

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD041619\  
 Data File : VD061815.D  
 Acq On : 16 Apr 2019 17:03  
 Operator : VA/AP  
 Sample : K2358-10  
 Misc : 4.88g/5ml/MSVOA D/SOIL  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 MSVOA\_D  
 ClientSampleId :  
 TP-7-S2

Integration Parameters: RTEINT.P

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

Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\VOASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
 Title : SW846 8260

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 1.569       | 12            | 15          | 28           | rVB2     | 86610          | 256849        | 18.13%          | 2.137%        |
| 2         | 1.854       | 40            | 44          | 51           | rBV      | 36332          | 67519         | 4.77%           | 0.562%        |
| 3         | 2.651       | 121           | 125         | 129          | rVB2     | 6242           | 15255         | 1.08%           | 0.127%        |
| 4         | 3.232       | 178           | 184         | 194          | rBV      | 28113          | 77022         | 5.44%           | 0.641%        |
| 5         | 3.400       | 197           | 201         | 206          | rBV      | 27237          | 59858         | 4.22%           | 0.498%        |
| 6         | 3.833       | 240           | 245         | 259          | rBV      | 101335         | 316312        | 22.32%          | 2.632%        |
| 7         | 4.246       | 283           | 287         | 293          | rBV4     | 8348           | 25392         | 1.79%           | 0.211%        |
| 8         | 4.719       | 330           | 335         | 342          | rVB3     | 11308          | 36105         | 2.55%           | 0.300%        |
| 9         | 6.077       | 466           | 473         | 478          | rBV4     | 8764           | 31175         | 2.20%           | 0.259%        |
| 10        | 6.944       | 554           | 561         | 567          | rBV      | 147590         | 422765        | 29.84%          | 3.517%        |
| 11        | 7.062       | 567           | 573         | 588          | rVB      | 407403         | 1170688       | 82.62%          | 9.740%        |
| 12        | 7.465       | 609           | 614         | 631          | rVB      | 193509         | 538551        | 38.01%          | 4.481%        |
| 13        | 8.233       | 686           | 692         | 706          | rBV      | 562883         | 1416909       | 100.00%         | 11.788%       |
| 14        | 10.419      | 909           | 914         | 921          | rBV      | 579427         | 1379594       | 97.37%          | 11.478%       |
| 15        | 10.517      | 921           | 924         | 930          | rVB      | 88563          | 182434        | 12.88%          | 1.518%        |
| 16        | 11.265      | 994           | 1000        | 1004         | rBB2     | 7107           | 16619         | 1.17%           | 0.138%        |
| 17        | 12.378      | 1109          | 1113        | 1120         | rVB      | 648409         | 1136811       | 80.23%          | 9.458%        |
| 18        | 12.525      | 1125          | 1128        | 1133         | rVB      | 9110           | 17516         | 1.24%           | 0.146%        |
| 19        | 12.703      | 1141          | 1146        | 1150         | rVB3     | 6965           | 16781         | 1.18%           | 0.140%        |
| 20        | 13.008      | 1174          | 1177        | 1179         | rBV3     | 10700          | 21180         | 1.49%           | 0.176%        |
| 21        | 13.057      | 1179          | 1182        | 1185         | rVB      | 21493          | 30282         | 2.14%           | 0.252%        |
| 22        | 13.224      | 1192          | 1199        | 1203         | rBV3     | 12282          | 29687         | 2.10%           | 0.247%        |
| 23        | 13.333      | 1203          | 1210        | 1215         | rVV7     | 18097          | 55934         | 3.95%           | 0.465%        |
| 24        | 13.480      | 1220          | 1225        | 1230         | rBV3     | 7680           | 16523         | 1.17%           | 0.137%        |
| 25        | 13.559      | 1230          | 1233        | 1237         | rBV      | 451252         | 634582        | 44.79%          | 5.280%        |
| 26        | 13.717      | 1246          | 1249        | 1254         | rVB2     | 53284          | 72715         | 5.13%           | 0.605%        |
| 27        | 13.825      | 1257          | 1260        | 1261         | rBV2     | 11319          | 18297         | 1.29%           | 0.152%        |
| 28        | 13.894      | 1264          | 1267        | 1269         | rBV3     | 20222          | 35937         | 2.54%           | 0.299%        |
| 29        | 13.933      | 1269          | 1271        | 1274         | rVV      | 60761          | 88034         | 6.21%           | 0.732%        |
| 30        | 13.972      | 1274          | 1275        | 1278         | rVB2     | 21262          | 23068         | 1.63%           | 0.192%        |
| 31        | 14.032      | 1278          | 1281        | 1283         | rBV3     | 11553          | 16625         | 1.17%           | 0.138%        |
| 32        | 14.091      | 1283          | 1287        | 1289         | rBV      | 38406          | 62588         | 4.42%           | 0.521%        |
| 33        | 14.209      | 1296          | 1299        | 1301         | rBV2     | 9022           | 19811         | 1.40%           | 0.165%        |
| 34        | 14.248      | 1301          | 1303        | 1310         | rVB5     | 21423          | 58386         | 4.12%           | 0.486%        |

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## Integration Parameters: RTEINT.P

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 Sampling : 1  
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Method : Z:\VOASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
 Title : SW846 8260

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 14.366 | 1310 | 1315 | 1320 | rBV6 | 18546  | 68975  | 4.87%  | 0.574% |
| 36 | 14.445 | 1320 | 1323 | 1325 | rVV  | 35190  | 54088  | 3.82%  | 0.450% |
| 37 | 14.524 | 1328 | 1331 | 1333 | rVV  | 424155 | 549221 | 38.76% | 4.569% |
| 38 | 14.563 | 1333 | 1335 | 1338 | rVV  | 193395 | 232513 | 16.41% | 1.934% |
| 39 | 14.652 | 1340 | 1344 | 1346 | rVV5 | 20929  | 50161  | 3.54%  | 0.417% |
| 40 | 14.711 | 1346 | 1350 | 1352 | rVV3 | 35659  | 60393  | 4.26%  | 0.502% |
| 41 | 14.740 | 1352 | 1353 | 1355 | rVV  | 21586  | 17274  | 1.22%  | 0.144% |
| 42 | 14.780 | 1355 | 1357 | 1359 | rVV  | 124310 | 169420 | 11.96% | 1.410% |
| 43 | 14.809 | 1359 | 1360 | 1363 | rVV3 | 73605  | 99371  | 7.01%  | 0.827% |
| 44 | 14.918 | 1367 | 1371 | 1373 | rVV  | 47516  | 91371  | 6.45%  | 0.760% |
| 45 | 14.977 | 1376 | 1377 | 1378 | rVV  | 17077  | 18714  | 1.32%  | 0.156% |
| 46 | 15.006 | 1378 | 1380 | 1383 | rVV  | 49263  | 77836  | 5.49%  | 0.648% |
| 47 | 15.055 | 1383 | 1385 | 1388 | rVV  | 78083  | 108732 | 7.67%  | 0.905% |
| 48 | 15.095 | 1388 | 1389 | 1391 | rVV2 | 21842  | 25542  | 1.80%  | 0.213% |
| 49 | 15.144 | 1391 | 1394 | 1398 | rVV3 | 104285 | 177966 | 12.56% | 1.481% |
| 50 | 15.223 | 1398 | 1402 | 1405 | rVV3 | 36937  | 87084  | 6.15%  | 0.725% |
| 51 | 15.272 | 1405 | 1407 | 1411 | rVV3 | 105582 | 174924 | 12.35% | 1.455% |
| 52 | 15.351 | 1413 | 1415 | 1416 | rVV  | 75344  | 93100  | 6.57%  | 0.775% |
| 53 | 15.380 | 1416 | 1418 | 1421 | rVV  | 158545 | 186965 | 13.20% | 1.556% |
| 54 | 15.429 | 1421 | 1423 | 1426 | rVV2 | 83507  | 128692 | 9.08%  | 1.071% |
| 55 | 15.469 | 1426 | 1427 | 1431 | rVV2 | 37838  | 77345  | 5.46%  | 0.643% |
| 56 | 15.538 | 1431 | 1434 | 1437 | rVV2 | 30887  | 66832  | 4.72%  | 0.556% |
| 57 | 15.587 | 1437 | 1439 | 1442 | rVV2 | 35364  | 55949  | 3.95%  | 0.465% |
| 58 | 15.636 | 1442 | 1444 | 1445 | rVV  | 38315  | 48252  | 3.41%  | 0.401% |
| 59 | 15.676 | 1445 | 1448 | 1450 | rVV  | 269752 | 378312 | 26.70% | 3.147% |
| 60 | 15.744 | 1453 | 1455 | 1457 | rVV2 | 31314  | 42180  | 2.98%  | 0.351% |
| 61 | 15.804 | 1457 | 1461 | 1464 | rVV2 | 76454  | 139831 | 9.87%  | 1.163% |
| 62 | 15.853 | 1464 | 1466 | 1468 | rVV2 | 15118  | 26413  | 1.86%  | 0.220% |
| 63 | 15.931 | 1468 | 1474 | 1476 | rVV3 | 49759  | 129901 | 9.17%  | 1.081% |
| 64 | 16.000 | 1476 | 1481 | 1483 | rVV2 | 51000  | 99577  | 7.03%  | 0.828% |
| 65 | 16.050 | 1483 | 1486 | 1487 | rVV2 | 10634  | 19841  | 1.40%  | 0.165% |
| 66 | 16.109 | 1490 | 1492 | 1495 | rBV3 | 14440  | 23639  | 1.67%  | 0.197% |
| 67 | 16.178 | 1497 | 1499 | 1501 | rVB  | 26145  | 28203  | 1.99%  | 0.235% |
| 68 | 16.237 | 1501 | 1505 | 1507 | rBV2 | 26389  | 43050  | 3.04%  | 0.358% |

