

Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
 Data File : VU038128.D  
 Acq On : 12 May 2020 16:45  
 Operator : JC/MD  
 Sample : L2570-07  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BFSP5

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

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\_U\METHOD\SOMUTR050720WMA.M

Title : TRACE VOA SOM01.0

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         | 1.613       | 73            | 85          | 94           | rBV2     | 278487         | 361841        | 8.55%           | 0.715%        |
| 2         | 1.932       | 176           | 184         | 201          | rVB      | 179668         | 243902        | 5.76%           | 0.482%        |
| 3         | 2.398       | 318           | 329         | 350          | rVB2     | 141733         | 263919        | 6.23%           | 0.521%        |
| 4         | 2.591       | 376           | 389         | 410          | rBV      | 296258         | 496188        | 11.72%          | 0.980%        |
| 5         | 3.073       | 528           | 539         | 545          | rBV3     | 32646          | 66186         | 1.56%           | 0.131%        |
| 6         | 3.394       | 620           | 639         | 657          | rBV2     | 187402         | 422012        | 9.97%           | 0.834%        |
| 7         | 3.909       | 782           | 799         | 819          | rBV      | 98974          | 239523        | 5.66%           | 0.473%        |
| 8         | 4.333       | 911           | 931         | 957          | rBV4     | 30531          | 93446         | 2.21%           | 0.185%        |
| 9         | 4.472       | 957           | 974         | 989          | rVV3     | 37438          | 94152         | 2.22%           | 0.186%        |
| 10        | 4.571       | 989           | 1005        | 1023         | rVV2     | 99273          | 236638        | 5.59%           | 0.468%        |
| 11        | 4.703       | 1023          | 1046        | 1110         | rVV3     | 524139         | 1704956       | 40.28%          | 3.369%        |
| 12        | 5.102       | 1152          | 1170        | 1182         | rBV      | 263657         | 585654        | 13.84%          | 1.157%        |
| 13        | 5.166       | 1182          | 1190        | 1212         | rVB      | 78829          | 187208        | 4.42%           | 0.370%        |
| 14        | 5.420       | 1253          | 1269        | 1281         | rVV3     | 191013         | 488051        | 11.53%          | 0.964%        |
| 15        | 5.494       | 1281          | 1292        | 1313         | rVB2     | 196646         | 447982        | 10.58%          | 0.885%        |
| 16        | 5.626       | 1319          | 1333        | 1340         | rBV4     | 64288          | 138915        | 3.28%           | 0.274%        |
| 17        | 5.668       | 1342          | 1346        | 1356         | rVV4     | 54484          | 99329         | 2.35%           | 0.196%        |
| 18        | 5.761       | 1356          | 1375        | 1381         | rVV2     | 564796         | 1402656       | 33.14%          | 2.771%        |
| 19        | 5.809       | 1381          | 1390        | 1404         | rVV      | 1145677        | 2410140       | 56.94%          | 4.762%        |
| 20        | 5.880       | 1404          | 1412        | 1425         | rVV2     | 150510         | 321673        | 7.60%           | 0.636%        |
| 21        | 5.954       | 1425          | 1435        | 1446         | rVV3     | 76534          | 163870        | 3.87%           | 0.324%        |
| 22        | 6.021       | 1446          | 1456        | 1471         | rVB3     | 46829          | 103789        | 2.45%           | 0.205%        |
| 23        | 6.279       | 1522          | 1536        | 1561         | rBV      | 452979         | 877343        | 20.73%          | 1.733%        |
| 24        | 6.536       | 1599          | 1616        | 1620         | rBV2     | 58929          | 108855        | 2.57%           | 0.215%        |
| 25        | 6.571       | 1622          | 1627        | 1640         | rVB3     | 77881          | 133620        | 3.16%           | 0.264%        |
| 26        | 6.719       | 1659          | 1673        | 1685         | rBV      | 346830         | 695135        | 16.42%          | 1.373%        |
| 27        | 6.890       | 1718          | 1726        | 1743         | rVB4     | 24320          | 47590         | 1.12%           | 0.094%        |
| 28        | 7.590       | 1929          | 1944        | 1965         | rBV      | 185080         | 340524        | 8.04%           | 0.673%        |
| 29        | 7.922       | 2021          | 2047        | 2060         | rBV      | 686088         | 1280521       | 30.25%          | 2.530%        |
| 30        | 7.992       | 2060          | 2069        | 2082         | rVV      | 158187         | 286485        | 6.77%           | 0.566%        |
| 31        | 8.201       | 2121          | 2134        | 2144         | rBV      | 120402         | 201163        | 4.75%           | 0.397%        |
| 32        | 8.262       | 2144          | 2153        | 2176         | rVB3     | 42641          | 84488         | 2.00%           | 0.167%        |
| 33        | 8.388       | 2176          | 2192        | 2205         | rBV3     | 30647          | 57330         | 1.35%           | 0.113%        |
| 34        | 8.658       | 2260          | 2276        | 2307         | rBV      | 1075020        | 2181440       | 51.53%          | 4.310%        |

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 Sample : L2570-07  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BFSP5

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.1  
 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\_U\METHOD\SOMUTR050720WMA.M  
 Title : TRACE VOA SOM01.0

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 35 | 8.986  | 2371 | 2378 | 2392 | rVB5 | 30728   | 54192   | 1.28%   | 0.107% |
| 36 | 9.311  | 2464 | 2479 | 2484 | rBV5 | 27536   | 53291   | 1.26%   | 0.105% |
| 37 | 9.343  | 2484 | 2489 | 2502 | rVB6 | 32485   | 56397   | 1.33%   | 0.111% |
| 38 | 9.465  | 2505 | 2527 | 2553 | rBV2 | 1414705 | 3471493 | 82.01%  | 6.859% |
| 39 | 9.587  | 2553 | 2565 | 2590 | rVV  | 871877  | 1589044 | 37.54%  | 3.140% |
| 40 | 9.709  | 2590 | 2603 | 2618 | rVV  | 2479604 | 4233116 | 100.00% | 8.364% |
| 41 | 9.777  | 2618 | 2624 | 2645 | rVB2 | 85000   | 154438  | 3.65%   | 0.305% |
| 42 | 10.114 | 2697 | 2729 | 2752 | rBV  | 592071  | 1047384 | 24.74%  | 2.069% |
| 43 | 10.497 | 2836 | 2848 | 2860 | rBV  | 255086  | 414838  | 9.80%   | 0.820% |
| 44 | 10.581 | 2862 | 2874 | 2894 | rVB  | 859316  | 1427722 | 33.73%  | 2.821% |
| 45 | 10.770 | 2922 | 2933 | 2945 | rBV  | 279014  | 468635  | 11.07%  | 0.926% |
| 46 | 10.918 | 2963 | 2979 | 2997 | rVB  | 611204  | 1264391 | 29.87%  | 2.498% |
| 47 | 11.008 | 2997 | 3007 | 3013 | rBV2 | 131419  | 233201  | 5.51%   | 0.461% |
| 48 | 11.102 | 3026 | 3036 | 3046 | rVB  | 117080  | 182810  | 4.32%   | 0.361% |
| 49 | 11.243 | 3071 | 3080 | 3088 | rVB  | 35308   | 50663   | 1.20%   | 0.100% |
| 50 | 11.304 | 3088 | 3099 | 3111 | rBV  | 209997  | 338013  | 7.98%   | 0.668% |
| 51 | 11.478 | 3140 | 3153 | 3168 | rVB  | 637300  | 1029475 | 24.32%  | 2.034% |
| 52 | 11.568 | 3168 | 3181 | 3189 | rBV2 | 32479   | 50570   | 1.19%   | 0.100% |
| 53 | 11.622 | 3189 | 3198 | 3202 | rBV  | 62107   | 92399   | 2.18%   | 0.183% |
| 54 | 11.655 | 3203 | 3208 | 3218 | rVB  | 69264   | 99262   | 2.34%   | 0.196% |
| 55 | 11.709 | 3219 | 3225 | 3232 | rVB3 | 38469   | 51688   | 1.22%   | 0.102% |
| 56 | 11.758 | 3232 | 3240 | 3247 | rVB2 | 48360   | 67301   | 1.59%   | 0.133% |
| 57 | 11.822 | 3247 | 3260 | 3272 | rBV2 | 721327  | 1379510 | 32.59%  | 2.726% |
| 58 | 11.905 | 3280 | 3286 | 3295 | rVB  | 235405  | 352900  | 8.34%   | 0.697% |
| 59 | 11.957 | 3295 | 3302 | 3313 | rVB  | 70840   | 106573  | 2.52%   | 0.211% |
| 60 | 12.024 | 3313 | 3323 | 3331 | rBV2 | 73656   | 116115  | 2.74%   | 0.229% |
| 61 | 12.105 | 3331 | 3348 | 3367 | rBV2 | 800887  | 1395314 | 32.96%  | 2.757% |
| 62 | 12.220 | 3367 | 3384 | 3400 | rVB4 | 1377677 | 3370647 | 79.63%  | 6.660% |
| 63 | 12.311 | 3401 | 3412 | 3422 | rBV5 | 36402   | 61398   | 1.45%   | 0.121% |
| 64 | 12.388 | 3423 | 3436 | 3450 | rVB  | 63673   | 132116  | 3.12%   | 0.261% |
| 65 | 12.478 | 3453 | 3464 | 3470 | rBV  | 85837   | 146874  | 3.47%   | 0.290% |
| 66 | 12.555 | 3478 | 3488 | 3495 | rBV  | 392711  | 673023  | 15.90%  | 1.330% |
| 67 | 12.626 | 3504 | 3510 | 3523 | rVB3 | 49062   | 79201   | 1.87%   | 0.156% |
| 68 | 12.696 | 3523 | 3532 | 3552 | rVB4 | 176051  | 442179  | 10.45%  | 0.874% |
| 69 | 12.793 | 3552 | 3562 | 3567 | rBV5 | 47557   | 81931   | 1.94%   | 0.162% |
| 70 | 12.838 | 3567 | 3576 | 3585 | rVV  | 345051  | 536406  | 12.67%  | 1.060% |
| 71 | 12.893 | 3585 | 3593 | 3606 | rVB3 | 61603   | 121859  | 2.88%   | 0.241% |

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 Sample : L2570-07  
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 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BFSP5

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.1  
 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\_U\METHOD\SOMUTR050720WMA.M  
 Title : TRACE VOA SOM01.0

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 72 | 12.992 | 3606 | 3624 | 3632 | rBV  | 199776  | 347564  | 8.21%  | 0.687% |
| 73 | 13.047 | 3632 | 3641 | 3652 | rVB2 | 188616  | 309281  | 7.31%  | 0.611% |
| 74 | 13.130 | 3658 | 3667 | 3671 | rBV4 | 40841   | 61158   | 1.44%  | 0.121% |
| 75 | 13.243 | 3690 | 3702 | 3708 | rBV3 | 22497   | 48231   | 1.14%  | 0.095% |
| 76 | 13.314 | 3715 | 3724 | 3738 | rVB  | 241055  | 365162  | 8.63%  | 0.721% |
| 77 | 13.433 | 3751 | 3761 | 3765 | rBV3 | 34398   | 56296   | 1.33%  | 0.111% |
| 78 | 13.478 | 3765 | 3775 | 3789 | rVB2 | 479641  | 839678  | 19.84% | 1.659% |
| 79 | 13.587 | 3802 | 3809 | 3814 | rBV2 | 37918   | 55784   | 1.32%  | 0.110% |
| 80 | 13.632 | 3815 | 3823 | 3841 | rVB4 | 114121  | 247266  | 5.84%  | 0.489% |
| 81 | 13.880 | 3886 | 3900 | 3909 | rBV  | 649319  | 1152640 | 27.23% | 2.277% |
| 82 | 13.979 | 3922 | 3931 | 3932 | rBV5 | 42401   | 51564   | 1.22%  | 0.102% |
| 83 | 14.114 | 3963 | 3973 | 3993 | rVB  | 52731   | 117588  | 2.78%  | 0.232% |
| 84 | 14.317 | 4027 | 4036 | 4046 | rBV5 | 26462   | 63936   | 1.51%  | 0.126% |
| 85 | 14.519 | 4088 | 4099 | 4106 | rBV2 | 33627   | 53257   | 1.26%  | 0.105% |
| 86 | 14.597 | 4106 | 4123 | 4139 | rVB  | 219616  | 388899  | 9.19%  | 0.768% |
| 87 | 14.847 | 4178 | 4201 | 4217 | rBV  | 142972  | 312449  | 7.38%  | 0.617% |
| 88 | 15.044 | 4246 | 4262 | 4270 | rBV6 | 62151   | 164518  | 3.89%  | 0.325% |
| 89 | 15.121 | 4271 | 4286 | 4296 | rBV2 | 219682  | 433398  | 10.24% | 0.856% |
| 90 | 15.314 | 4336 | 4346 | 4356 | rBV5 | 39150   | 74395   | 1.76%  | 0.147% |
| 91 | 15.597 | 4420 | 4434 | 4446 | rBV2 | 116009  | 194562  | 4.60%  | 0.384% |
| 92 | 15.847 | 4501 | 4512 | 4526 | rBV4 | 66971   | 119536  | 2.82%  | 0.236% |
| 93 | 16.002 | 4543 | 4560 | 4571 | rBV2 | 52274   | 118297  | 2.79%  | 0.234% |
| 94 | 16.146 | 4588 | 4605 | 4619 | rBV  | 1273857 | 2102963 | 49.68% | 4.155% |
| 95 | 16.844 | 4811 | 4822 | 4843 | rBV2 | 214904  | 419398  | 9.91%  | 0.829% |
| 96 | 16.966 | 4850 | 4860 | 4870 | rBV2 | 136314  | 222662  | 5.26%  | 0.440% |

