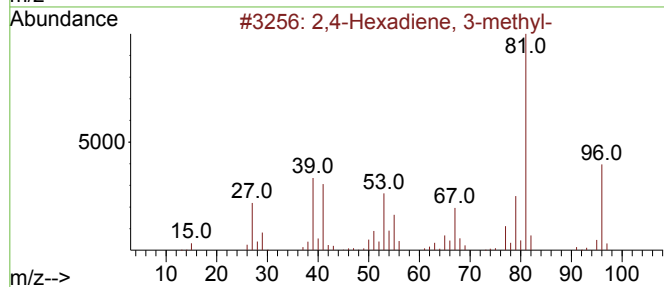
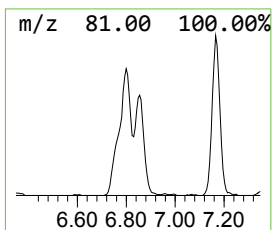
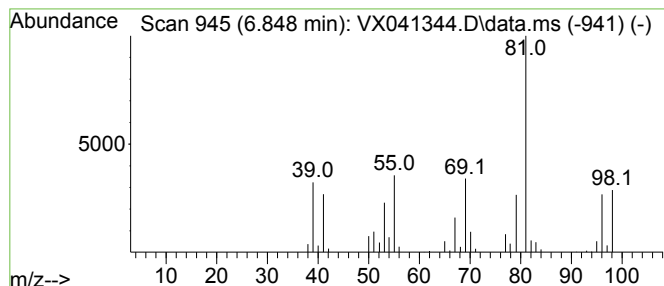
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 2,4-Hexadiene, 3-methyl- Concentration Rank 14

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 6.848 | 10.84 ug/l | 140918 | 1,4-Difluorobenzene | 6.763 |

| Hit# | of | 5 | Tentative ID                | MW | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|----|---------|-------------|------|
| 1    |    |   | 2,4-Hexadiene, 3-methyl-    | 96 | C7H12   | 028823-42-9 | 74   |
| 2    |    |   | Cyclopentene, 4,4-dimethyl- | 96 | C7H12   | 019037-72-0 | 64   |
| 3    |    |   | Cyclohexene, 3-methyl-      | 96 | C7H12   | 000591-48-0 | 64   |
| 4    |    |   | Furan, 2-ethyl-             | 96 | C6H8O   | 003208-16-0 | 58   |
| 5    |    |   | 2,4-Heptadiene, (E,E)-      | 96 | C7H12   | 002384-94-3 | 58   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

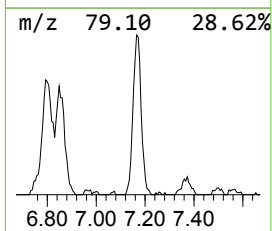
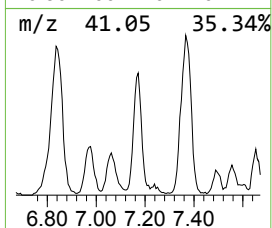
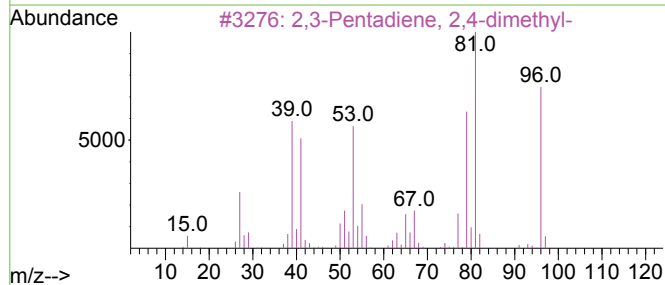
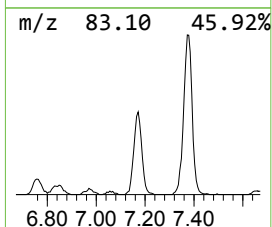
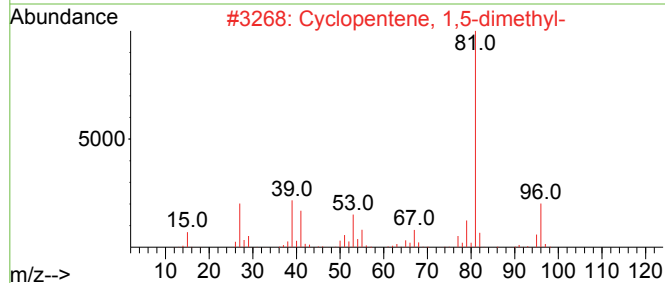
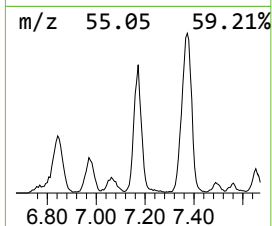
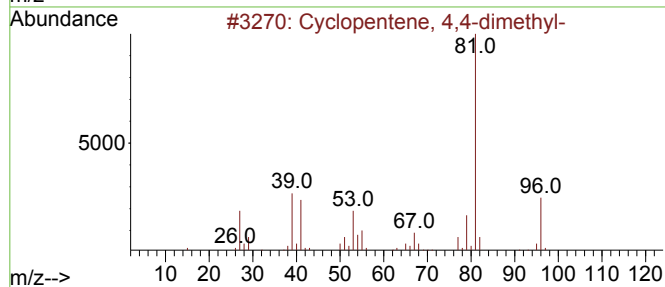
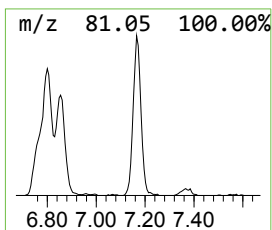
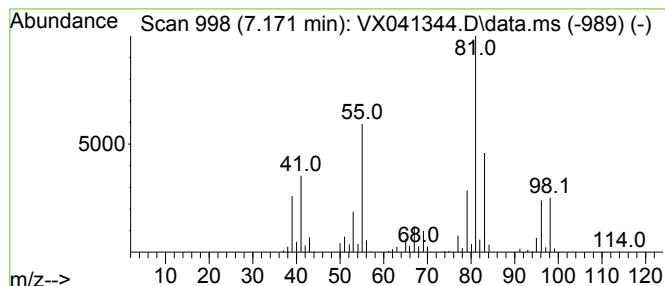
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Cyclopentene, 4,4-dimethyl- Concentration Rank 12