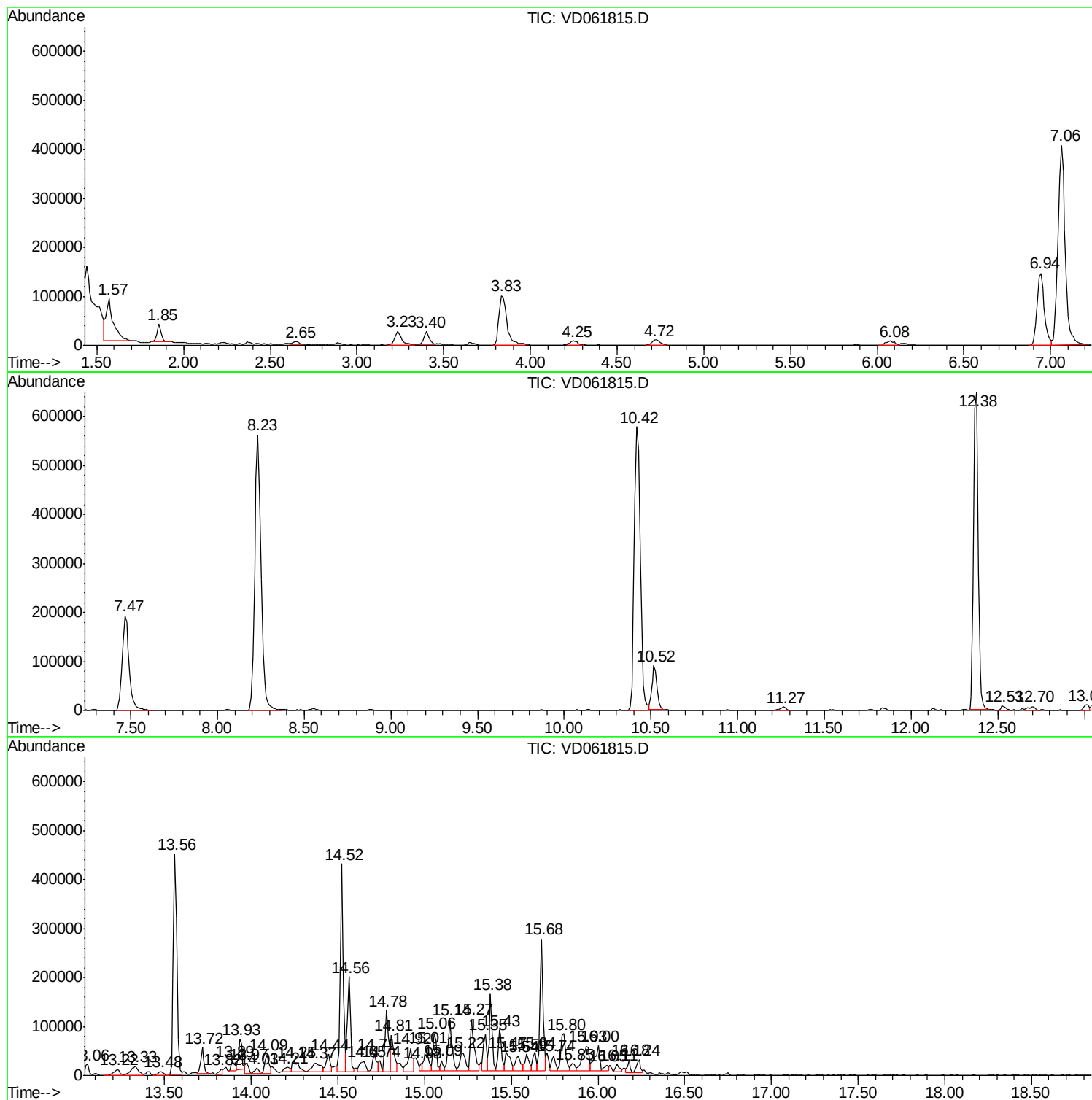
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ALS Vial : 14 Sample Multiplier: 1

Instrument :  
MSVOA\_D  
ClientSampled :  
TP-7-S2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041619\  
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ALS Vial : 14 Sample Multiplier: 1

Instrument :  
MSVOA\_D  
ClientSampleId :  
TP-7-S2

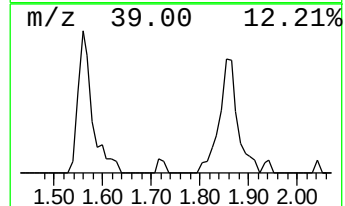
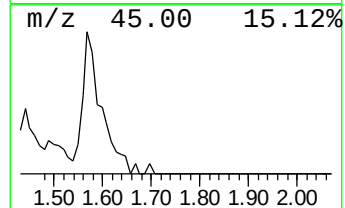
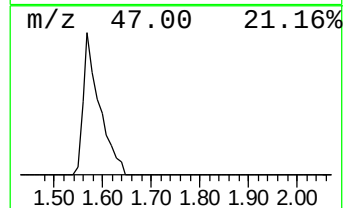
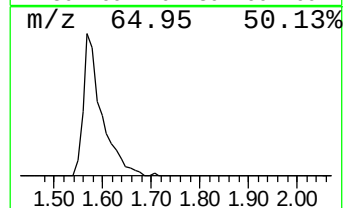
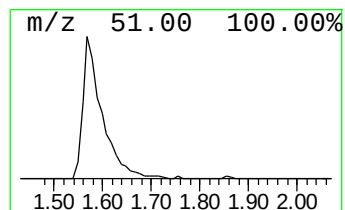
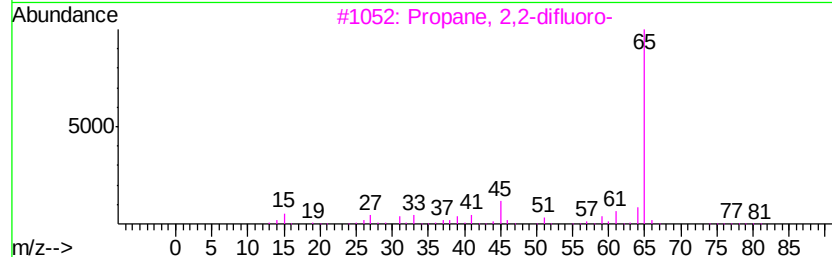
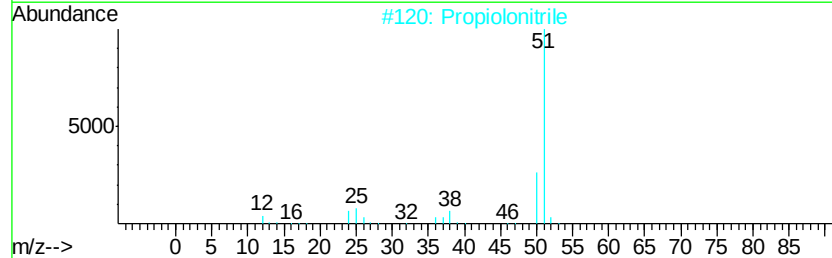
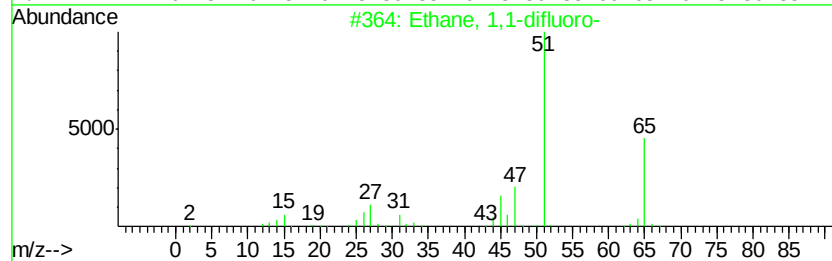
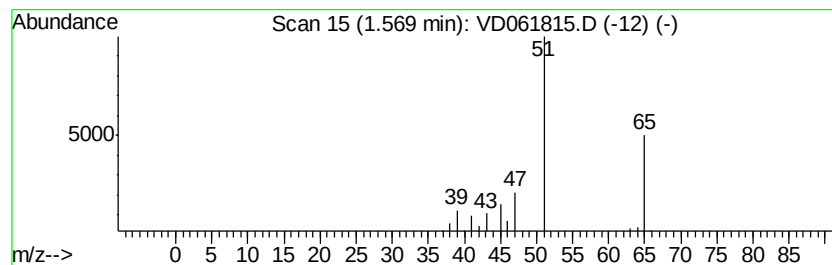
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
Quant Title : SW846 8260