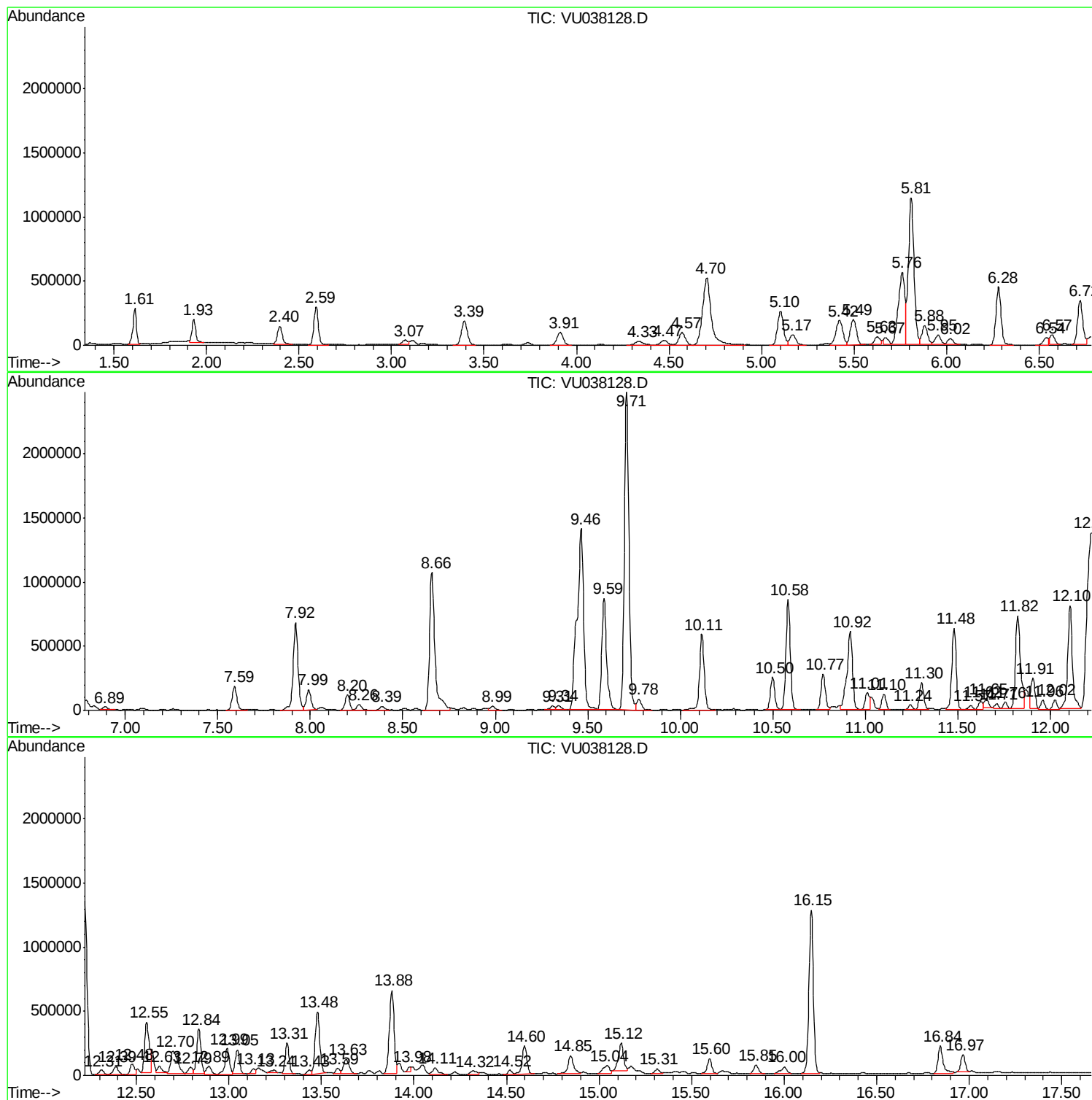
Sum of corrected areas: 50613375

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ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
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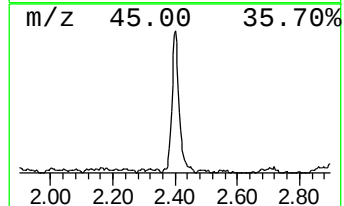
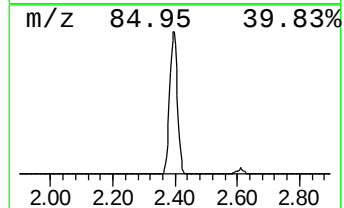
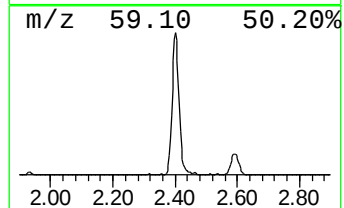
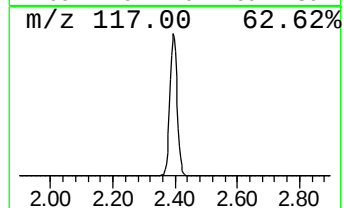
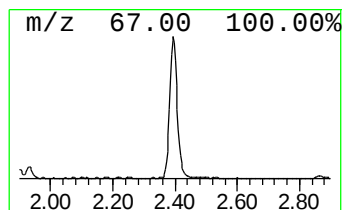
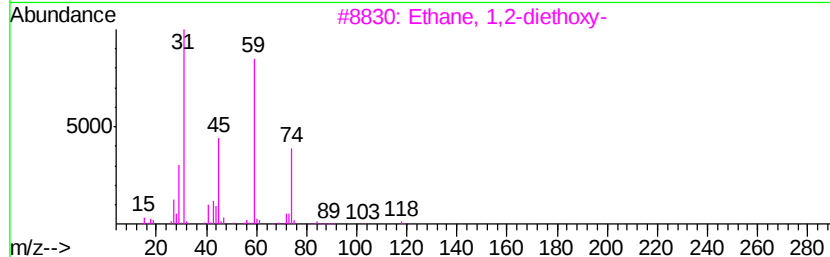
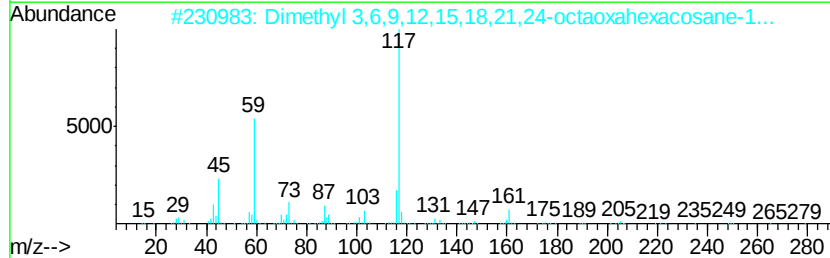
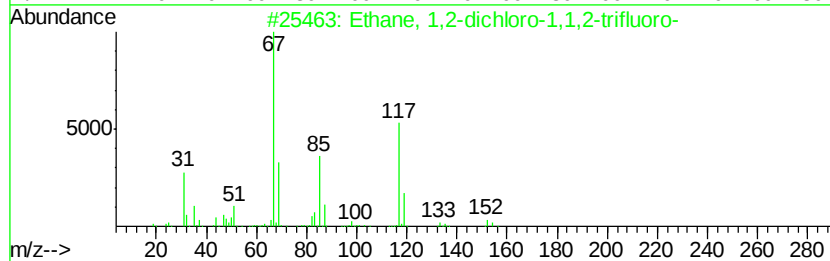
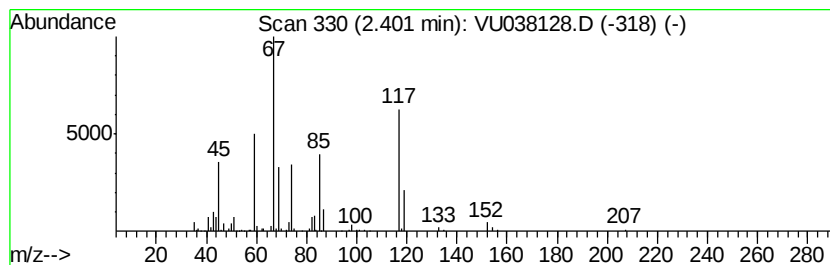
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 1 unknown-01 Concentration Rank 15

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 2.40 | 1.50 ug/L | 263919 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Ethane, 1,2-dichloro-1,1,2-trifl... | 152 | C2HCl2F3  | 000354-23-4  | 43   |
| 2    |    |   | Dimethyl 3,6,9,12,15,18,21,24-oc... | 470 | C20H38O12 | 1000366-80-8 | 16   |
| 3    |    |   | Ethane, 1,2-diethoxy-               | 118 | C6H14O2   | 000629-14-1  | 14   |
| 4    |    |   | Ethyl ether                         | 74  | C4H10O    | 000060-29-7  | 11   |
| 5    |    |   | Ethyl ether                         | 74  | C4H10O    | 000060-29-7  | 10   |



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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

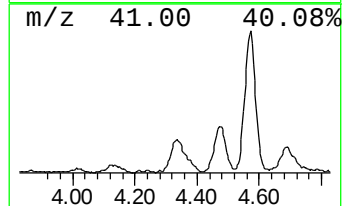
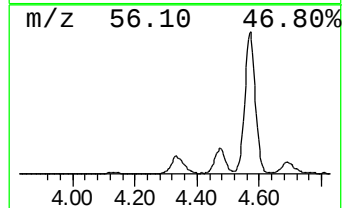
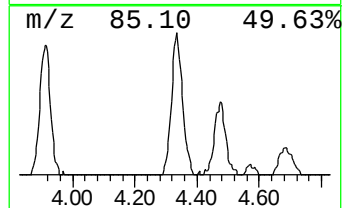
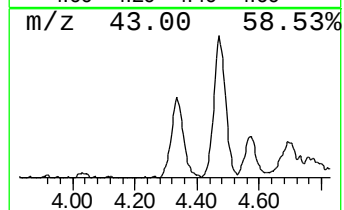
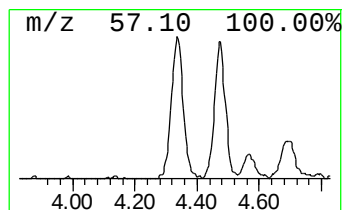
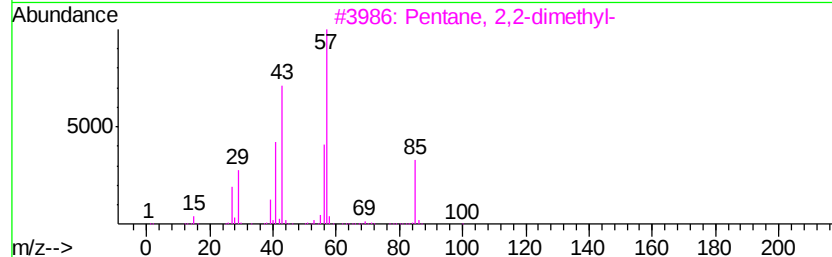
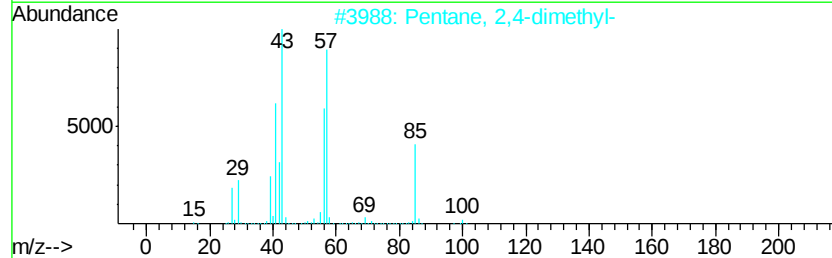
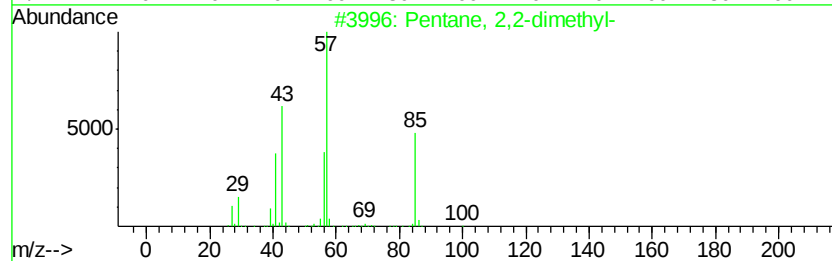
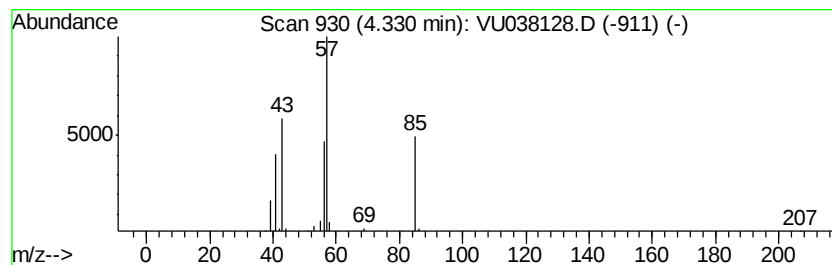
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T. | EstConc   | Area  | Relative to ISTD    | R.T. |
|------|-----------|-------|---------------------|------|
| 4.33 | 0.53 ug/L | 93446 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentane, 2,2-dimethyl-              | 100 | C7H16   | 000590-35-2 | 90   |
| 2    |    | Pentane, 2,4-dimethyl-              | 100 | C7H16   | 000108-08-7 | 74   |
| 3    |    | Pentane, 2,2-dimethyl-              | 100 | C7H16   | 000590-35-2 | 74   |
| 4    |    | Butanoic acid, 2-methyl-, 2-meth... | 158 | C9H18O2 | 002445-67-2 | 72   |
| 5    |    | Hexane, 2,4-dimethyl-               | 114 | C8H18   | 000589-43-5 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