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 7.171 | 12.86 ug/l | 167106 | 1,4-Difluorobenzene | 6.763 |

| Hit# | of | 5 | Tentative ID                       | MW | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|----|---------|-------------|------|
| 1    |    |   | Cyclopentene, 4,4-dimethyl-        | 96 | C7H12   | 019037-72-0 | 60   |
| 2    |    |   | Cyclopentene, 1,5-dimethyl-        | 96 | C7H12   | 016491-15-9 | 60   |
| 3    |    |   | 2,3-Pentadiene, 2,4-dimethyl-      | 96 | C7H12   | 001000-87-9 | 53   |
| 4    |    |   | Furan, 2-ethyl-                    | 96 | C6H8O   | 003208-16-0 | 50   |
| 5    |    |   | Cyclobutane, (1-methylethylidene)- | 96 | C7H12   | 001528-22-9 | 49   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

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

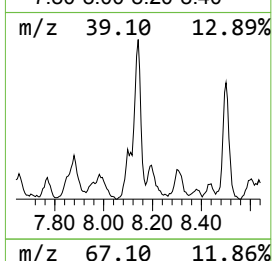
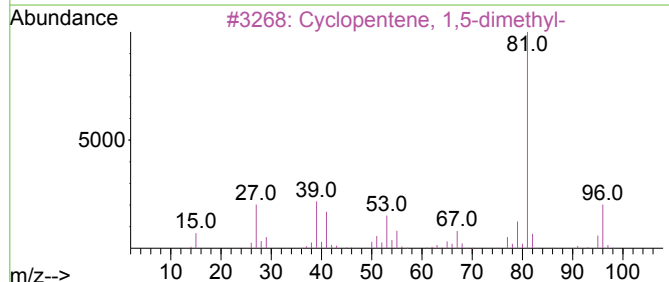
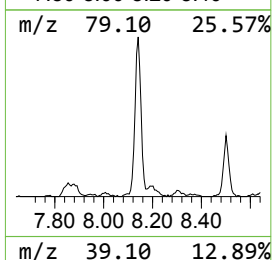
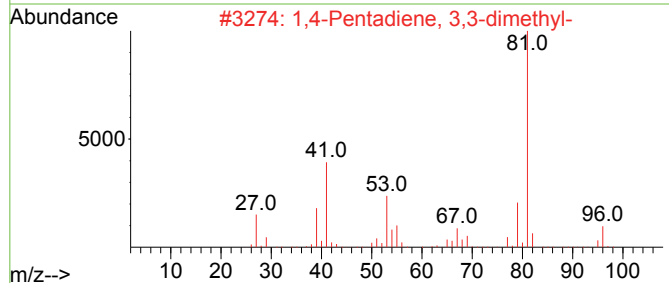
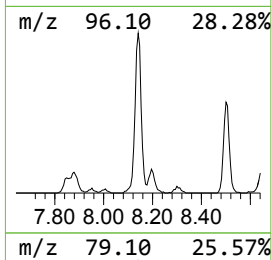
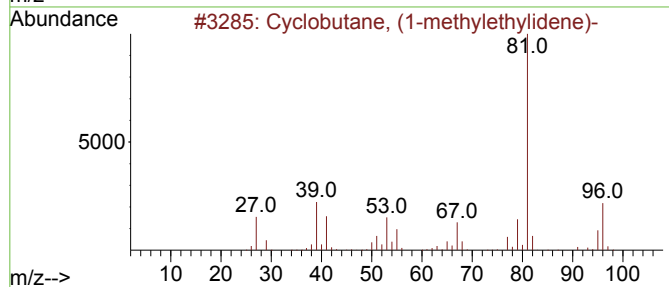
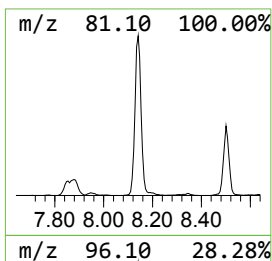
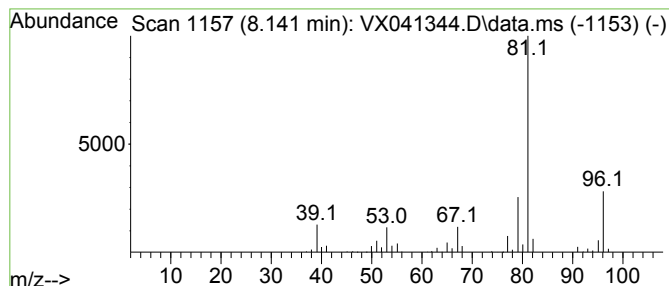
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Peak Number 7 Cyclobutane, (1-methylethyl... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 8.141 | 16.79 ug/l | 218158 | 1,4-Difluorobenzene | 6.763 |

| Hit# | of 5 | Tentative ID                        | MW | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|----|---------|-------------|------|
| 1    |      | Cyclobutane, (1-methylethylidene)-  | 96 | C7H12   | 001528-22-9 | 86   |
| 2    |      | 1,4-Pentadiene, 3,3-dimethyl-       | 96 | C7H12   | 001112-35-2 | 78   |
| 3    |      | Cyclopentene, 1,5-dimethyl-         | 96 | C7H12   | 016491-15-9 | 78   |
| 4    |      | Bicyclo[2.1.0]pentane, 1,4-dimet... | 96 | C7H12   | 017065-18-8 | 72   |
| 5    |      | Cyclohexene, 4-methyl-              | 96 | C7H12   | 000591-47-9 | 72   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