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

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Peak Number 1 Ethane, 1,1-difluoro- Concentration Rank 9

| R.T. | EstConc    | Area   | Relative to ISTD   | R.T. |
|------|------------|--------|--------------------|------|
| 1.57 | 10.97 ug/l | 256849 | Pentafluorobenzene | 7.06 |

| Hit# | of | Tentative ID           | MW | MolForm | CAS#        | Qual |
|------|----|------------------------|----|---------|-------------|------|
| 1    | 5  | Ethane, 1,1-difluoro-  | 66 | C2H4F2  | 000075-37-6 | 91   |
| 2    |    | Propiolonitrile        | 51 | C3HN    | 001070-71-9 | 3    |
| 3    |    | Propane, 2,2-difluoro- | 80 | C3H6F2  | 000420-45-1 | 2    |
| 4    |    | Difluoromethane        | 52 | CH2F2   | 000075-10-5 | 1    |
| 5    |    | 2-Chloroethanol        | 80 | C2H5ClO | 000107-07-3 | 1    |



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ALS Vial : 14 Sample Multiplier: 1

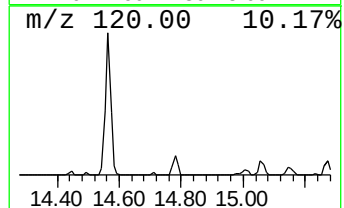
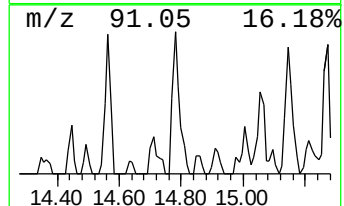
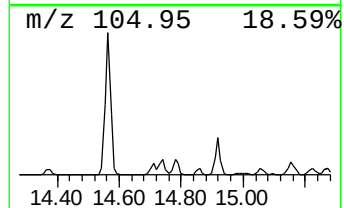
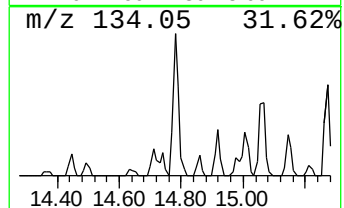
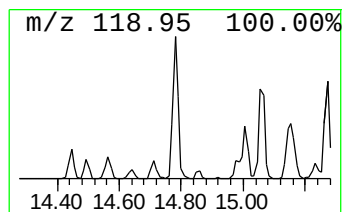
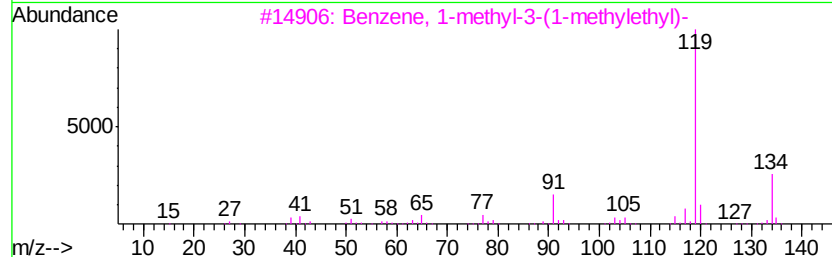
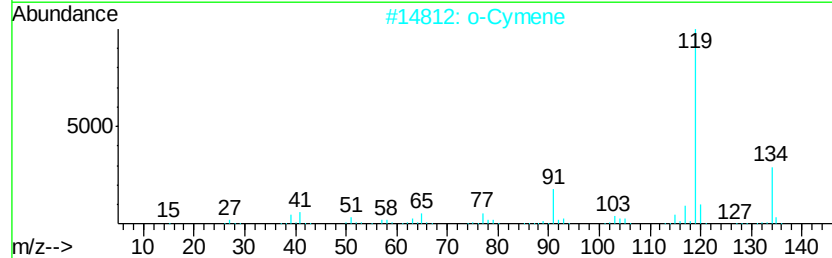
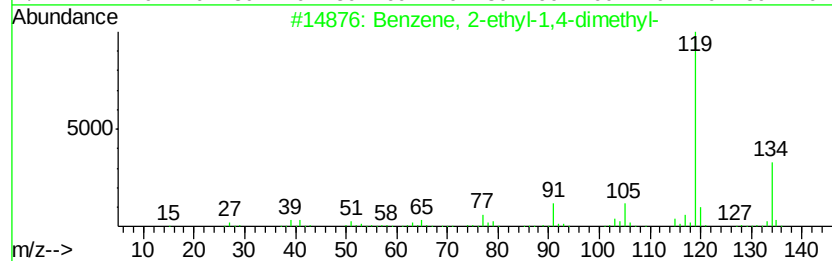
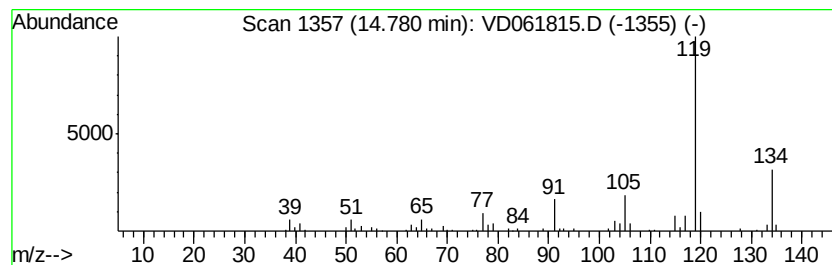
Instrument :  
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Quant Title : SW846 8260

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

\*\*\*\*\*  
Peak Number 2 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 5

| R.T.      | EstConc                              | Area   | Relative to ISTD       | R.T.        |      |
|-----------|--------------------------------------|--------|------------------------|-------------|------|
| 14.78     | 15.42 ug/l                           | 169420 | 1,4-Dichlorobenzene-d4 | 14.52       |      |
| Hit# of 5 | Tentative ID                         | MW     | MolForm                | CAS#        | Qual |
| 1         | Benzene, 2-ethyl-1,4-dimethyl-       | 134    | C10H14                 | 001758-88-9 | 94   |
| 2         | o-Cymene                             | 134    | C10H14                 | 000527-84-4 | 94   |
| 3         | Benzene, 1-methyl-3-(1-methylethyl)- | 134    | C10H14                 | 000535-77-3 | 93   |
| 4         | Benzene, 1-ethyl-2,3-dimethyl-       | 134    | C10H14                 | 000933-98-2 | 91   |
| 5         | Benzene, 4-ethyl-1,2-dimethyl-       | 134    | C10H14                 | 000934-80-5 | 91   |