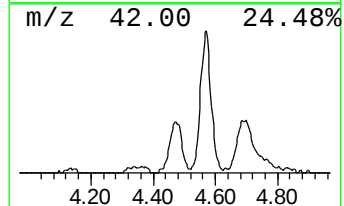
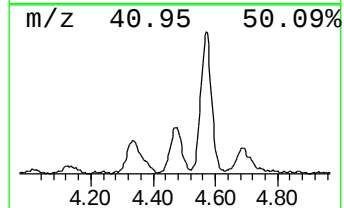
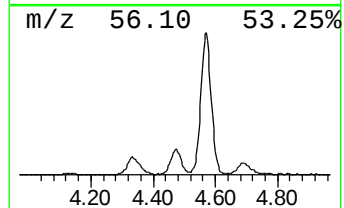
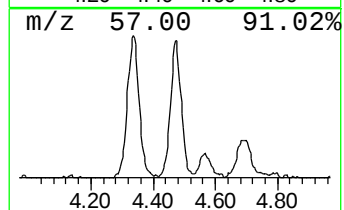
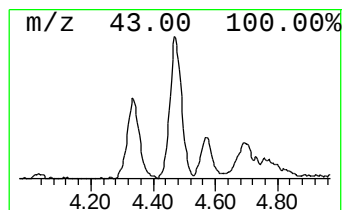
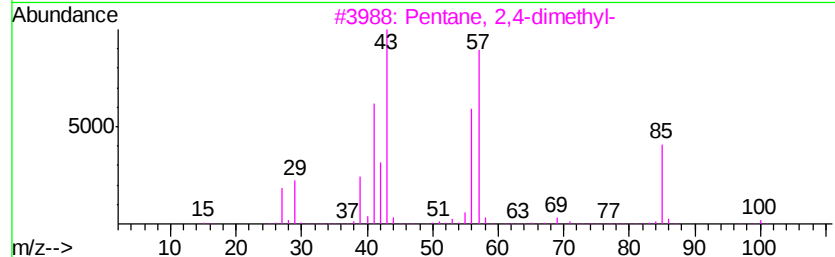
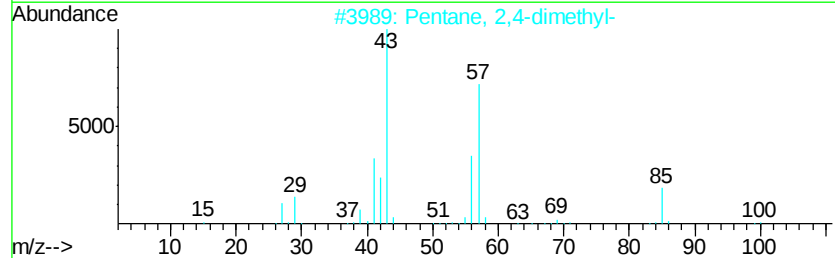
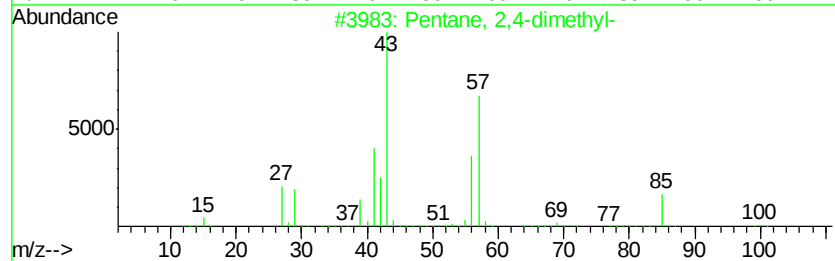
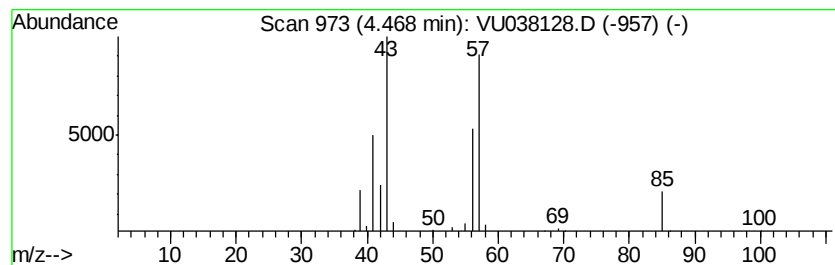
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 35

| R.T. | EstConc   | Area  | Relative to ISTD    | R.T. |
|------|-----------|-------|---------------------|------|
| 4.47 | 0.54 ug/L | 94152 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentane, 2,4-dimethyl-     | 100 | C7H16   | 000108-08-7 | 74   |
| 2    |    | Pentane, 2,4-dimethyl-     | 100 | C7H16   | 000108-08-7 | 52   |
| 3    |    | Pentane, 2,4-dimethyl-     | 100 | C7H16   | 000108-08-7 | 52   |
| 4    |    | Butane, 2,2,3-trimethyl-   | 100 | C7H16   | 000464-06-2 | 50   |
| 5    |    | 2,2,6,6-Tetramethylheptane | 156 | C11H24  | 040117-45-1 | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

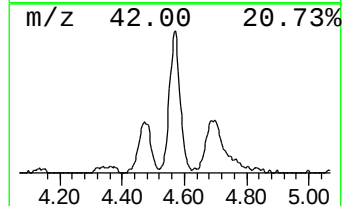
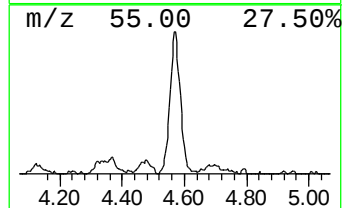
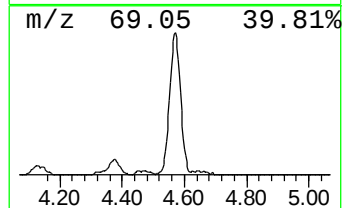
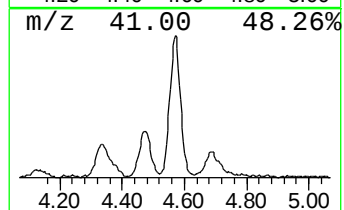
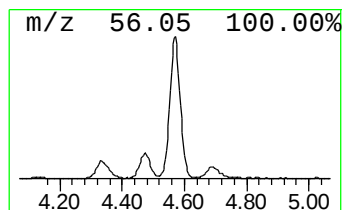
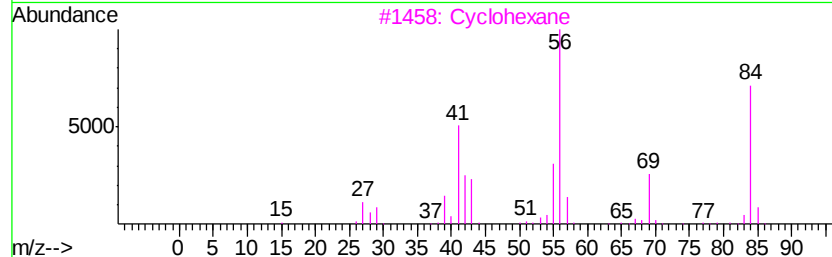
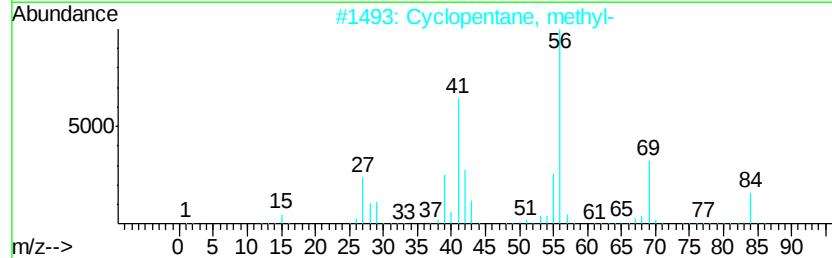
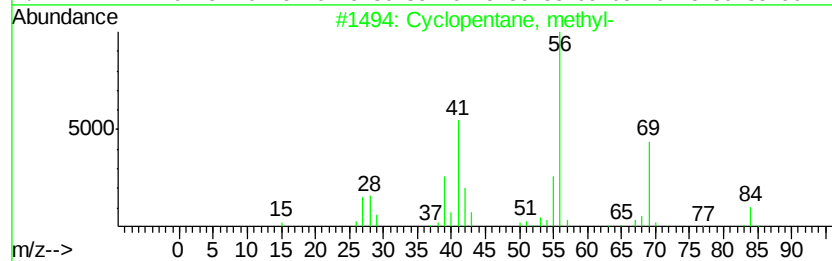
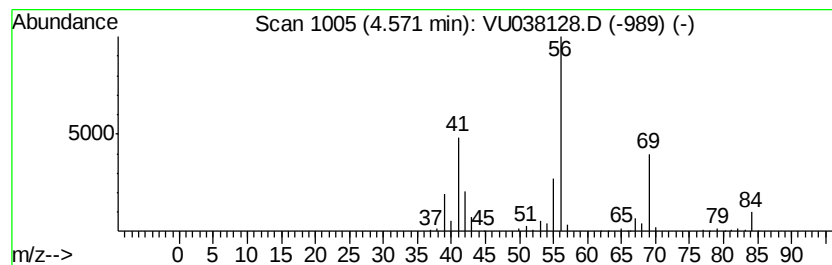
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Cyclic4.57 Concentration Rank 17

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 4.57 | 1.35 ug/L | 236638 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID          | MW | MolForm | CAS#        | Qual |
|------|----|-----------------------|----|---------|-------------|------|
| 1    | 5  | Cyclopentane, methyl- | 84 | C6H12   | 000096-37-7 | 91   |
| 2    |    | Cyclopentane, methyl- | 84 | C6H12   | 000096-37-7 | 86   |
| 3    |    | Cyclohexane           | 84 | C6H12   | 000110-82-7 | 83   |
| 4    |    | Cyclopropane, propyl- | 84 | C6H12   | 002415-72-7 | 78   |
| 5    |    | Cyclobutane, ethyl-   | 84 | C6H12   | 004806-61-5 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