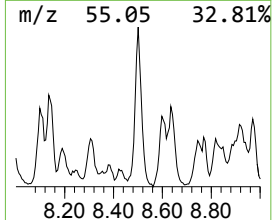
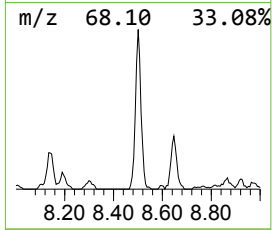
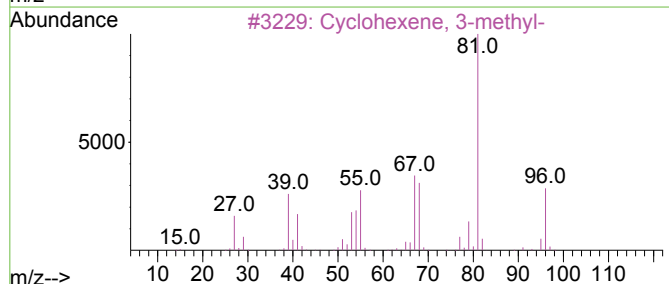
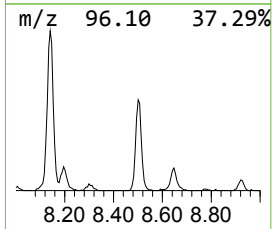
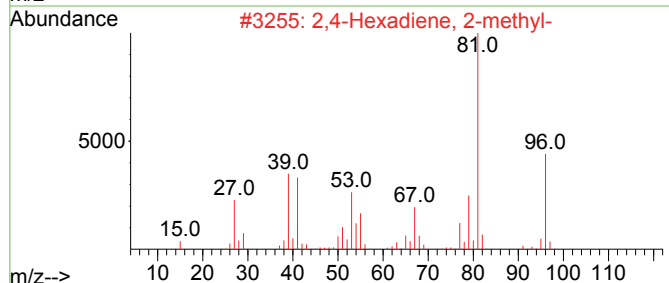
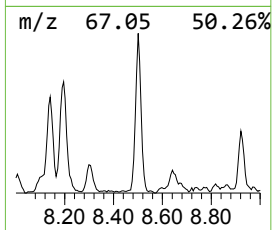
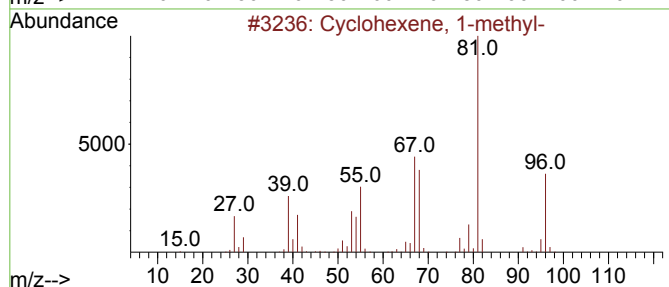
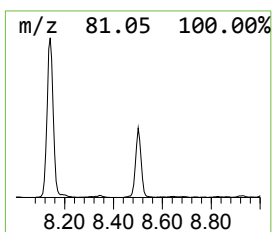
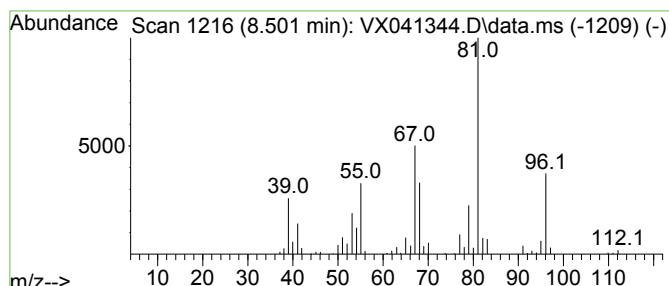
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Cyclohexene, 1-methyl- Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.   |
|-------|------------|--------|------------------|--------|
| 8.501 | 10.51 ug/l | 152880 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID             | MW | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|----|---------|-------------|------|
| 1    |    |   | Cyclohexene, 1-methyl-   | 96 | C7H12   | 000591-49-1 | 94   |
| 2    |    |   | 2,4-Hexadiene, 2-methyl- | 96 | C7H12   | 028823-41-8 | 93   |
| 3    |    |   | Cyclohexene, 3-methyl-   | 96 | C7H12   | 000591-48-0 | 90   |
| 4    |    |   | Cyclohexene, 4-methyl-   | 96 | C7H12   | 000591-47-9 | 90   |
| 5    |    |   | Cyclohexane, methylene-  | 96 | C7H12   | 001192-37-6 | 74   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

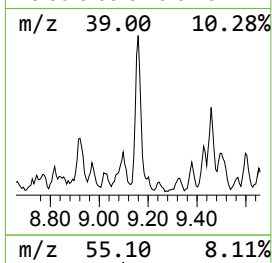
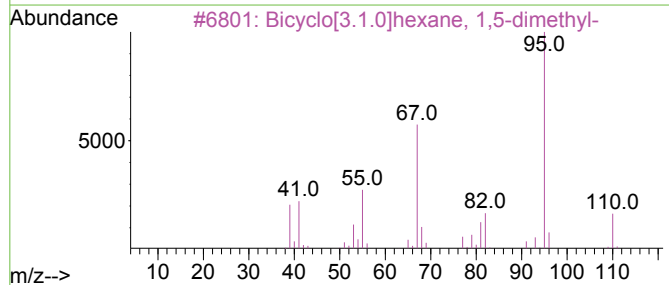
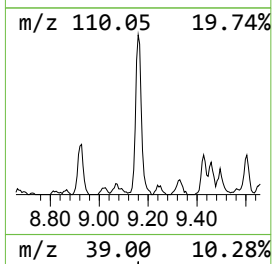
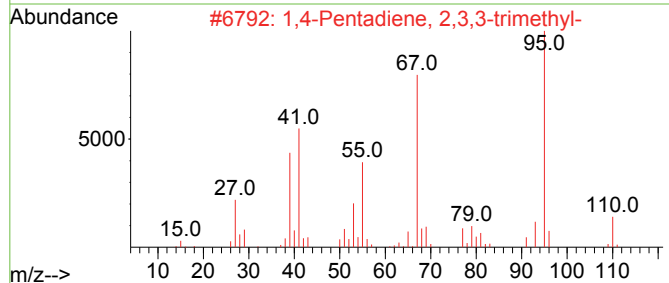
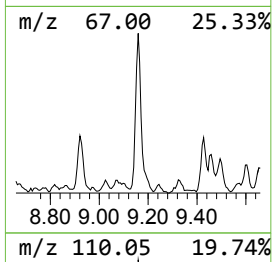
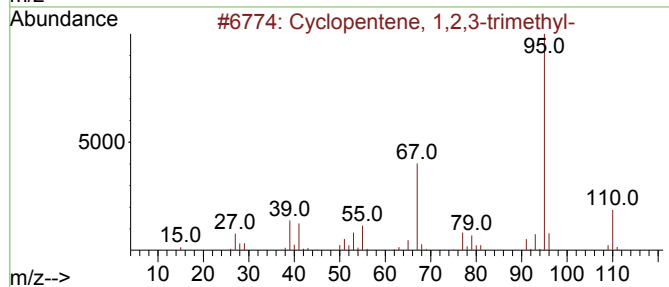
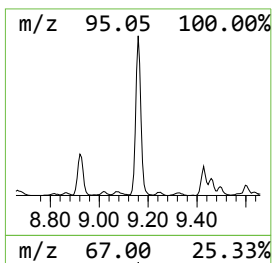
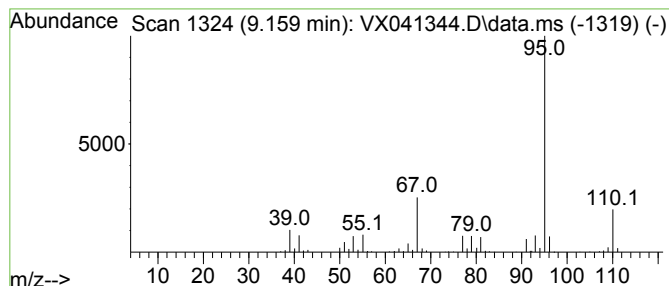
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Cyclopentene, 1,2,3-trimethyl- Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.   |
|-------|------------|--------|------------------|--------|
| 9.159 | 11.79 ug/l | 171471 | Chlorobenzene-d5 | 10.055 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Cyclopentene, 1,2,3-trimethyl-      | 110 | C8H14   | 000473-91-6  | 91   |
| 2    |    |   | 1,4-Pentadiene, 2,3,3-trimethyl-    | 110 | C8H14   | 000756-02-5  | 86   |
| 3    |    |   | Bicyclo[3.1.0]hexane, 1,5-dimethyl- | 110 | C8H14   | 1010142-17-5 | 83   |
| 4    |    |   | Cyclohexene, 3,5-dimethyl-          | 110 | C8H14   | 000823-17-6  | 72   |
| 5    |    |   | 5,5-Dimethyl-1,3-hexadiene          | 110 | C8H14   | 001515-79-3  | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