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Quant Title : SW846 8260

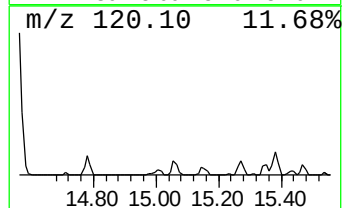
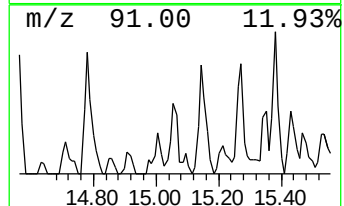
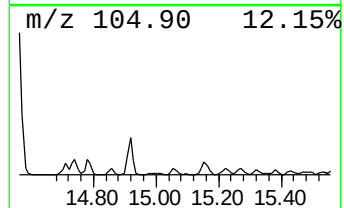
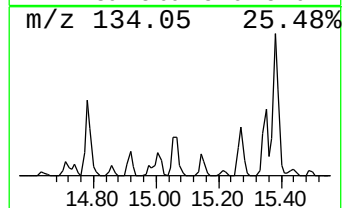
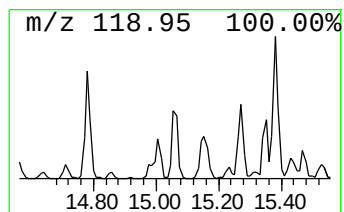
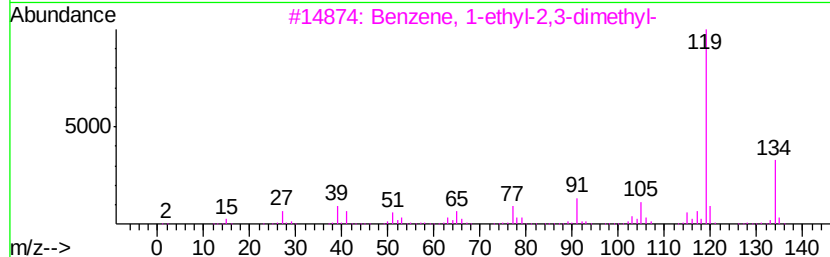
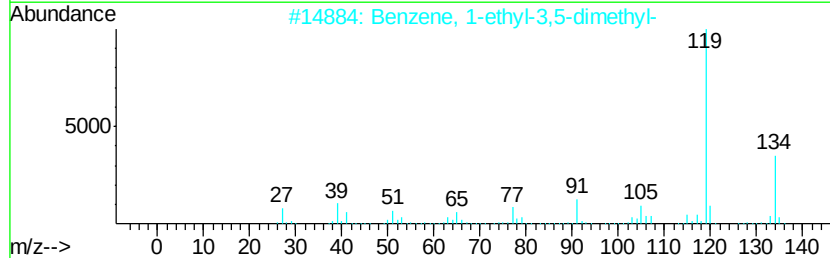
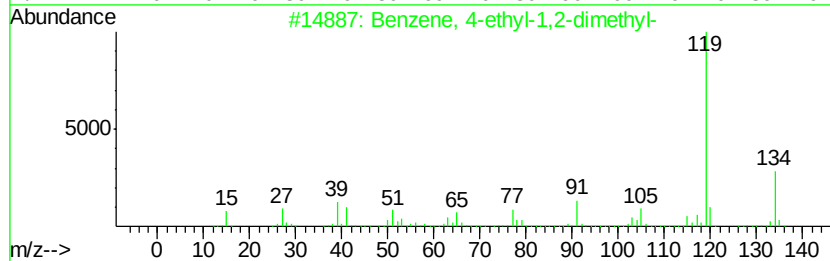
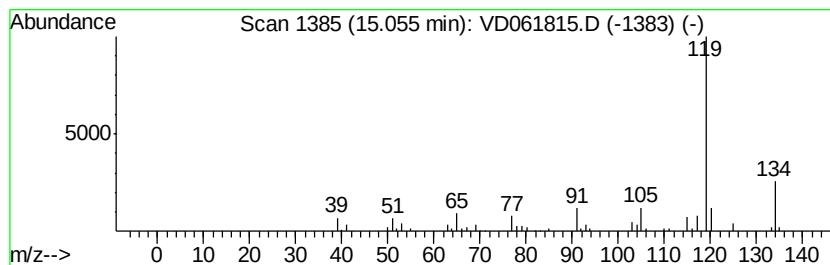
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 10

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 15.06 | 9.90 ug/l | 108732 | 1,4-Dichlorobenzene-d4 | 14.52 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 90   |
| 2    |    | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 90   |
| 3    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 90   |
| 4    |    | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 87   |
| 5    |    | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 87   |



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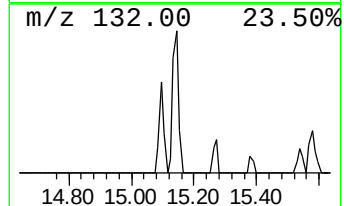
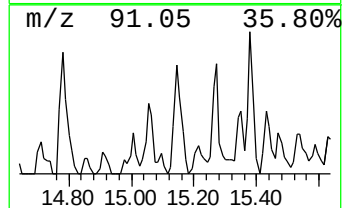
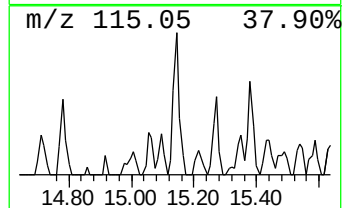
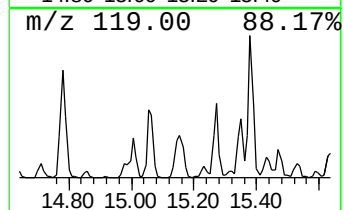
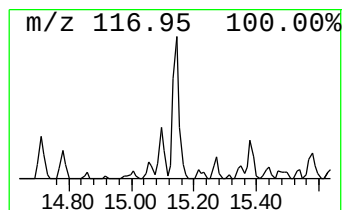
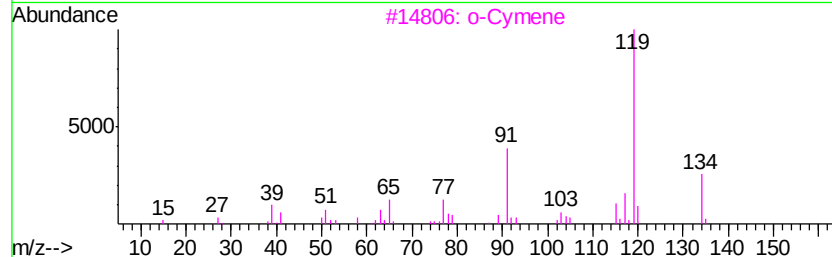
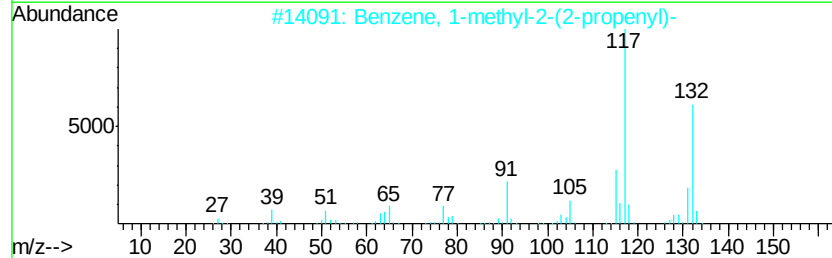
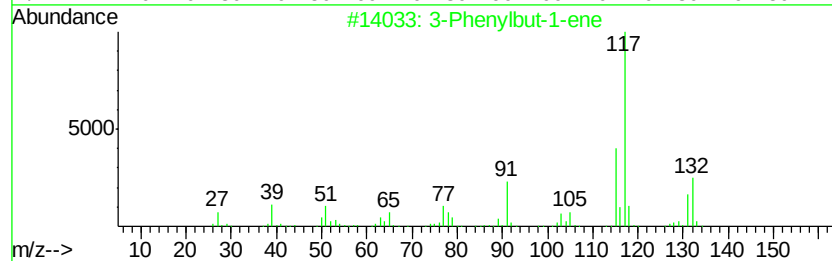
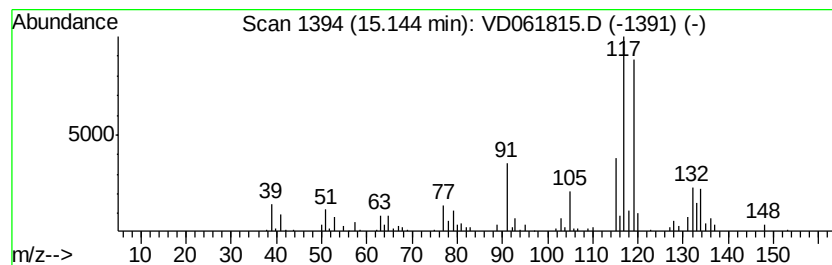
Instrument :  
MSVOA\_D  
ClientSampleId :  
TP-7-S2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
Quant Title : SW846 8260

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

\*\*\*\*\*  
Peak Number 4 3-Phenylbut-1-ene Concentration Rank 3

| R.T.    | EstConc                           | Area         | Relative to ISTD       | R.T.      |
|---------|-----------------------------------|--------------|------------------------|-----------|
| 15.14   | 16.20 ug/l                        | 177966       | 1,4-Dichlorobenzene-d4 | 14.52     |
| Hit# of | 5                                 | Tentative ID | MW MolForm             | CAS# Qual |
| 1       | 3-Phenylbut-1-ene                 | 132 C10H12   | 000934-10-1            | 50        |
| 2       | Benzene, 1-methyl-2-(2-propenyl)- | 132 C10H12   | 001587-04-8            | 46        |
| 3       | o-Cymene                          | 134 C10H14   | 000527-84-4            | 46        |
| 4       | Benzene, tert-butyl-              | 134 C10H14   | 000098-06-6            | 42        |
| 5       | Benzene, 1-ethyl-2,3-dimethyl-    | 134 C10H14   | 000933-98-2            | 41        |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041619\  
Data File : VD061815.D  
Acq On : 16 Apr 2019 17:03  
Operator : VA/AP  
Sample : K2358-10  
Misc : 4.88g/5ml/MSVOA D/SOIL  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
MSVOA\_D  
ClientSampleId :  
TP-7-S2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
Quant Title : SW846 8260