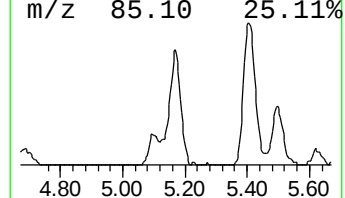
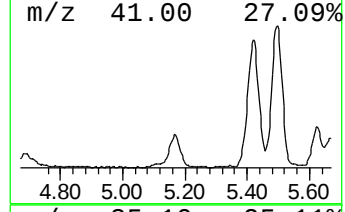
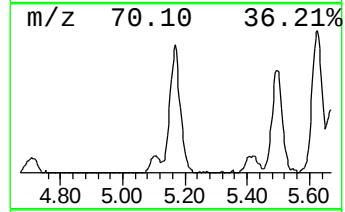
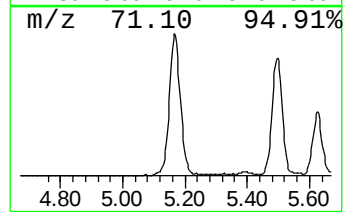
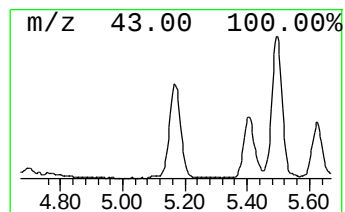
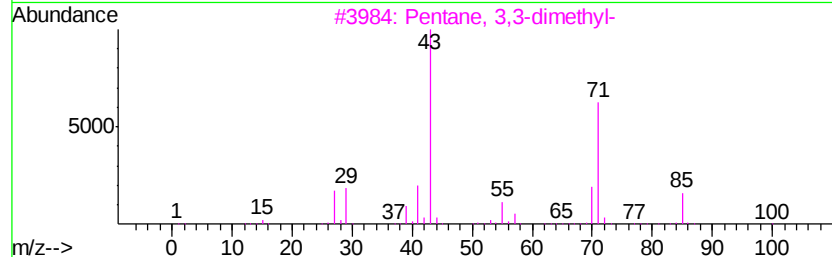
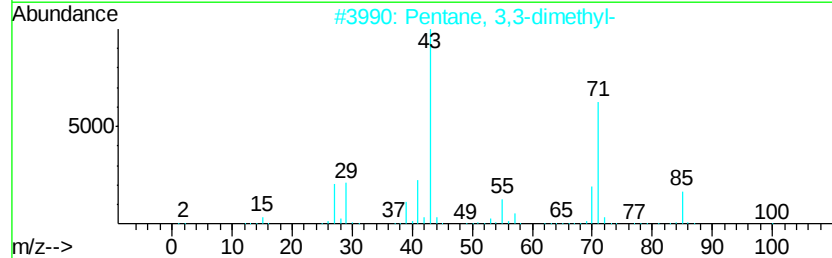
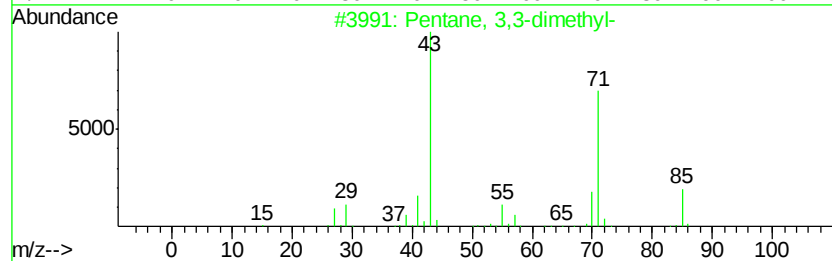
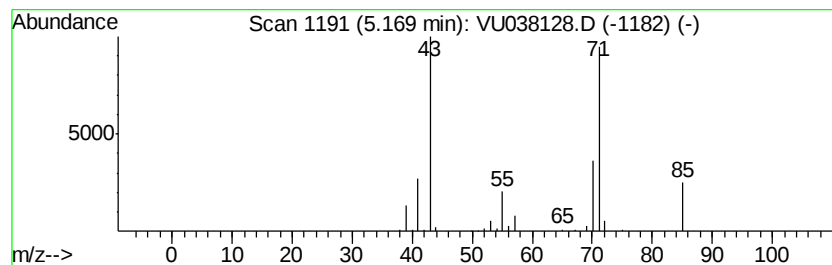
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 5.17 | 1.07 ug/L | 187208 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentane, 3,3-dimethyl-    | 100 | C7H16   | 000562-49-2 | 78   |
| 2    |    | Pentane, 3,3-dimethyl-    | 100 | C7H16   | 000562-49-2 | 56   |
| 3    |    | Pentane, 3,3-dimethyl-    | 100 | C7H16   | 000562-49-2 | 50   |
| 4    |    | Pentane, 2,3,4-trimethyl- | 114 | C8H18   | 000565-75-3 | 42   |
| 5    |    | Pentane, 2,3,4-trimethyl- | 114 | C8H18   | 000565-75-3 | 42   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

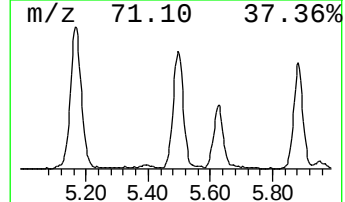
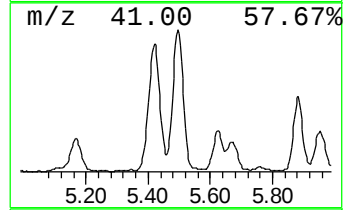
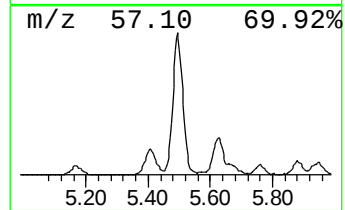
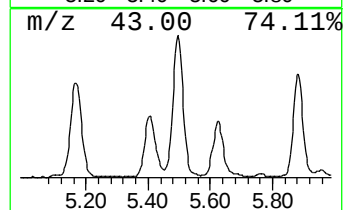
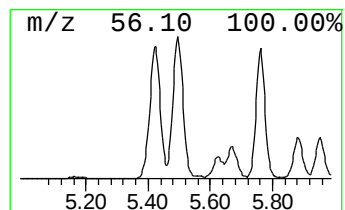
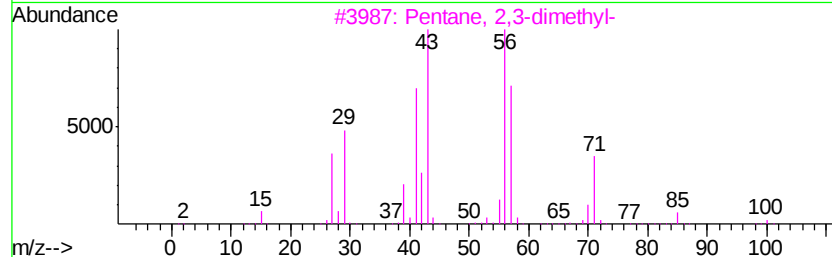
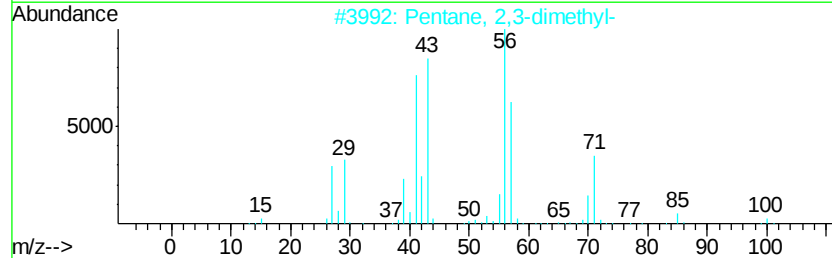
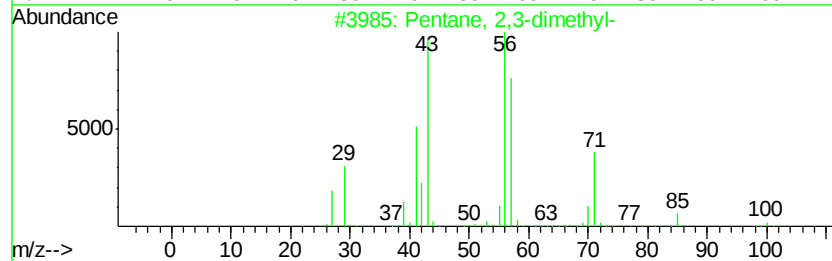
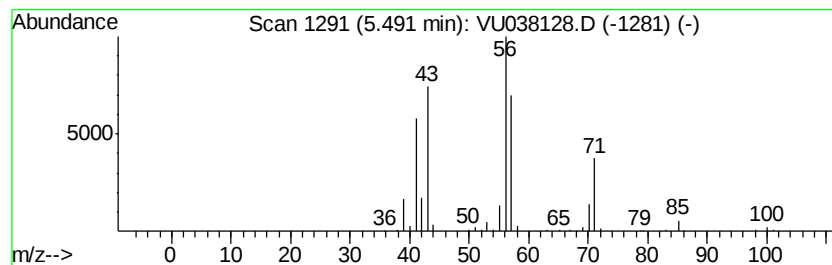
Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

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

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

| R.T.      | EstConc                   | Area   | Relative to ISTD    | R.T.        |      |
|-----------|---------------------------|--------|---------------------|-------------|------|
| 5.49      | 2.55 ug/L                 | 447982 | 1,4-Difluorobenzene | 6.28        |      |
| Hit# of 5 | Tentative ID              | MW     | MolForm             | CAS#        | Qual |
| 1         | Pentane, 2,3-dimethyl-    | 100    | C7H16               | 000565-59-3 | 94   |
| 2         | Pentane, 2,3-dimethyl-    | 100    | C7H16               | 000565-59-3 | 91   |
| 3         | Pentane, 2,3-dimethyl-    | 100    | C7H16               | 000565-59-3 | 74   |
| 4         | Pentane, 2,3-dimethyl-    | 100    | C7H16               | 000565-59-3 | 64   |
| 5         | Heptane, 2,2,4-trimethyl- | 142    | C10H22              | 014720-74-2 | 45   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

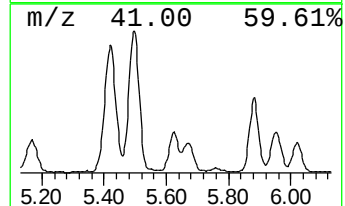
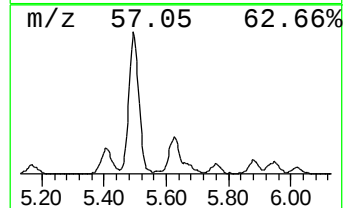
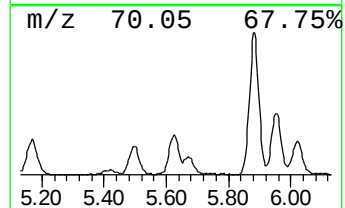
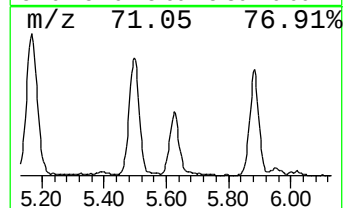
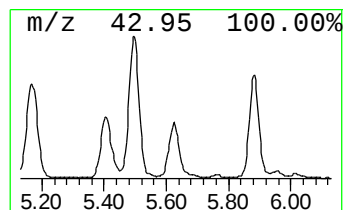
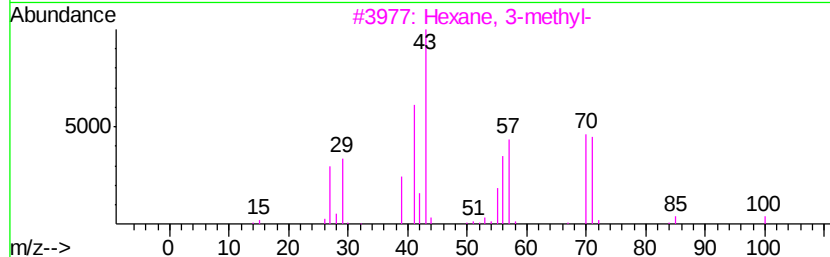
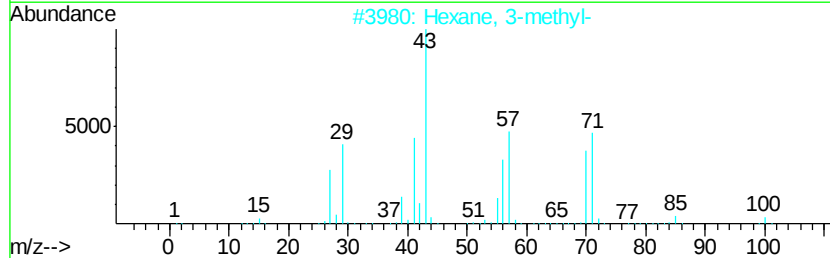
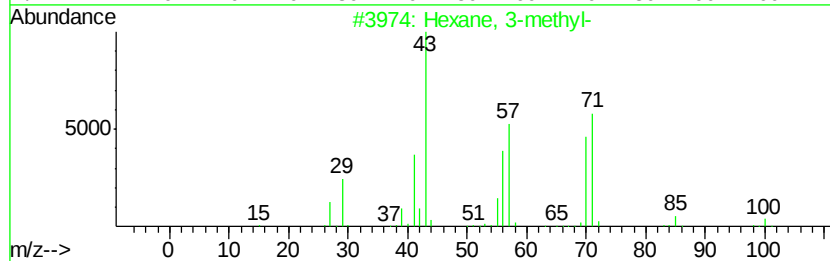
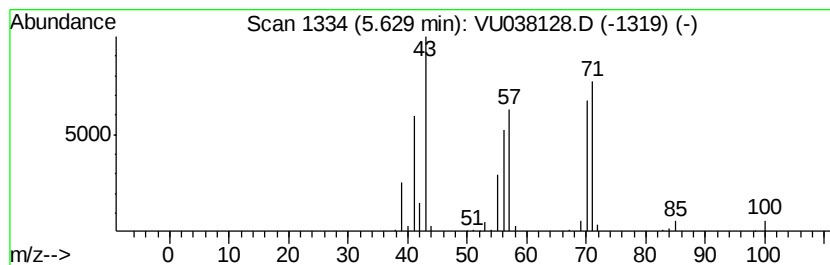
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 29

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 5.63 | 0.79 ug/L | 138915 | 1,4-Difluorobenzene | 6.28 |

| Hit# of | 5                      | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|------------------------|--------------|-----|---------|-------------|------|
| 1       | Hexane, 3-methyl-      |              | 100 | C7H16   | 000589-34-4 | 91   |
| 2       | Hexane, 3-methyl-      |              | 100 | C7H16   | 000589-34-4 | 87   |
| 3       | Hexane, 3-methyl-      |              | 100 | C7H16   | 000589-34-4 | 72   |
| 4       | Heptane, 4-methyl-     |              | 114 | C8H18   | 000589-53-7 | 59   |
| 5       | Heptane, 3,4-dimethyl- |              | 128 | C9H20   | 000922-28-1 | 56   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
BFSP5