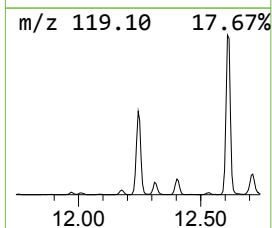
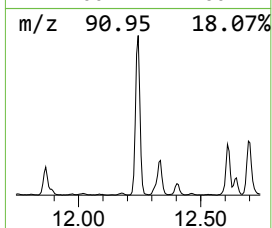
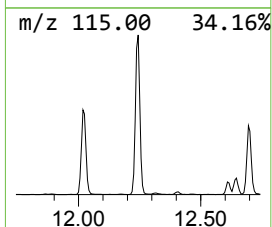
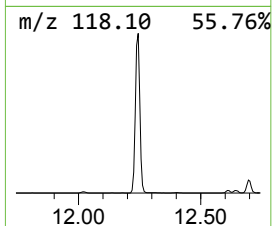
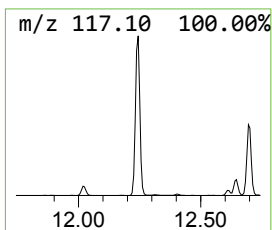
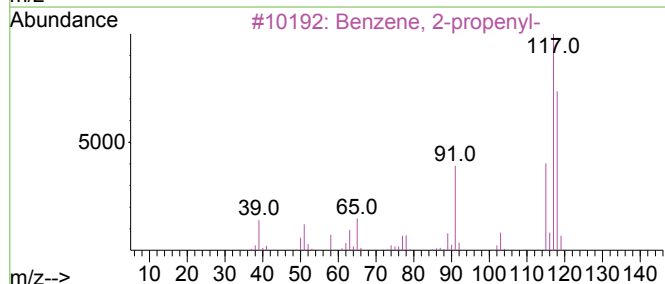
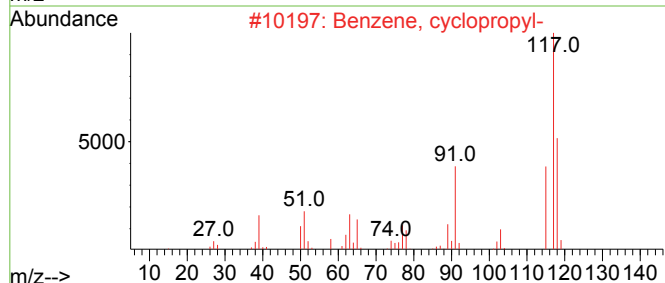
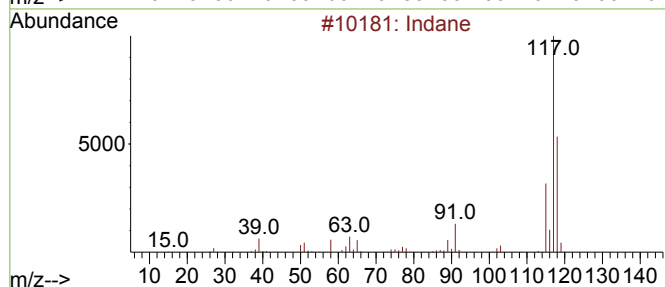
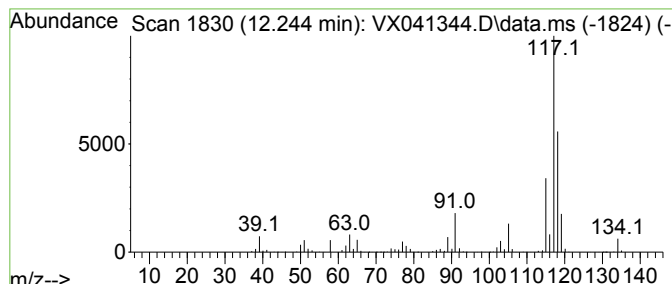
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Quant Title : SW846 8260

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

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Peak Number 10 Indane Concentration Rank 1

| R.T.      | EstConc                      | Area    | Relative to ISTD       | R.T.        |      |
|-----------|------------------------------|---------|------------------------|-------------|------|
| 12.244    | 87.50 ug/l                   | 1232360 | 1,4-Dichlorobenzene-d4 | 12.024      |      |
| Hit# of 5 | Tentative ID                 | MW      | MolForm                | CAS#        | Qual |
| 1         | Indane                       | 118     | C9H10                  | 000496-11-7 | 93   |
| 2         | Benzene, cyclopropyl-        | 118     | C9H10                  | 000873-49-4 | 76   |
| 3         | Benzene, 2-propenyl-         | 118     | C9H10                  | 000300-57-2 | 76   |
| 4         | Delatacyclene                | 118     | C9H10                  | 007785-10-6 | 58   |
| 5         | Benzene, 1-ethenyl-2-methyl- | 118     | C9H10                  | 000611-15-4 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