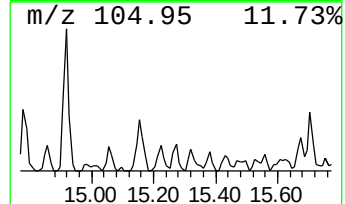
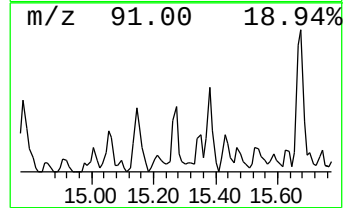
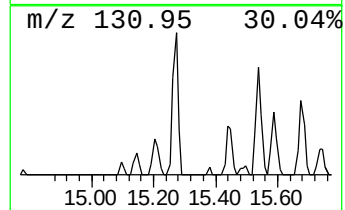
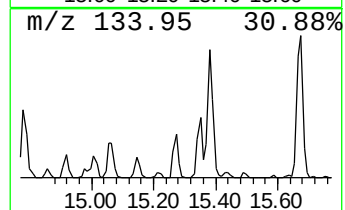
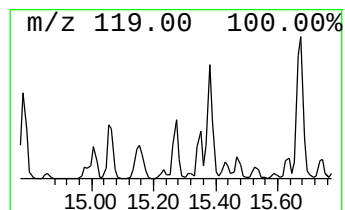
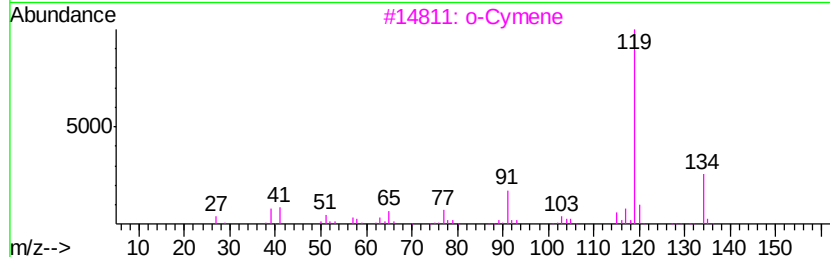
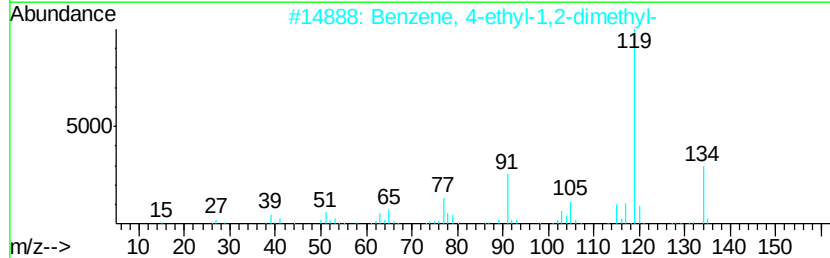
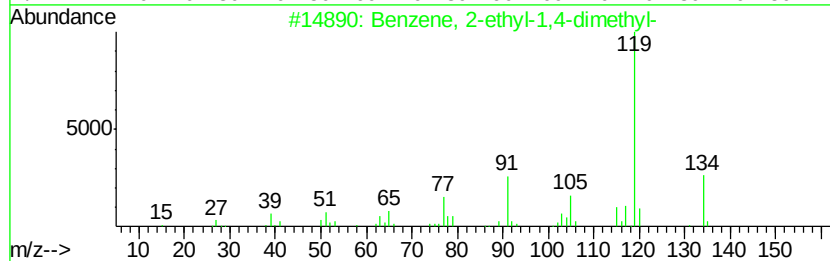
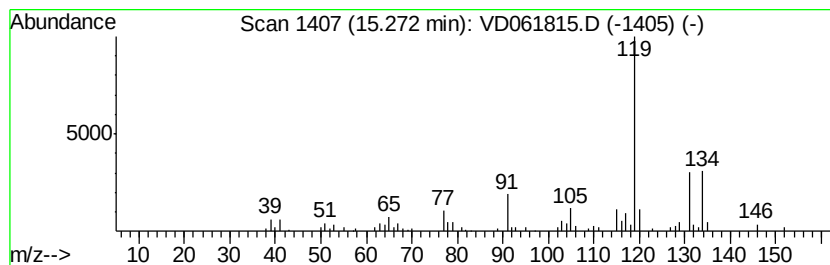
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 15.27 | 15.92 ug/l | 174924 | 1,4-Dichlorobenzene-d4 | 14.52 |

| Hit# of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|---------|-------------------------------------|-----|---------|-------------|------|
| 1       | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 95   |
| 2       | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 94   |
| 3       | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 94   |
| 4       | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 93   |
| 5       | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 93   |





Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041619\  
Data File : VD061815.D  
Acq On : 16 Apr 2019 17:03  
Operator : VA/AP  
Sample : K2358-10  
Misc : 4.88g/5ml/MSVOA D/SOIL  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
MSVOA\_D  
ClientSampleId :  
TP-7-S2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
Quant Title : SW846 8260

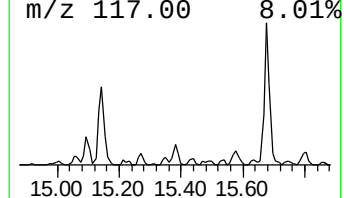
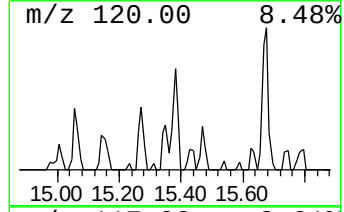
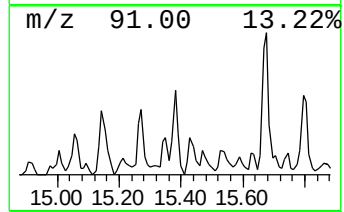
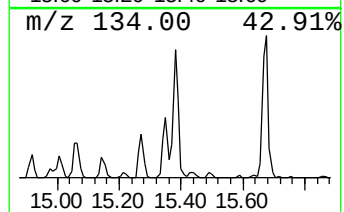
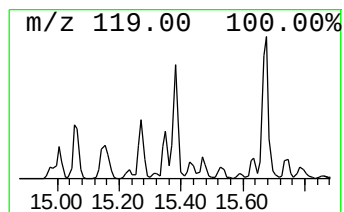
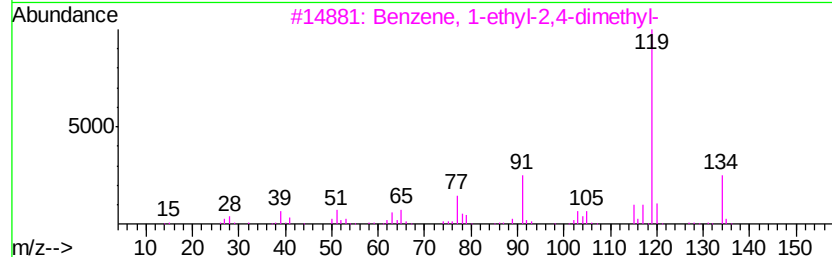
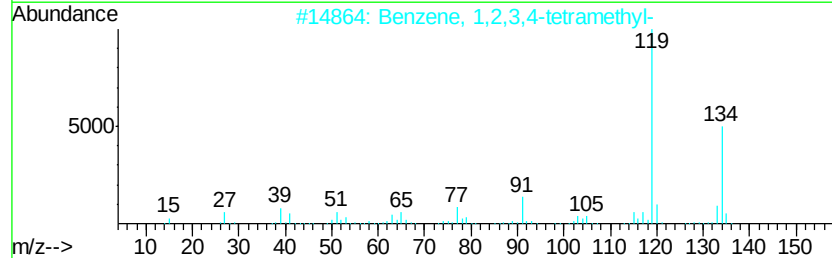
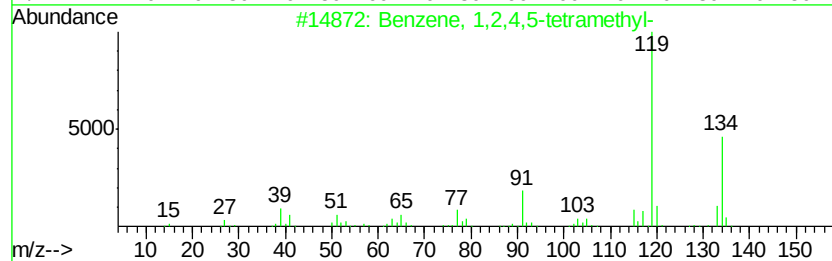
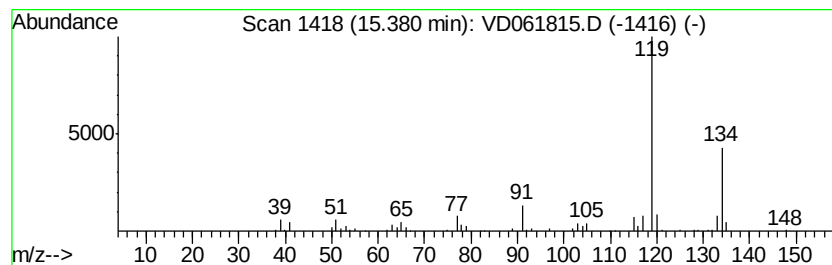
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 2