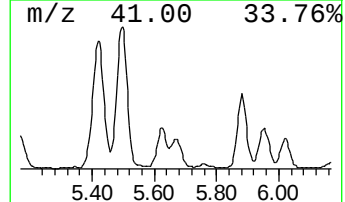
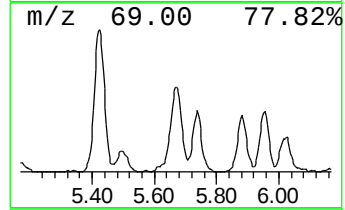
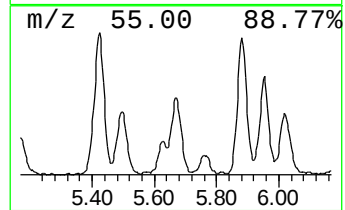
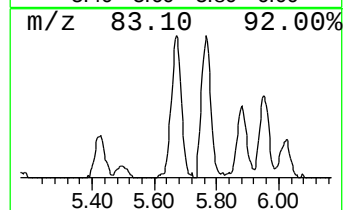
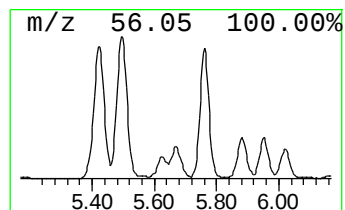
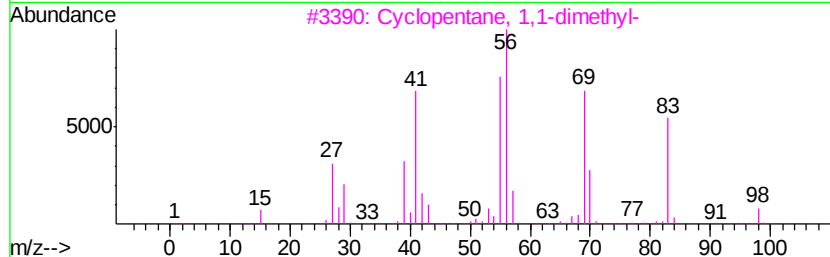
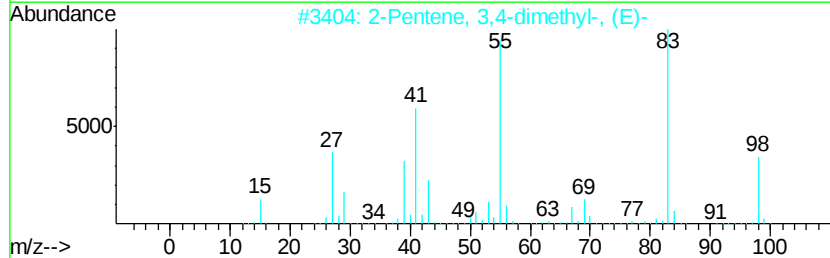
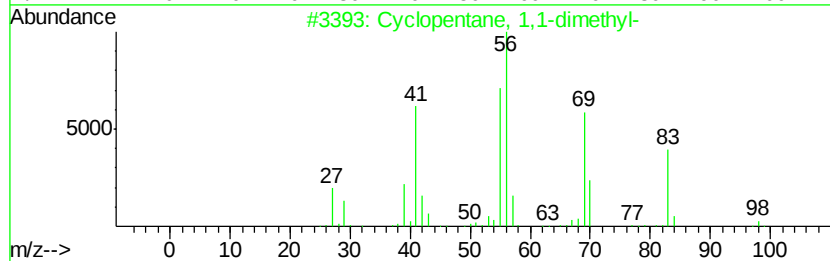
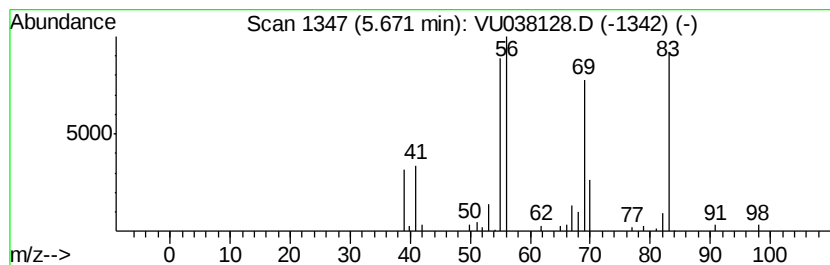
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Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 8 (DEL) Alkane: Cyclic5.67 Concentration Rank 34

| R.T. | EstConc   | Area  | Relative to ISTD    | R.T. |
|------|-----------|-------|---------------------|------|
| 5.67 | 0.57 ug/L | 99329 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopentane, 1,1-dimethyl-         | 98  | C7H14   | 001638-26-2 | 45   |
| 2    |    | 2-Pentene, 3,4-dimethyl-, (E)-      | 98  | C7H14   | 004914-92-5 | 43   |
| 3    |    | Cyclopentane, 1,1-dimethyl-         | 98  | C7H14   | 001638-26-2 | 36   |
| 4    |    | Cyclopropane, 2-bromo-1,1,3-trim... | 162 | C6H11Br | 036617-00-2 | 35   |
| 5    |    | Cyclopentane, 1-ethyl-1-methyl-     | 112 | C8H16   | 016747-50-5 | 28   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

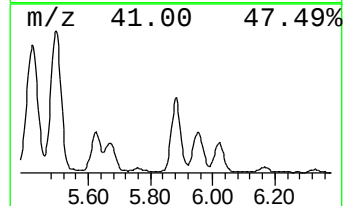
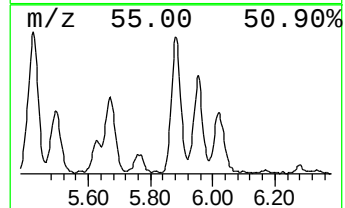
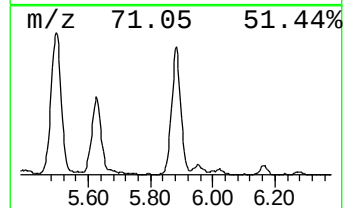
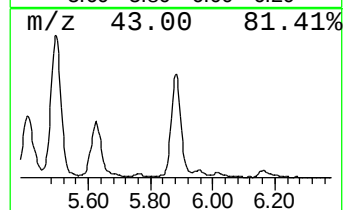
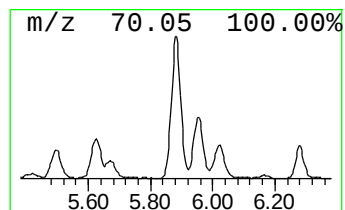
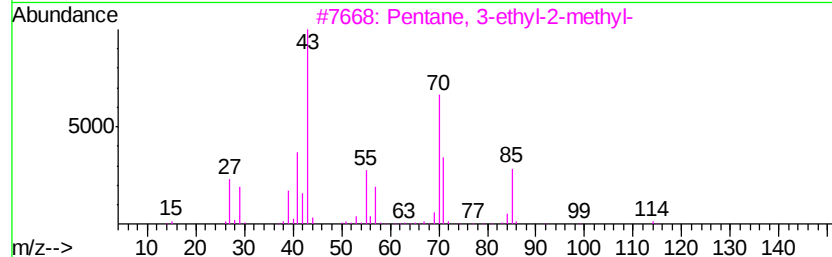
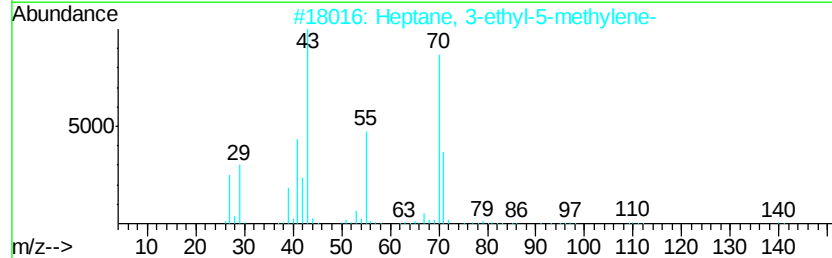
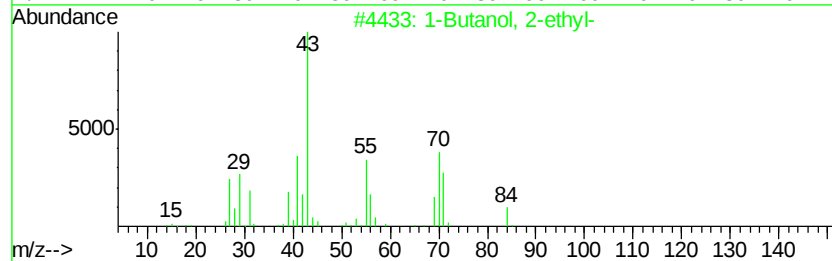
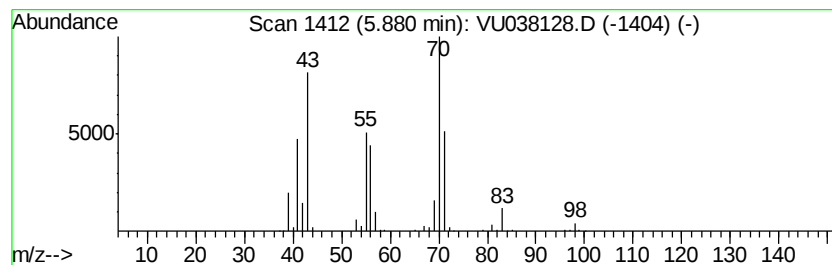
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MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 9 1-Butanol, 2-ethyl- Concentration Rank 11

| R.T.    | EstConc   | Area                          | Relative to ISTD    | R.T.    |              |      |
|---------|-----------|-------------------------------|---------------------|---------|--------------|------|
| 5.88    | 1.83 ug/L | 321673                        | 1,4-Difluorobenzene | 6.28    |              |      |
| Hit# of | 5         | Tentative ID                  | MW                  | MolForm | CAS#         | Qual |
| 1       |           | 1-Butanol, 2-ethyl-           | 102                 | C6H14O  | 000097-95-0  | 50   |
| 2       |           | Heptane, 3-ethyl-5-methylene- | 140                 | C10H20  | 1000160-41-7 | 50   |
| 3       |           | Pentane, 3-ethyl-2-methyl-    | 114                 | C8H18   | 000609-26-7  | 45   |
| 4       |           | Tridecane, 3-methylene-       | 196                 | C14H28  | 019780-34-8  | 40   |
| 5       |           | 1-Pentanol, 3,4-dimethyl-     | 116                 | C7H16O  | 006570-87-2  | 38   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

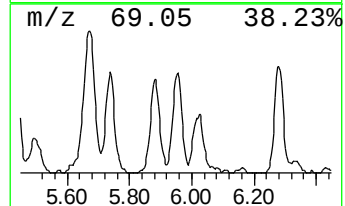
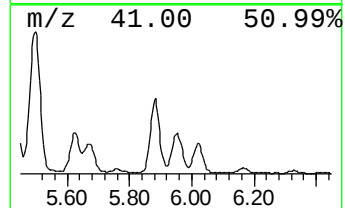
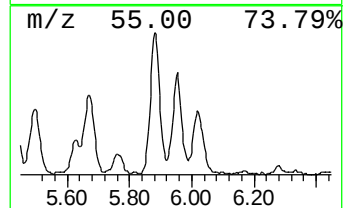
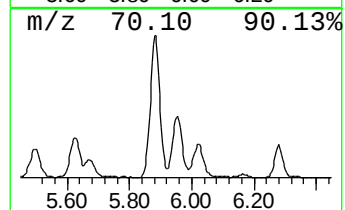
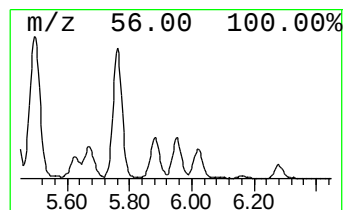
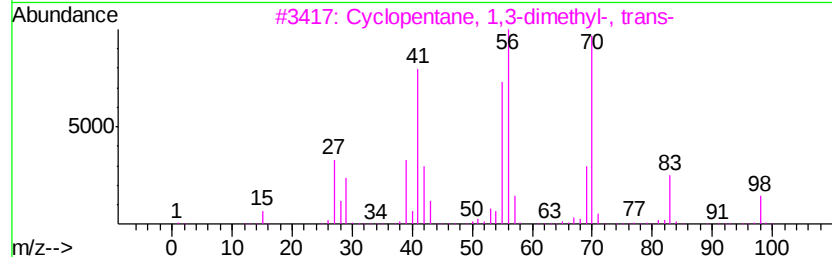
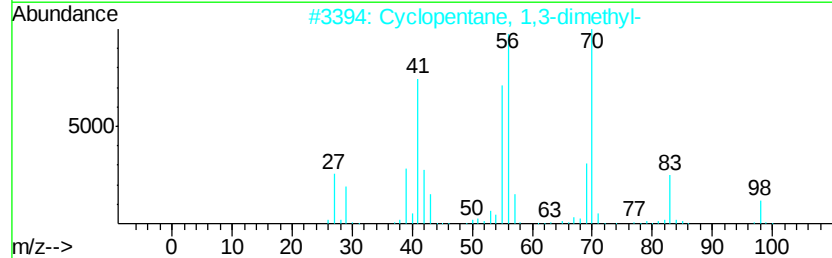
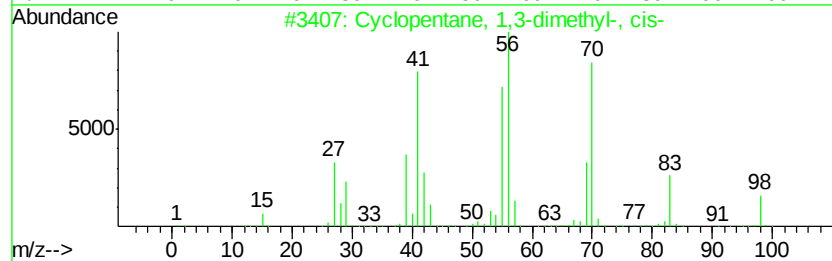
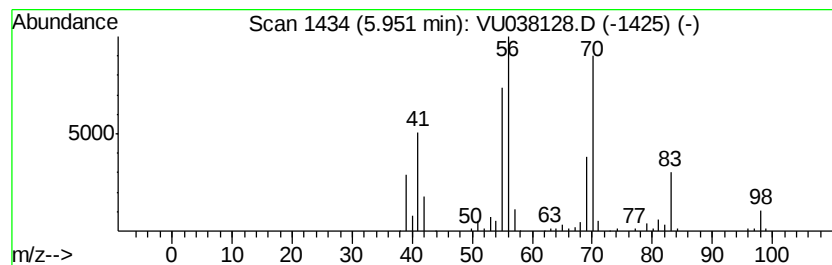
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Cyclic5.95 Concentration Rank 25

| R.T. | EstConc   | Area   | Relative to ISTD    | R.T. |
|------|-----------|--------|---------------------|------|
| 5.95 | 0.93 ug/L | 163870 | 1,4-Difluorobenzene | 6.28 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopentane, 1,3-dimethyl-, cis-   | 98  | C7H14   | 002532-58-3 | 91   |
| 2    |    | Cyclopentane, 1,3-dimethyl-         | 98  | C7H14   | 002453-00-1 | 87   |
| 3    |    | Cyclopentane, 1,3-dimethyl-, trans- | 98  | C7H14   | 001759-58-6 | 78   |
| 4    |    | Cyclobutanone, 3-ethyl-             | 98  | C6H10O  | 056335-73-0 | 64   |
| 5    |    | 2H-Pyran-2,6(3H)-dione, dihydro-... | 142 | C7H10O3 | 004160-82-1 | 59   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