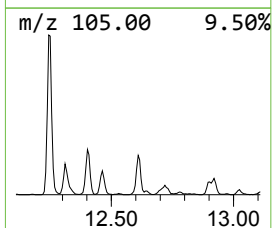
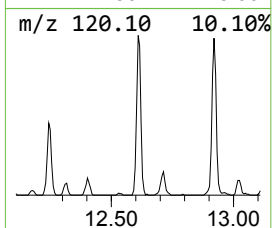
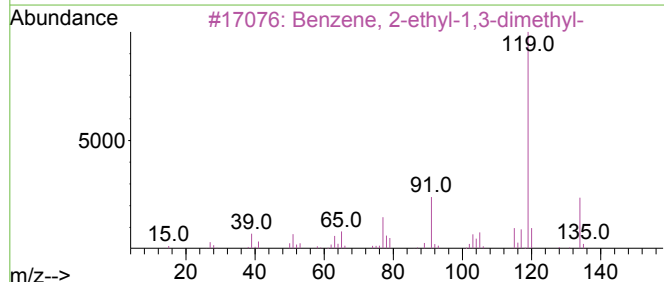
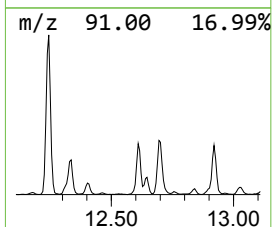
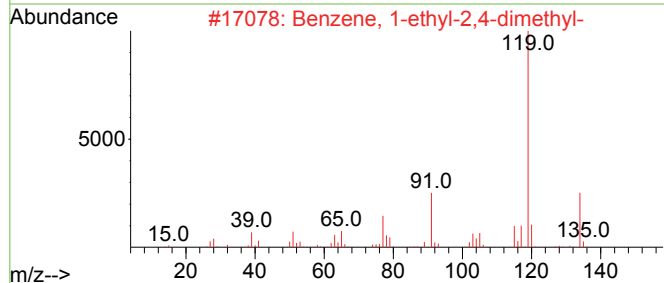
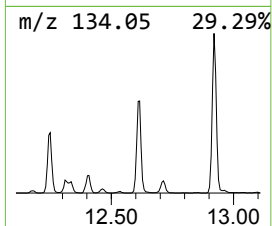
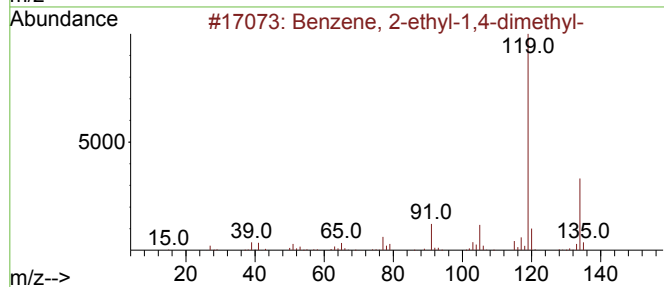
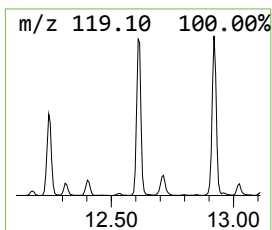
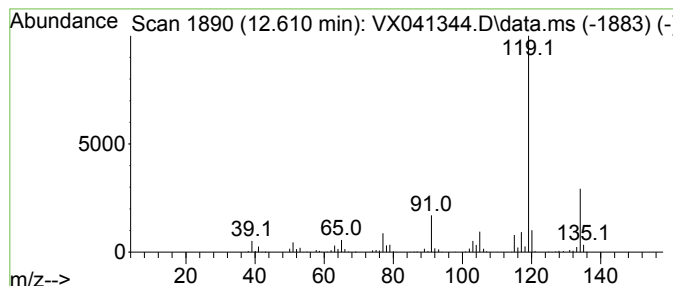
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 6

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 12.610 | 22.07 ug/l | 310847 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 95   |
| 2    |    |   | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 95   |
| 3    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 95   |
| 4    |    |   | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 95   |
| 5    |    |   | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 95   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

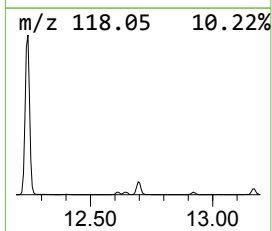
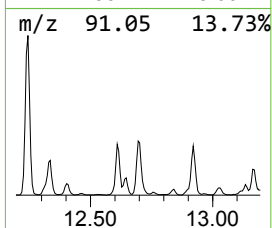
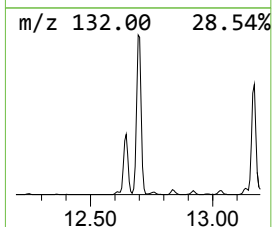
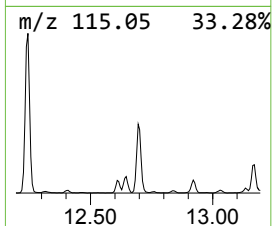
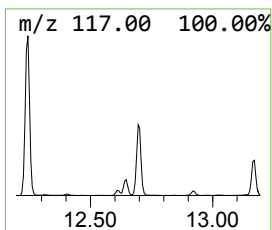
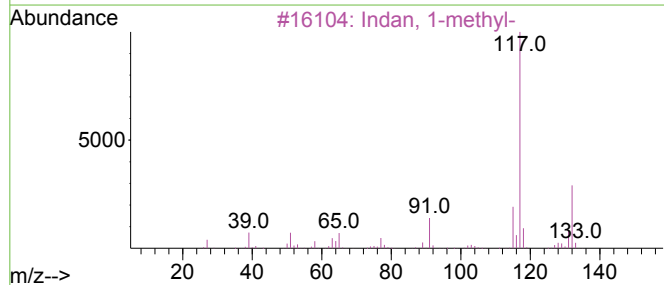
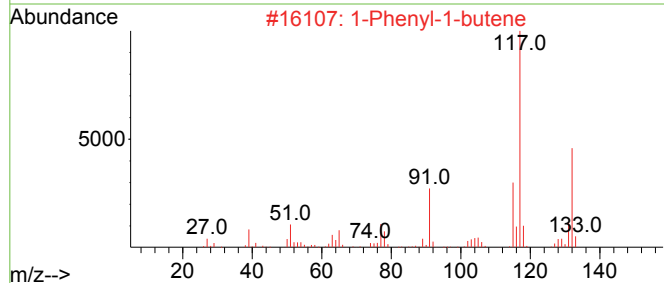
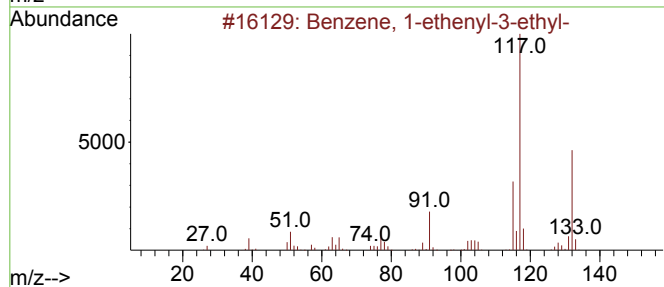
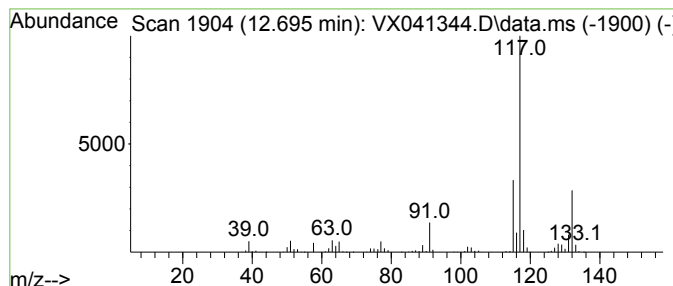
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Quant Title : SW846 8260