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 15.38 | 17.02 ug/l | 186965 | 1,4-Dichlorobenzene-d4 | 14.52 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 95   |
| 2    |    | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 94   |
| 3    |    | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 94   |
| 4    |    | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 94   |
| 5    |    | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041619\  
Data File : VD061815.D  
Acq On : 16 Apr 2019 17:03  
Operator : VA/AP  
Sample : K2358-10  
Misc : 4.88g/5ml/MSVOA D/SOIL  
ALS Vial : 14 Sample Multiplier: 1

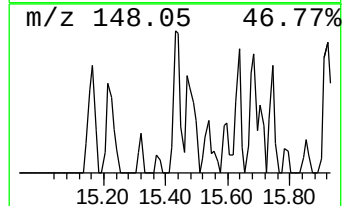
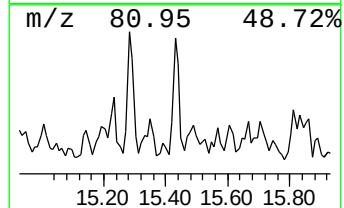
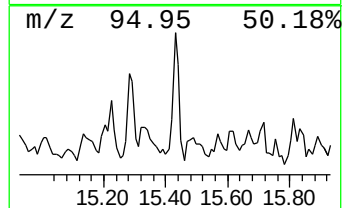
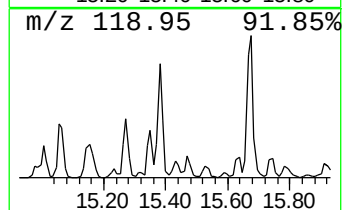
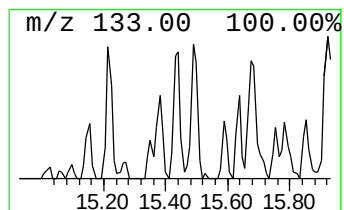
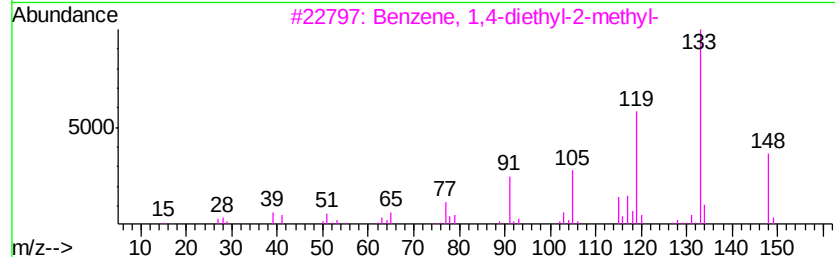
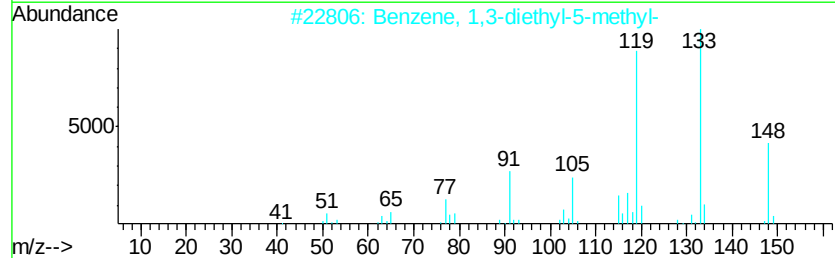
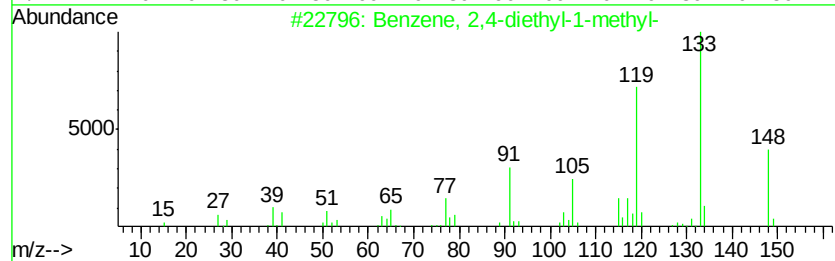
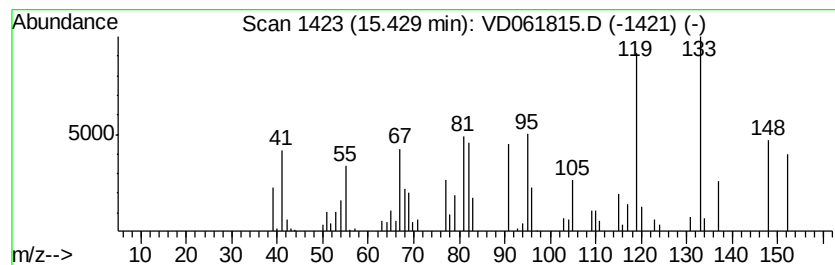
Instrument :  
MSVOA\_D  
ClientSampleId :  
TP-7-S2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
Quant Title : SW846 8260

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

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Peak Number 7 Benzene, 2,4-diethyl-1-methyl- Concentration Rank 8

| R.T.    | EstConc    | Area                                | Relative to ISTD       | R.T.            |
|---------|------------|-------------------------------------|------------------------|-----------------|
| 15.43   | 11.72 ug/l | 128692                              | 1,4-Dichlorobenzene-d4 | 14.52           |
| Hit# of | 5          | Tentative ID                        | MW MolForm             | CAS# Qual       |
| 1       |            | Benzene, 2,4-diethyl-1-methyl-      | 148 C11H16             | 001758-85-6 50  |
| 2       |            | Benzene, 1,3-diethyl-5-methyl-      | 148 C11H16             | 002050-24-0 50  |
| 3       |            | Benzene, 1,4-diethyl-2-methyl-      | 148 C11H16             | 013632-94-5 30  |
| 4       |            | 3,4-Dihydro-2-quinoxalinal          | 148 C8H8N2O            | 1000332-36-8 18 |
| 5       |            | 5H-Cyclopentapyrazine, 6,7-dihyd... | 148 C9H12N2            | 038917-61-2 18  |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041619\  
Data File : VD061815.D  
Acq On : 16 Apr 2019 17:03  
Operator : VA/AP  
Sample : K2358-10  
Misc : 4.88g/5ml/MSVOA D/SOIL  
ALS Vial : 14 Sample Multiplier: 1