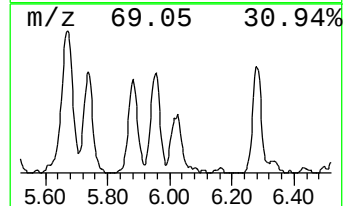
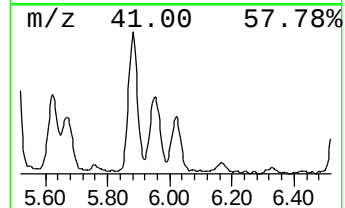
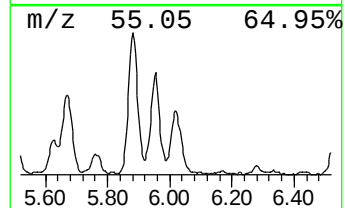
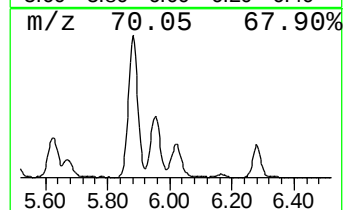
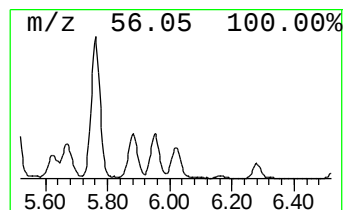
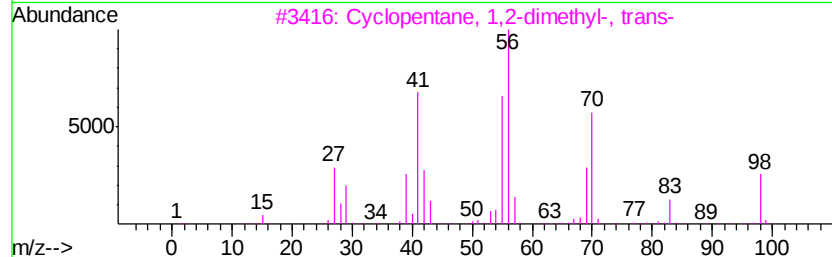
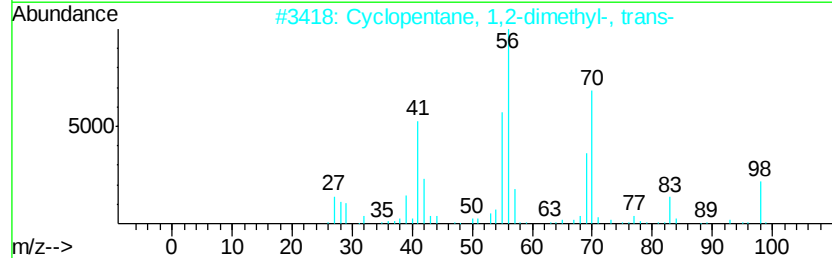
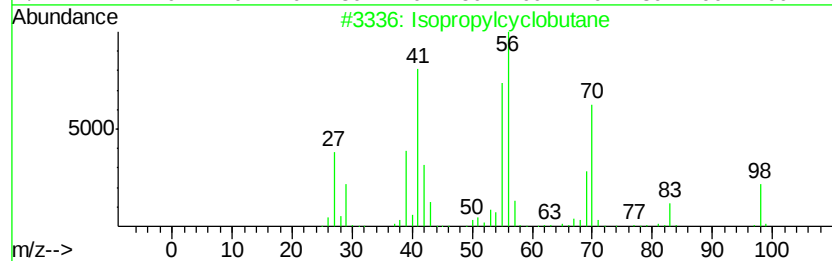
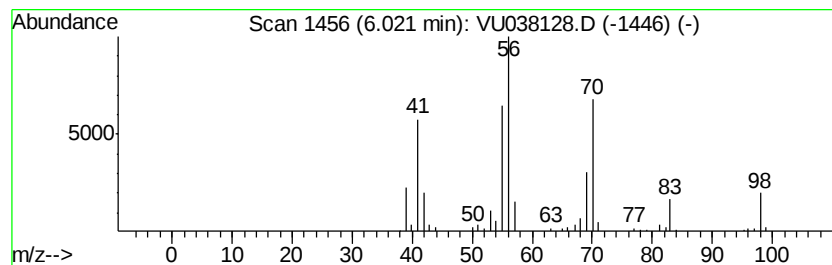
Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.    | EstConc   | Area                                | Relative to ISTD    | R.T.           |
|---------|-----------|-------------------------------------|---------------------|----------------|
| 6.02    | 0.59 ug/L | 103789                              | 1,4-Difluorobenzene | 6.28           |
| Hit# of | 5         | Tentative ID                        | MW MolForm          | CAS# Qual      |
| 1       |           | Isopropylcyclobutane                | 98 C7H14            | 000872-56-0 93 |
| 2       |           | Cyclopentane, 1,2-dimethyl-, trans- | 98 C7H14            | 000822-50-4 91 |
| 3       |           | Cyclopentane, 1,2-dimethyl-, trans- | 98 C7H14            | 000822-50-4 91 |
| 4       |           | Cyclopentane, 1,2-dimethyl-         | 98 C7H14            | 002452-99-5 91 |
| 5       |           | Cyclopentane, 1,2-dimethyl-, cis-   | 98 C7H14            | 001192-18-3 76 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

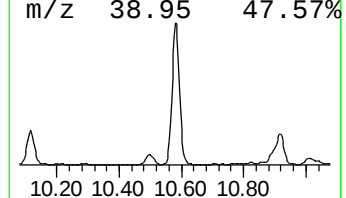
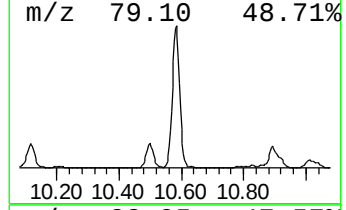
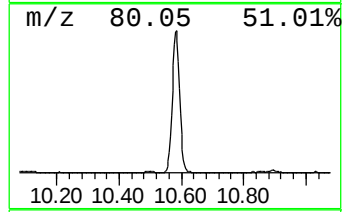
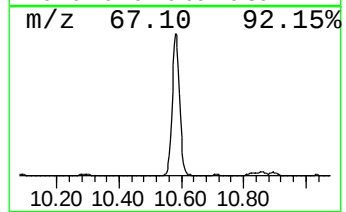
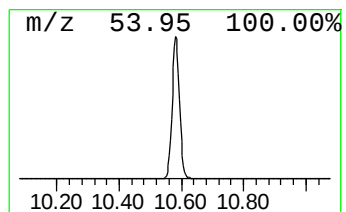
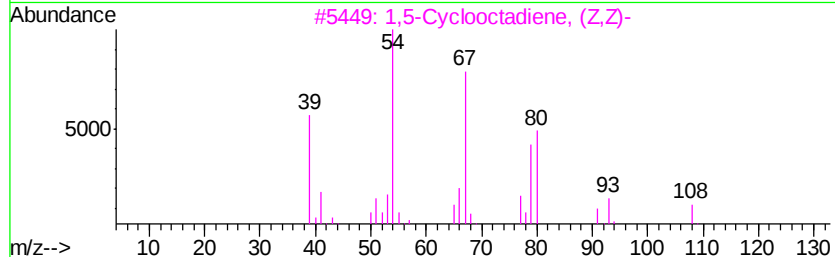
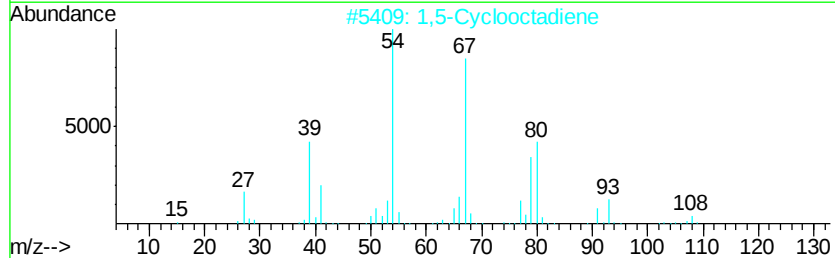
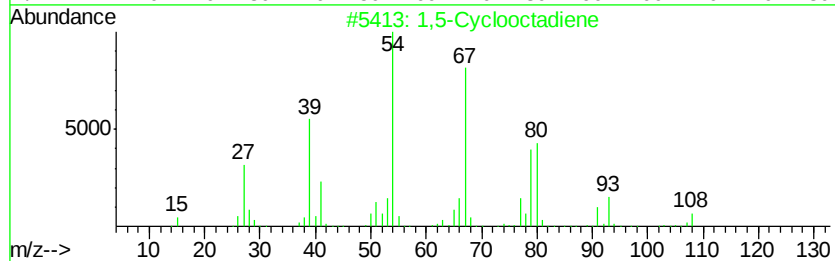
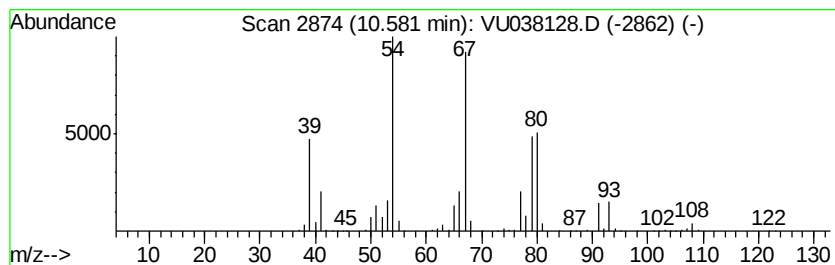
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 1,5-Cyclooctadiene Concentration Rank 8

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T. |
|-------|-----------|---------|------------------|------|
| 10.58 | 2.06 ug/L | 1427720 | Chlorobenzene-d5 | 9.44 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | 1,5-Cyclooctadiene         | 108 | C8H12   | 000111-78-4 | 95   |
| 2    |    |   | 1,5-Cyclooctadiene         | 108 | C8H12   | 000111-78-4 | 94   |
| 3    |    |   | 1,5-Cyclooctadiene, (Z,Z)- | 108 | C8H12   | 001552-12-1 | 90   |
| 4    |    |   | 4-Cyanocyclohexene         | 107 | C7H9N   | 000100-45-8 | 83   |
| 5    |    |   | 1,5-Cyclooctadiene, (E,Z)- | 108 | C8H12   | 005259-71-2 | 56   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