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

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Peak Number 12 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 3

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 12.695 | 34.52 ug/l | 486215 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethenyl-3-ethyl-         | 132 | C10H12  | 007525-62-4 | 91   |
| 2    |    |   | 1-Phenyl-1-butene                   | 132 | C10H12  | 000824-90-8 | 90   |
| 3    |    |   | Indan, 1-methyl-                    | 132 | C10H12  | 000767-58-8 | 90   |
| 4    |    |   | 1-Methyl-2-phenylcyclopropane       | 132 | C10H12  | 003145-76-4 | 87   |
| 5    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 87   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

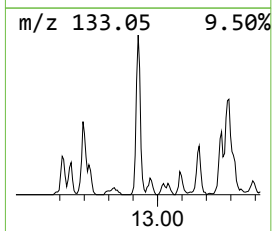
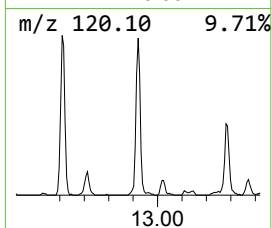
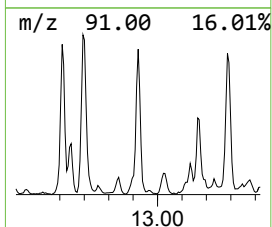
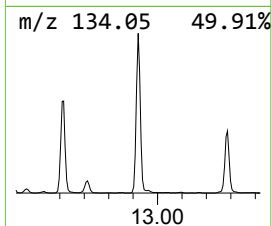
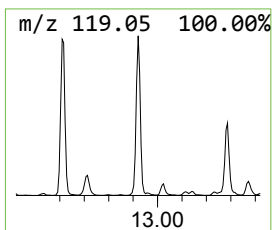
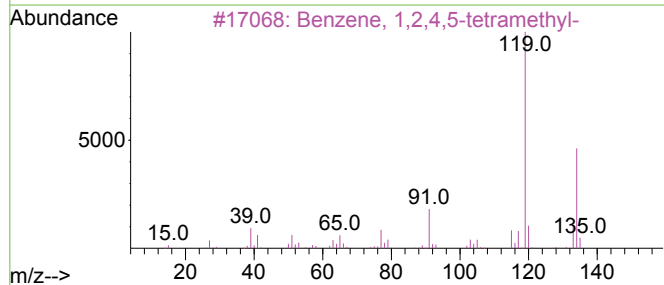
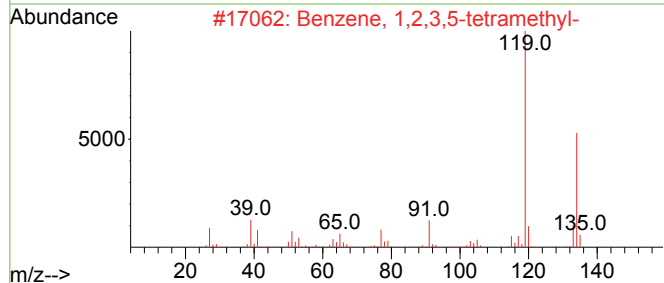
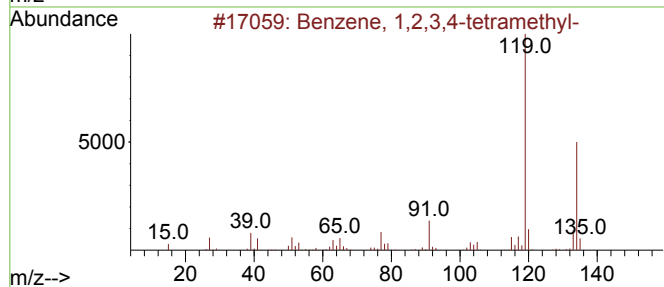
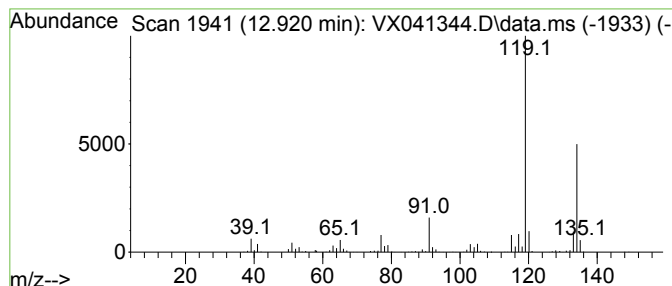
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Quant Title : SW846 8260

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

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Peak Number 13 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 4

| R.T.      | EstConc                        | Area   | Relative to ISTD       | R.T.        |      |
|-----------|--------------------------------|--------|------------------------|-------------|------|
| 12.921    | 24.72 ug/l                     | 348205 | 1,4-Dichlorobenzene-d4 | 12.024      |      |
| Hit# of 5 | Tentative ID                   | MW     | MolForm                | CAS#        | Qual |
| 1         | Benzene, 1,2,3,4-tetramethyl-  | 134    | C10H14                 | 000488-23-3 | 97   |
| 2         | Benzene, 1,2,3,5-tetramethyl-  | 134    | C10H14                 | 000527-53-7 | 97   |
| 3         | Benzene, 1,2,4,5-tetramethyl-  | 134    | C10H14                 | 000095-93-2 | 96   |
| 4         | Benzene, 2-ethyl-1,3-dimethyl- | 134    | C10H14                 | 002870-04-4 | 94   |
| 5         | Benzene, 1-ethyl-2,4-dimethyl- | 134    | C10H14                 | 000874-41-9 | 93   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