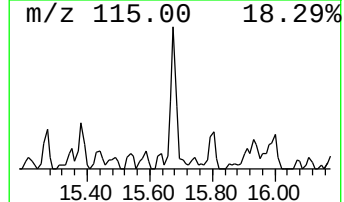
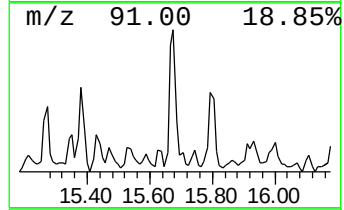
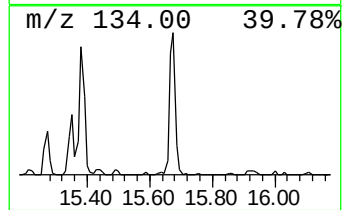
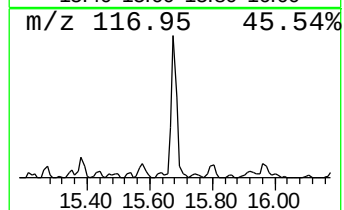
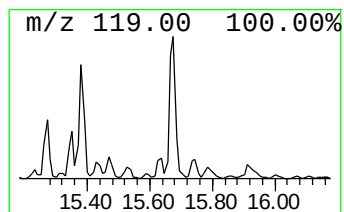
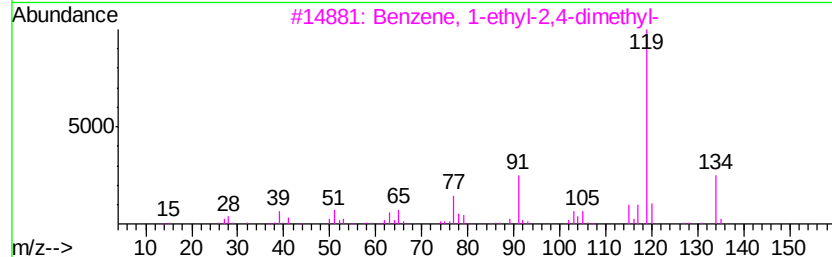
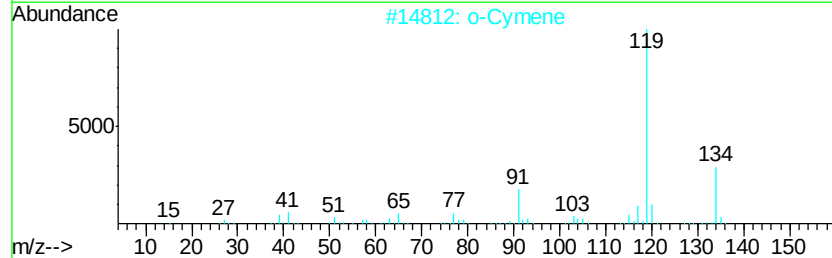
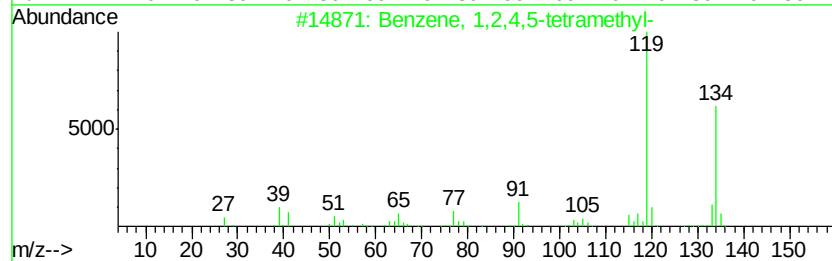
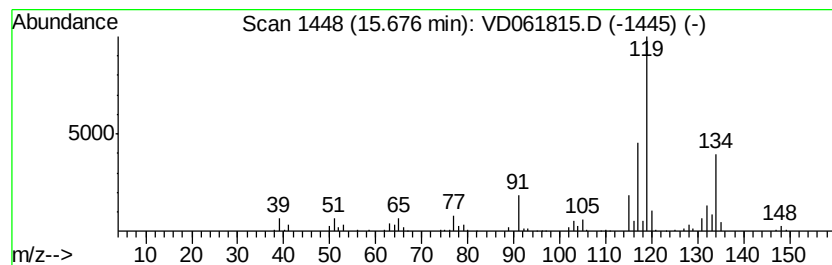
Instrument :  
MSVOA\_D  
ClientSampleId :  
TP-7-S2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
Quant Title : SW846 8260

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

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Peak Number 8 o-Cymene Concentration Rank 1

| R.T.      | EstConc                        | Area   | Relative to ISTD       | R.T.        |      |
|-----------|--------------------------------|--------|------------------------|-------------|------|
| 15.68     | 34.44 ug/l                     | 378312 | 1,4-Dichlorobenzene-d4 | 14.52       |      |
| Hit# of 5 | Tentative ID                   | MW     | MolForm                | CAS#        | Qual |
| 1         | Benzene, 1,2,4,5-tetramethyl-  | 134    | C10H14                 | 000095-93-2 | 76   |
| 2         | o-Cymene                       | 134    | C10H14                 | 000527-84-4 | 70   |
| 3         | Benzene, 1-ethyl-2,4-dimethyl- | 134    | C10H14                 | 000874-41-9 | 70   |
| 4         | Benzene, 1,2,3,4-tetramethyl-  | 134    | C10H14                 | 000488-23-3 | 70   |
| 5         | Benzene, 1-ethyl-2,3-dimethyl- | 134    | C10H14                 | 000933-98-2 | 68   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041619\  
Data File : VD061815.D  
Acq On : 16 Apr 2019 17:03  
Operator : VA/AP  
Sample : K2358-10  
Misc : 4.88g/5ml/MSVOA D/SOIL  
ALS Vial : 14 Sample Multiplier: 1

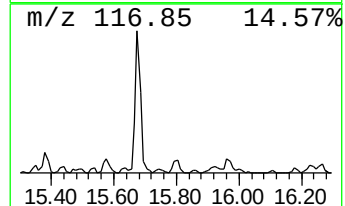
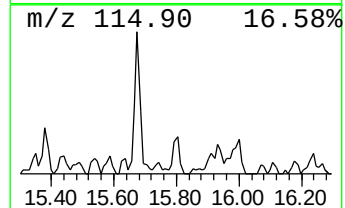
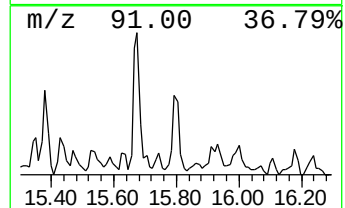
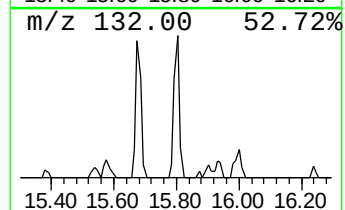
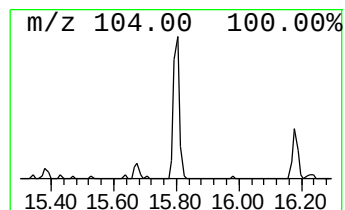
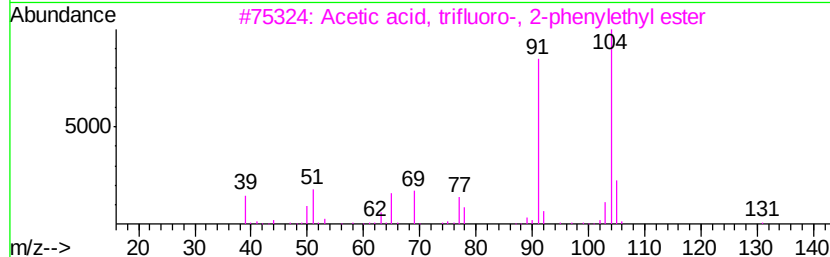
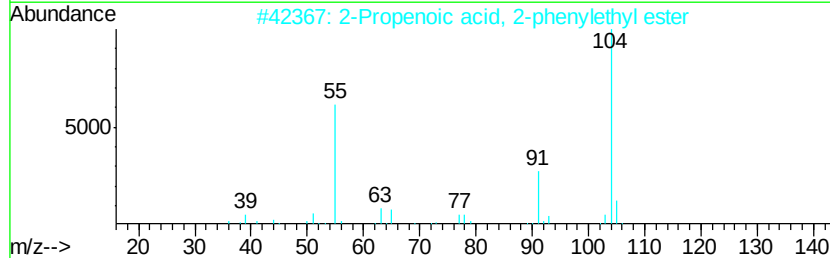
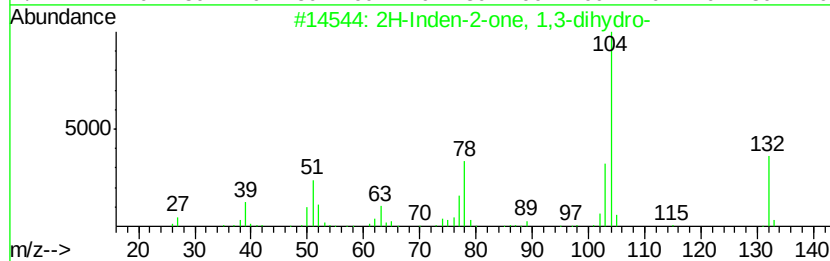
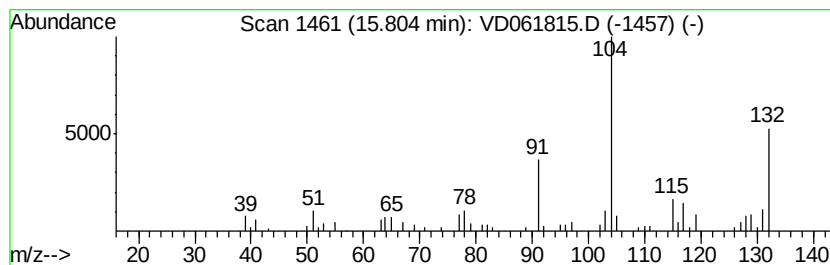
Instrument :  
MSVOA\_D  
ClientSampleId :  
TP-7-S2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
Quant Title : SW846 8260