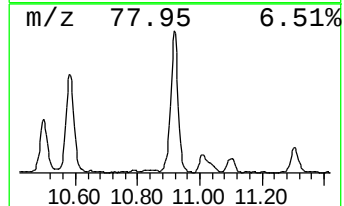
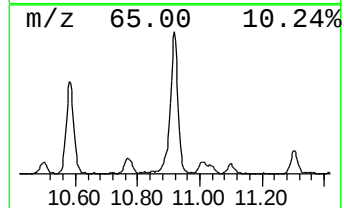
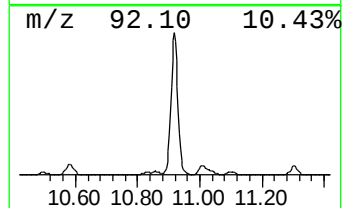
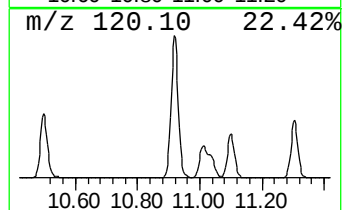
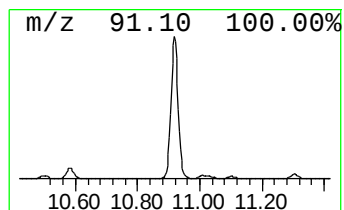
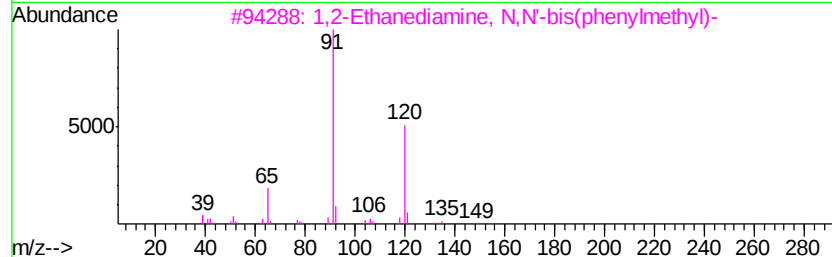
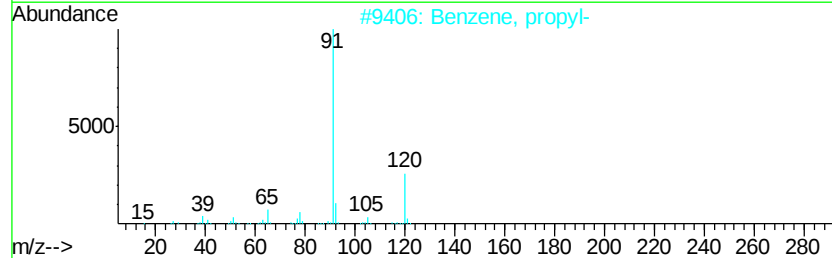
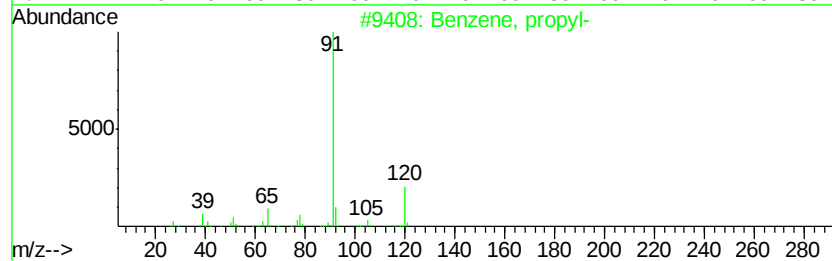
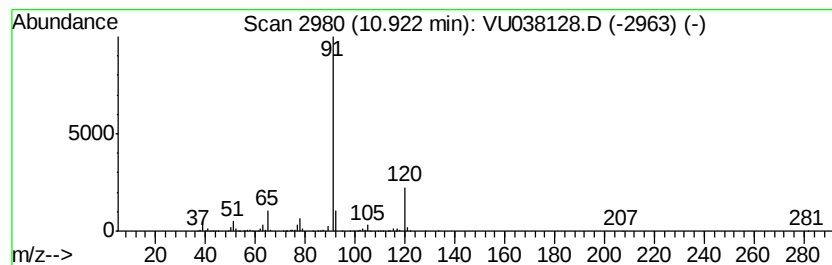
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Benzene, propyl- Concentration Rank 3

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 10.92 | 4.58 ug/L | 1264390 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Benzene, propyl-                    | 120 | C9H12    | 000103-65-1 | 90   |
| 2    |    |   | Benzene, propyl-                    | 120 | C9H12    | 000103-65-1 | 87   |
| 3    |    |   | 1,2-Ethanediamine, N,N'-bis(phen... | 240 | C16H20N2 | 000140-28-3 | 72   |
| 4    |    |   | Propanenitrile, 3-[(phenylmethyl... | 160 | C10H12N2 | 000706-03-6 | 72   |
| 5    |    |   | N-Benzyl-2-phenethylamine           | 211 | C15H17N  | 003647-71-0 | 59   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

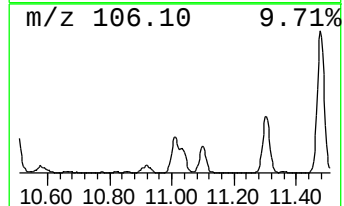
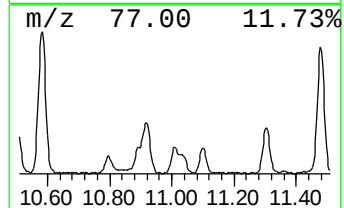
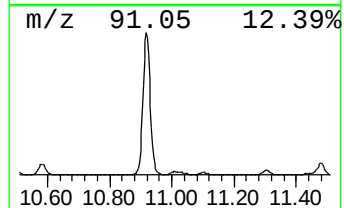
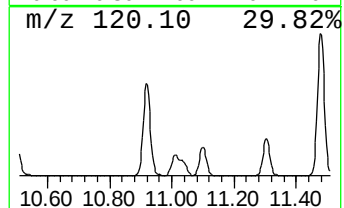
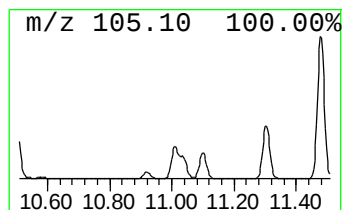
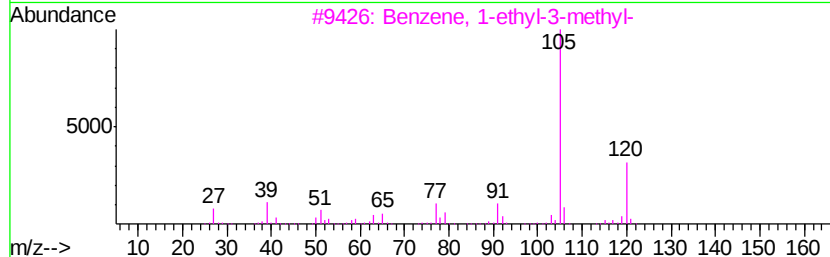
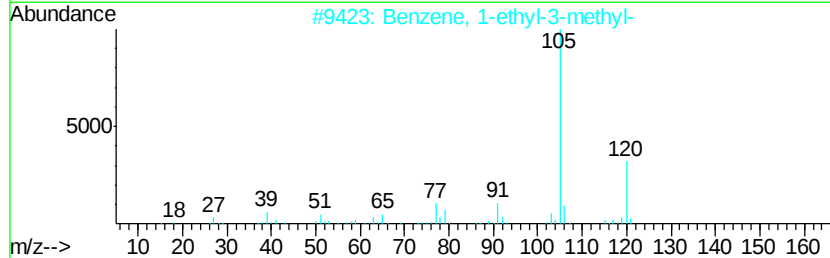
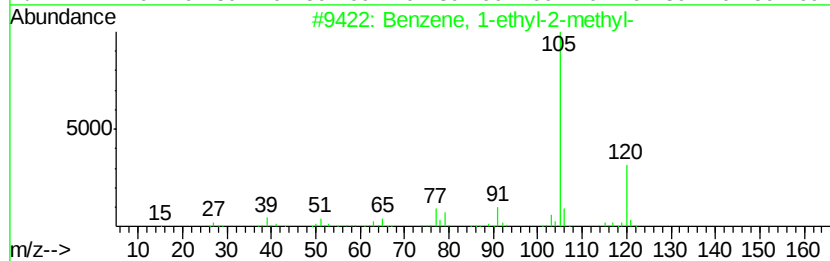
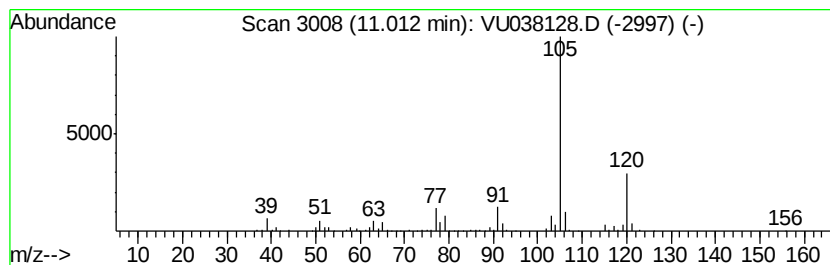
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Benzene, 1-ethyl-2-methyl- Concentration Rank 27

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 11.01 | 0.85 ug/L | 233201 | 1,4-Dichlorobenzene-d4 | 11.82 |

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

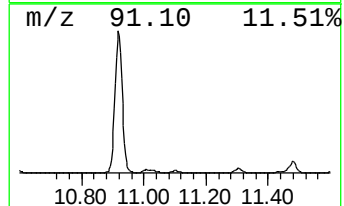
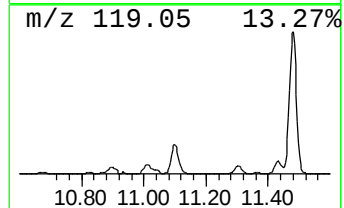
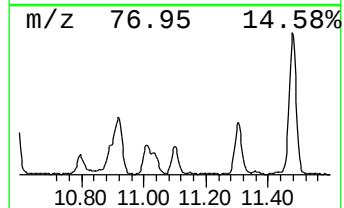
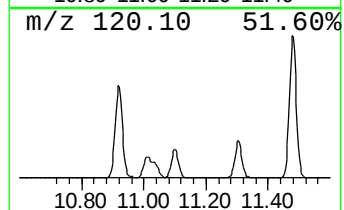
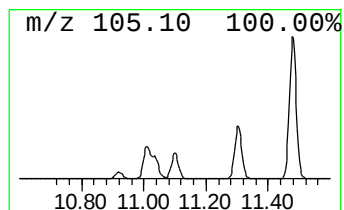
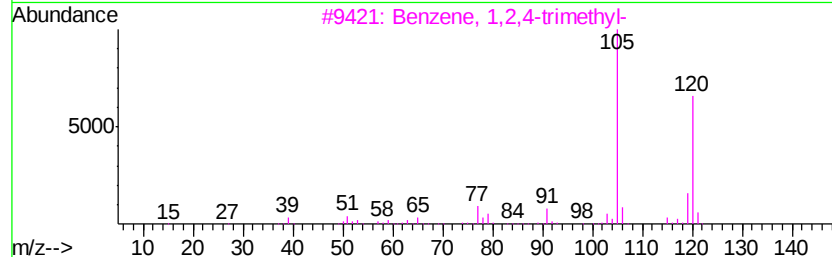
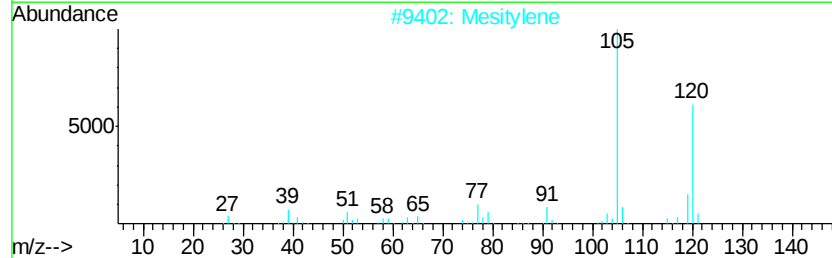
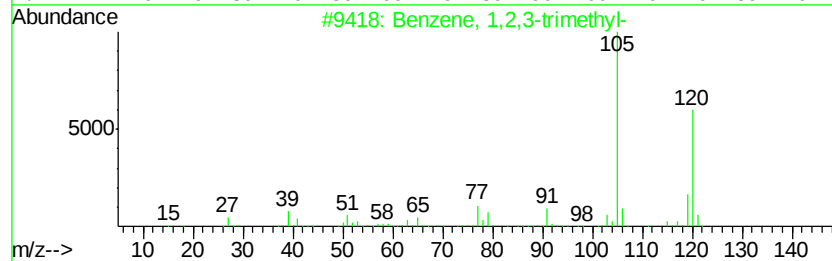
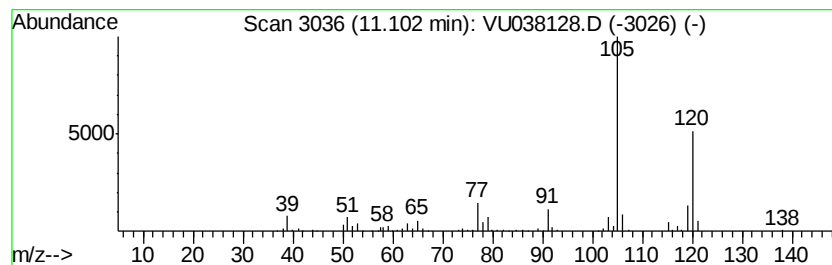
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Benzene, 1,2,3-trimethyl- Concentration Rank 31

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 11.10 | 0.66 ug/L | 182810 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 97   |
| 3    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 95   |
| 4    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 95   |
| 5    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 95   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