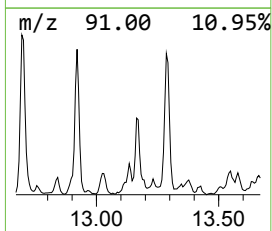
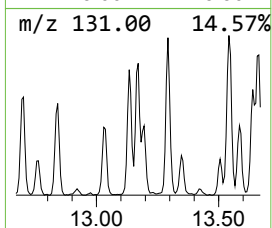
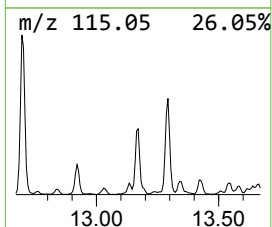
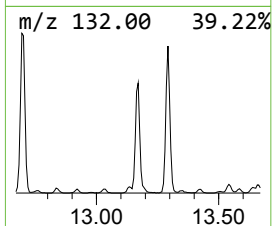
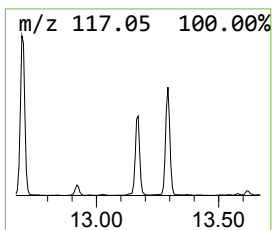
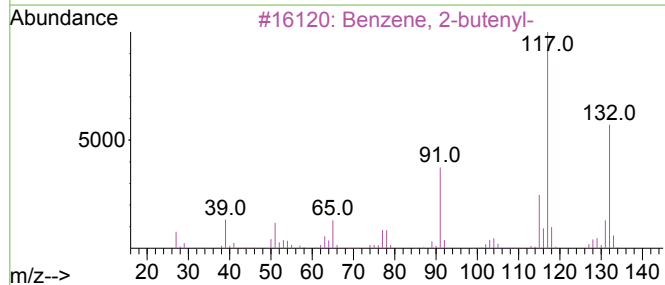
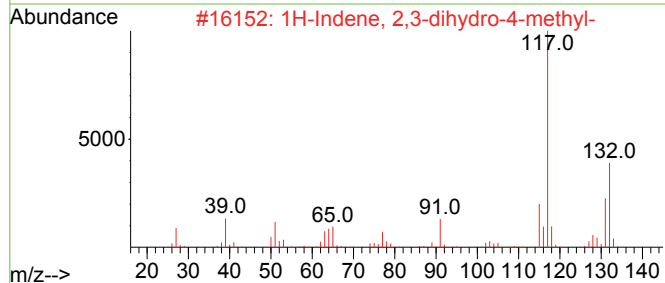
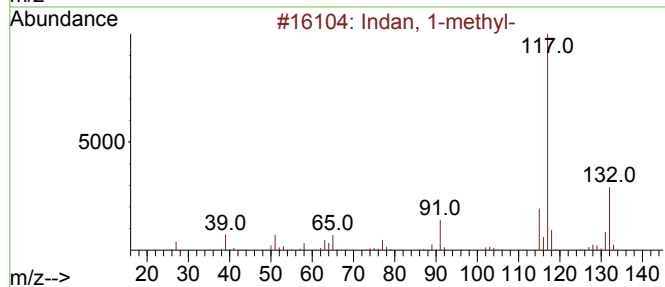
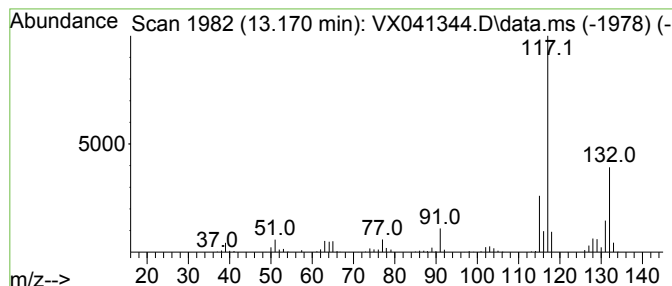
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Quant Title : SW846 8260

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

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Peak Number 14 Indan, 1-methyl- Concentration Rank 9

| R.T.   | EstConc    | Area   | Relative to ISTD       | R.T.   |
|--------|------------|--------|------------------------|--------|
| 13.170 | 19.50 ug/l | 274640 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Indan, 1-methyl-                  | 132 | C10H12  | 000767-58-8 | 94   |
| 2    |    |   | 1H-Indene, 2,3-dihydro-4-methyl-  | 132 | C10H12  | 000824-22-6 | 90   |
| 3    |    |   | Benzene, 2-butenyl-               | 132 | C10H12  | 001560-06-1 | 87   |
| 4    |    |   | Benzene, 1-ethenyl-4-ethyl-       | 132 | C10H12  | 003454-07-7 | 87   |
| 5    |    |   | Benzene, 1-methyl-2-(2-propenyl)- | 132 | C10H12  | 001587-04-8 | 87   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