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

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Peak Number 9 2H-Inden-2-one, 1,3-dihydro- Concentration Rank 6

| R.T.    | EstConc                              | Area           | Relative to ISTD       | R.T.      |
|---------|--------------------------------------|----------------|------------------------|-----------|
| 15.80   | 12.73 ug/l                           | 139831         | 1,4-Dichlorobenzene-d4 | 14.52     |
| Hit# of | 5                                    | Tentative ID   | MW MolForm             | CAS# Qual |
| 1       | 2H-Inden-2-one, 1,3-dihydro-         | 132 C9H8O      | 000615-13-4            | 40        |
| 2       | 2-Propenoic acid, 2-phenylethyl ...  | 176 C11H12O2   | 003530-36-7            | 38        |
| 3       | Acetic acid, trifluoro-, 2-phenyl... | 218 C10H9F3O2  | 055419-66-4            | 38        |
| 4       | Bromoacetic acid, 2-phenylethyl ...  | 242 C10H11BrO2 | 003785-33-9            | 38        |
| 5       | Acetic acid, 2-phenylethyl ester     | 164 C10H12O2   | 000103-45-7            | 38        |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041619\  
Data File : VD061815.D  
Acq On : 16 Apr 2019 17:03  
Operator : VA/AP  
Sample : K2358-10  
Misc : 4.88g/5ml/MSVOA D/SOIL  
ALS Vial : 14 Sample Multiplier: 1

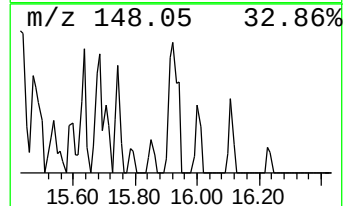
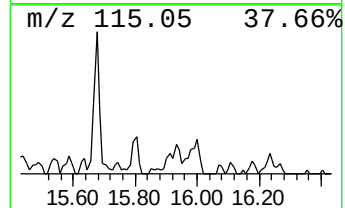
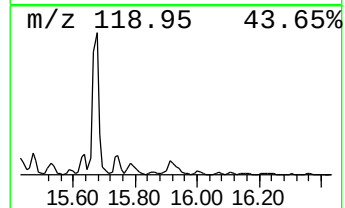
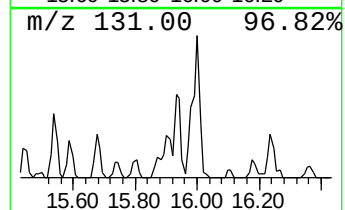
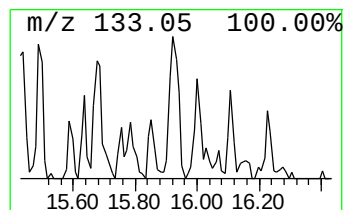
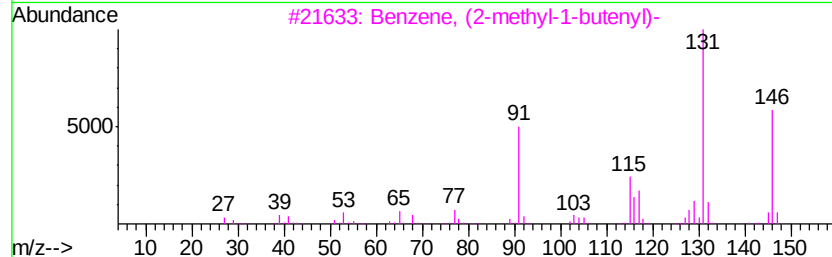
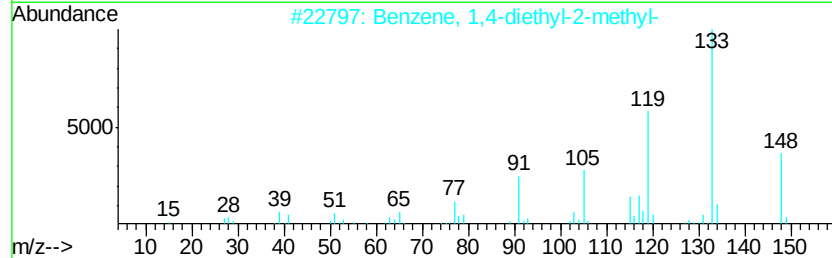
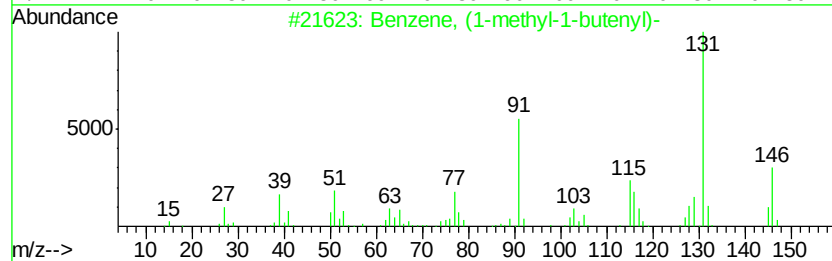
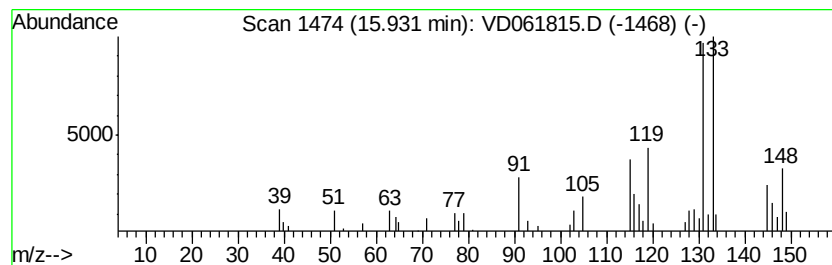
Instrument :  
MSVOA\_D  
ClientSampleId :  
TP-7-S2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
Quant Title : SW846 8260

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

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Peak Number 10 Benzene, (1-methyl-1-butenyl)- Concentration Rank 7

| R.T.      | EstConc                             | Area   | Relative to ISTD       | R.T.        |      |
|-----------|-------------------------------------|--------|------------------------|-------------|------|
| 15.93     | 11.83 ug/l                          | 129901 | 1,4-Dichlorobenzene-d4 | 14.52       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm                | CAS#        | Qual |
| 1         | Benzene, (1-methyl-1-butenyl)-      | 146    | C11H14                 | 053172-84-2 | 80   |
| 2         | Benzene, 1,4-diethyl-2-methyl-      | 148    | C11H16                 | 013632-94-5 | 45   |
| 3         | Benzene, (2-methyl-1-butenyl)-      | 146    | C11H14                 | 056253-64-6 | 42   |
| 4         | 1,3-Cyclopentadiene, 5-(trans-2-... | 146    | C11H14                 | 079209-36-2 | 38   |
| 5         | Benzene, 1-ethyl-4-(1-methylethyl)- | 148    | C11H16                 | 004218-48-8 | 38   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_D\DATA\VD041619\  
Data File : VD061815.D  
Acq On : 16 Apr 2019 17:03  
Operator : VA/AP  
Sample : K2358-10  
Misc : 4.88g/5ml/MSVOA\_D/SOIL  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
MSVOA\_D  
ClientSampleId :  
TP-7-S2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_D\METHOD\82D041519S.M  
Quant Title : SW846 8260

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Ethane, 1,1-diflu... | 1.57  | 11.0    | ug/l  | 256849   | 1                     | 7.06  | 1170690 | 50.0 |
| Benzene, 2-ethyl-... | 14.78 | 15.4    | ug/l  | 169420   | 4                     | 14.52 | 549221  | 50.0 |
| Benzene, 4-ethyl-... | 15.06 | 9.9     | ug/l  | 108732   | 4                     | 14.52 | 549221  | 50.0 |
| 3-Phenylbut-1-ene    | 15.14 | 16.2    | ug/l  | 177966   | 4                     | 14.52 | 549221  | 50.0 |
| Benzene, 1-ethyl-... | 15.27 | 15.9    | ug/l  | 174924   | 4                     | 14.52 | 549221  | 50.0 |
| Benzene, 1,2,4,5-... | 15.38 | 17.0    | ug/l  | 186965   | 4                     | 14.52 | 549221  | 50.0 |
| Benzene, 2,4-diet... | 15.43 | 11.7    | ug/l  | 128692   | 4                     | 14.52 | 549221  | 50.0 |
| o-Cymene             | 15.68 | 34.4    | ug/l  | 378312   | 4                     | 14.52 | 549221  | 50.0 |
| 2H-Inden-2-one, 1... | 15.80 | 12.7    | ug/l  | 139831   | 4                     | 14.52 | 549221  | 50.0 |
| Benzene, (1-methy... | 15.93 | 11.8    | ug/l  | 129901   | 4                     | 14.52 | 549221  | 50.0 |