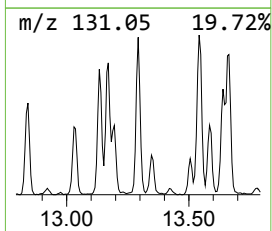
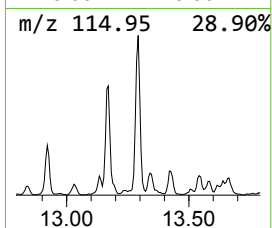
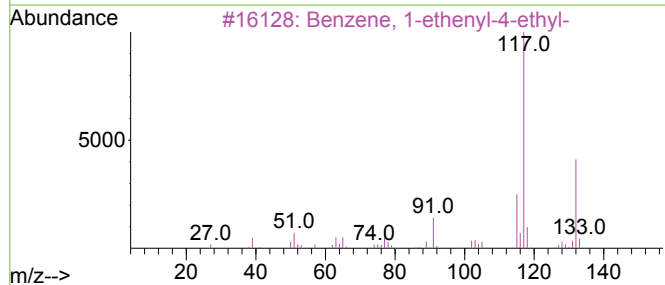
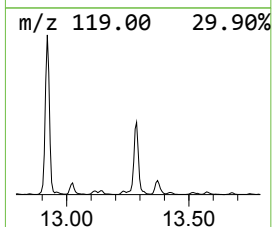
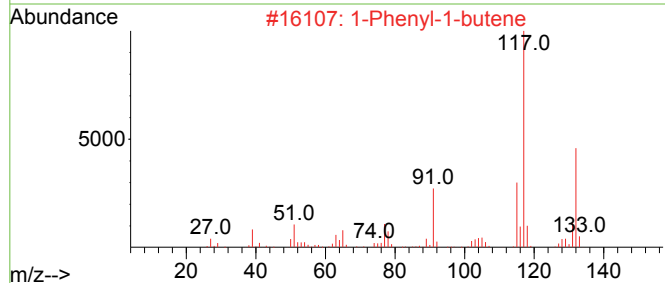
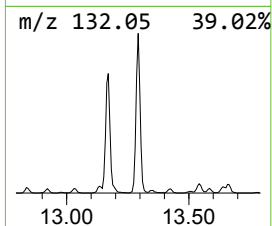
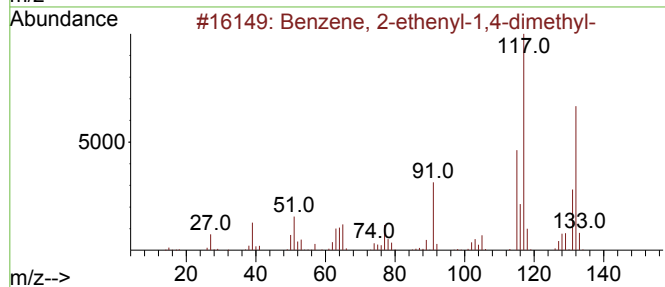
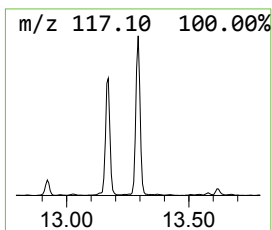
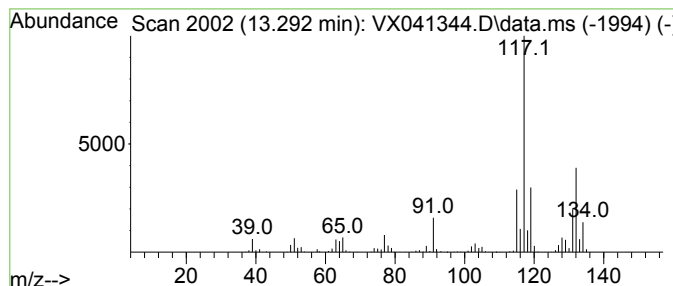
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Quant Title : SW846 8260

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

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Peak Number 15 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

| R.T.      | EstConc                           | Area   | Relative to ISTD       |         |             | R.T.   |
|-----------|-----------------------------------|--------|------------------------|---------|-------------|--------|
| 13.292    | 34.64 ug/l                        | 487835 | 1,4-Dichlorobenzene-d4 |         |             | 12.024 |
| Hit# of 5 | Tentative ID                      |        | MW                     | MolForm | CAS#        | Qual   |
| 1         | Benzene, 2-ethenyl-1,4-dimethyl-  |        | 132                    | C10H12  | 002039-89-6 | 91     |
| 2         | 1-Phenyl-1-butene                 |        | 132                    | C10H12  | 000824-90-8 | 90     |
| 3         | Benzene, 1-ethenyl-4-ethyl-       |        | 132                    | C10H12  | 003454-07-7 | 89     |
| 4         | 3-Phenylbut-1-ene                 |        | 132                    | C10H12  | 000934-10-1 | 87     |
| 5         | Benzene, 1-methyl-2-(2-propenyl)- |        | 132                    | C10H12  | 001587-04-8 | 81     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_X\Data\VX050224\  
Data File : VX041344.D  
Acq On : 02 May 2024 11:32  
Operator : JC/MD  
Sample : P2264-14  
Misc : 5.0mL/MSVOA\_X/WATER  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
MSVOA\_X  
ClientSampleId :  
T2-MW-3-8.0-042424

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_X\Method\82X043024W.M  
Quant Title : SW846 8260

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |        |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|--------|------|
|                    |        |         |       |          | #                     | RT     | Resp   | Conc |
| Pentane, 2-methyl- | 2.819  | 20.8    | ug/l  | 147031   | 1                     | 5.550  | 353680 | 50.0 |
| Pentane, 3-methyl- | 3.093  | 20.2    | ug/l  | 143086   | 1                     | 5.550  | 353680 | 50.0 |
| Cyclopentane, m... | 4.294  | 23.9    | ug/l  | 169036   | 1                     | 5.550  | 353680 | 50.0 |
| Hexane, 3-methyl-  | 5.818  | 13.4    | ug/l  | 94686    | 1                     | 5.550  | 353680 | 50.0 |
| 2,4-Hexadiene, ... | 6.848  | 10.8    | ug/l  | 140918   | 2                     | 6.763  | 649845 | 50.0 |
| Cyclopentene, 4... | 7.171  | 12.9    | ug/l  | 167106   | 2                     | 6.763  | 649845 | 50.0 |
| Cyclobutane, (1... | 8.141  | 16.8    | ug/l  | 218158   | 2                     | 6.763  | 649845 | 50.0 |
| Cyclohexene, 1-... | 8.501  | 10.5    | ug/l  | 152880   | 3                     | 10.055 | 727063 | 50.0 |
| Cyclopentene, 1... | 9.159  | 11.8    | ug/l  | 171471   | 3                     | 10.055 | 727063 | 50.0 |
| Indane             | 12.244 | 87.5    | ug/l  | 1232360  | 4                     | 12.024 | 704201 | 50.0 |
| Benzene, 2-ethy... | 12.610 | 22.1    | ug/l  | 310847   | 4                     | 12.024 | 704201 | 50.0 |
| Benzene, 1-ethe... | 12.695 | 34.5    | ug/l  | 486215   | 4                     | 12.024 | 704201 | 50.0 |
| Benzene, 1,2,3,... | 12.921 | 24.7    | ug/l  | 348205   | 4                     | 12.024 | 704201 | 50.0 |
| Indan, 1-methyl-   | 13.170 | 19.5    | ug/l  | 274640   | 4                     | 12.024 | 704201 | 50.0 |
| Benzene, 2-ethe... | 13.292 | 34.6    | ug/l  | 487835   | 4                     | 12.024 | 704201 | 50.0 |