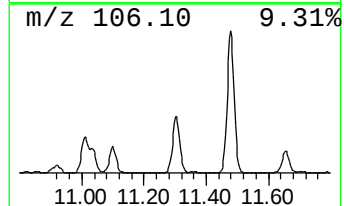
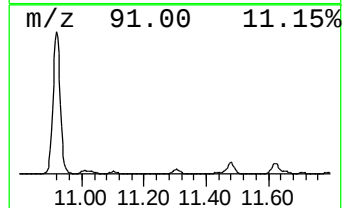
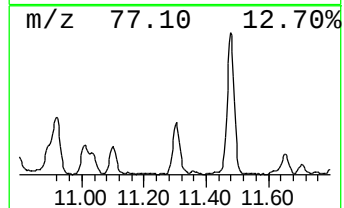
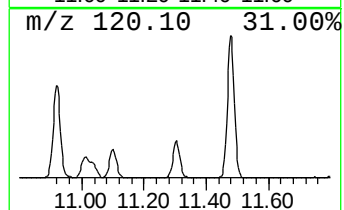
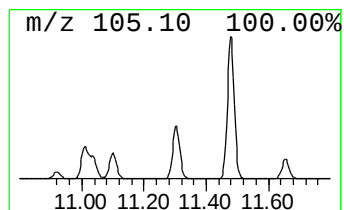
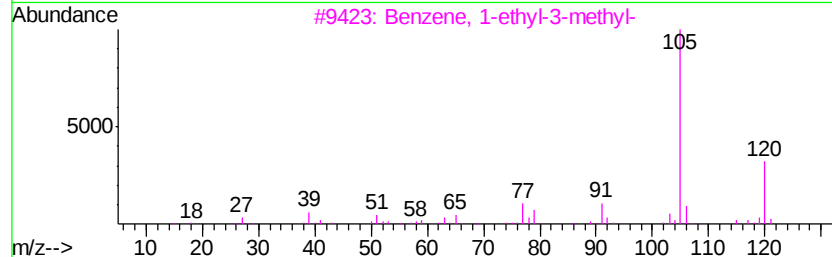
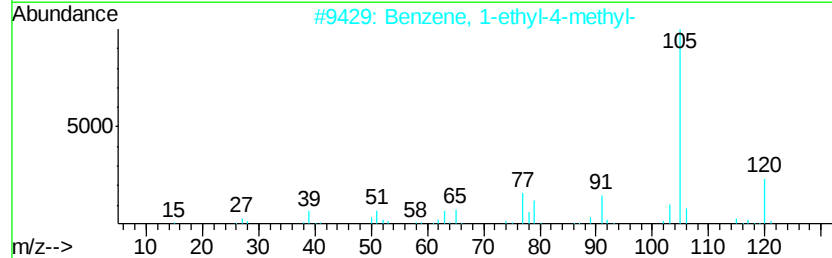
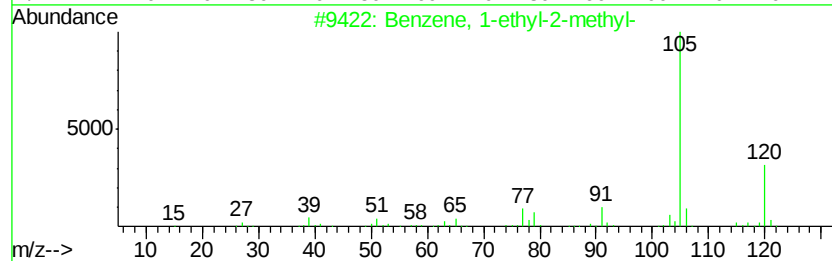
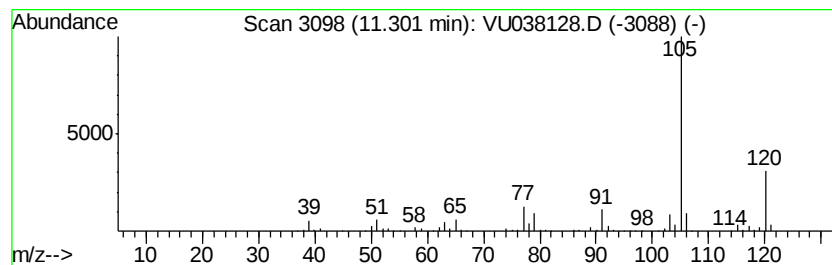
Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

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

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Peak Number 17 Benzene, 1-ethyl-4-methyl- Concentration Rank 21

| R.T.    | EstConc   | Area                       | Relative to ISTD       | R.T.           |
|---------|-----------|----------------------------|------------------------|----------------|
| 11.30   | 1.23 ug/L | 338013                     | 1,4-Dichlorobenzene-d4 | 11.82          |
| Hit# of | 5         | Tentative ID               | MW MolForm             | CAS# Qual      |
| 1       |           | Benzene, 1-ethyl-2-methyl- | 120 C9H12              | 000611-14-3 95 |
| 2       |           | Benzene, 1-ethyl-4-methyl- | 120 C9H12              | 000622-96-8 94 |
| 3       |           | Benzene, 1-ethyl-3-methyl- | 120 C9H12              | 000620-14-4 91 |
| 4       |           | Benzene, 1,2,3-trimethyl-  | 120 C9H12              | 000526-73-8 91 |
| 5       |           | Mesitylene                 | 120 C9H12              | 000108-67-8 91 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

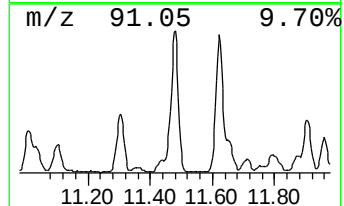
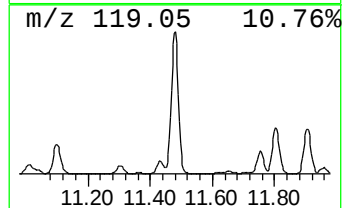
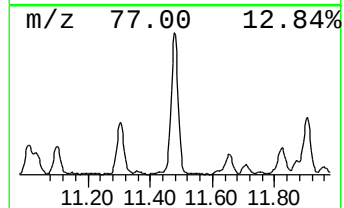
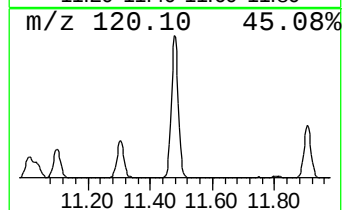
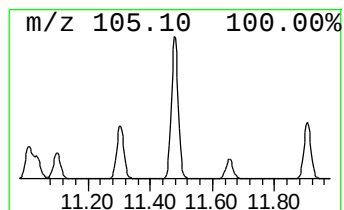
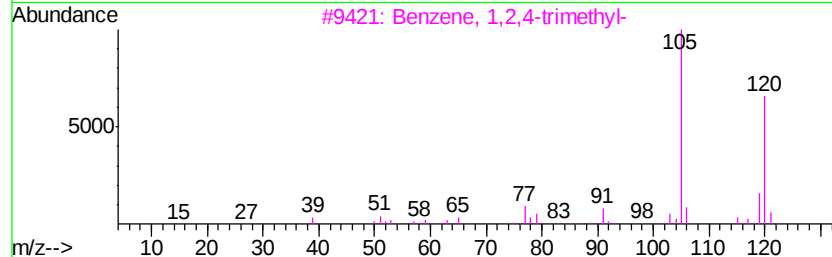
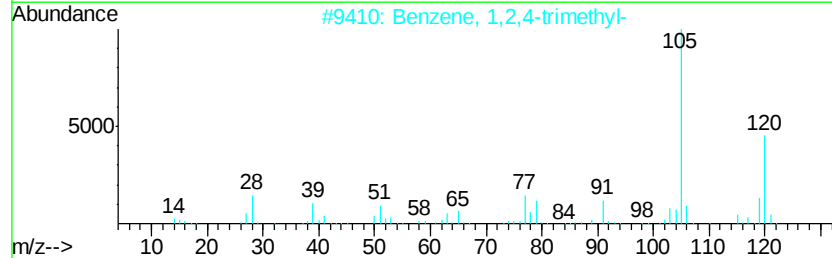
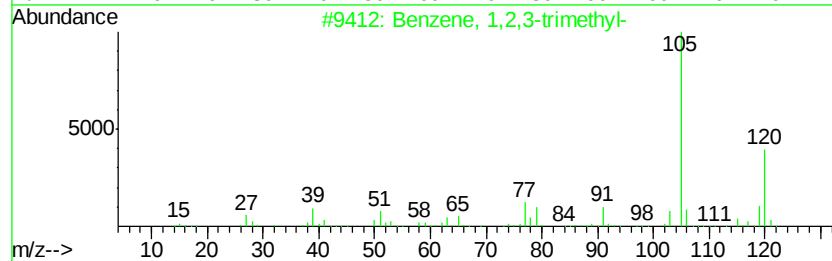
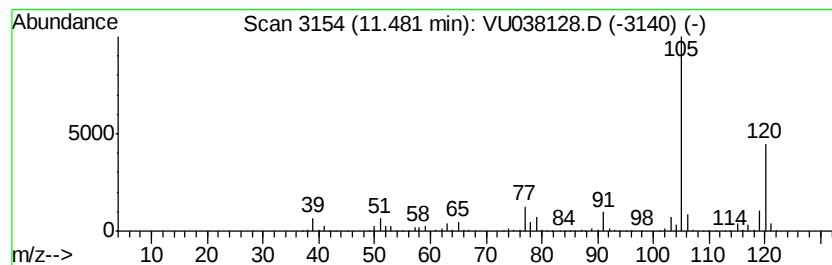
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Benzene, 1,2,4-trimethyl- Concentration Rank 4

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 11.48 | 3.73 ug/L | 1029480 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 94   |
| 2    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 3    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 4    |    | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 94   |
| 5    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

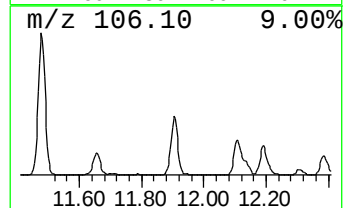
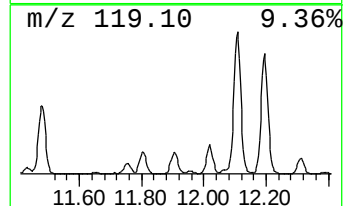
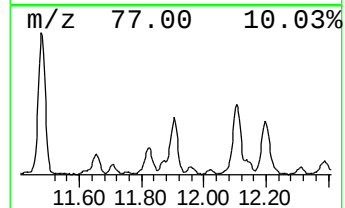
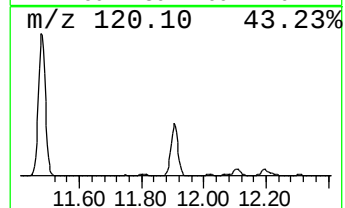
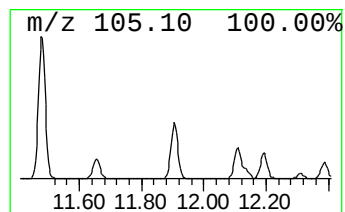
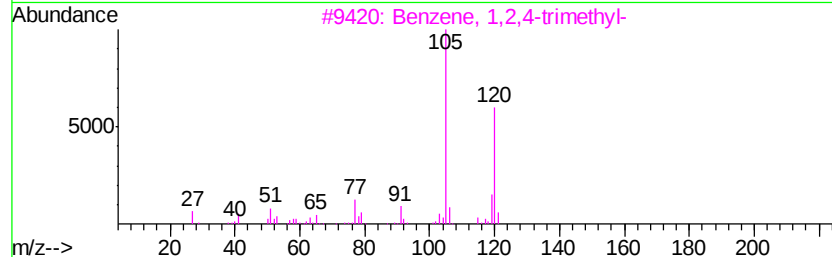
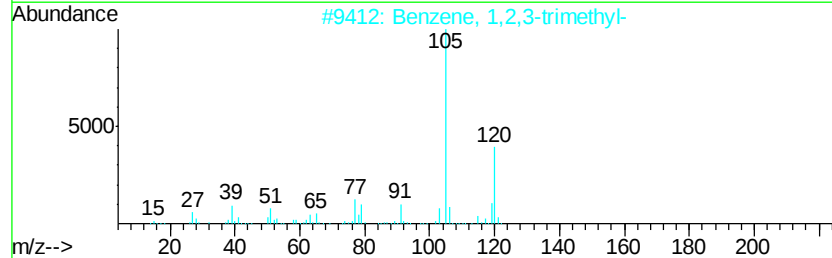
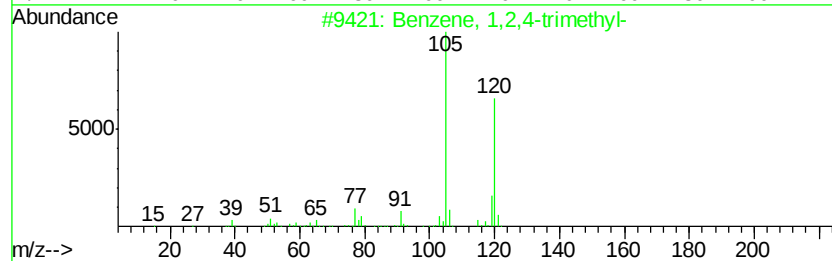
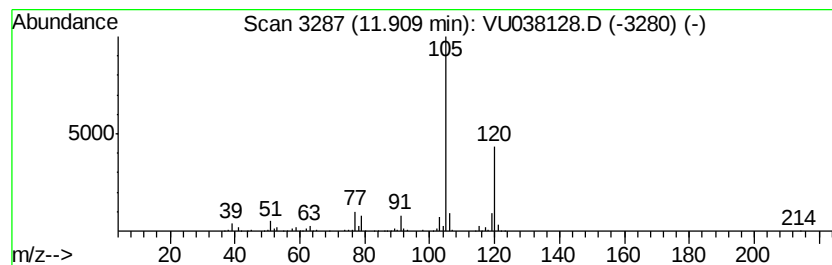
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Mesitylene Concentration Rank 19

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 11.91 | 1.28 ug/L | 352900 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 2    |    | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 91   |
| 3    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 91   |
| 4    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 91   |
| 5    |    | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

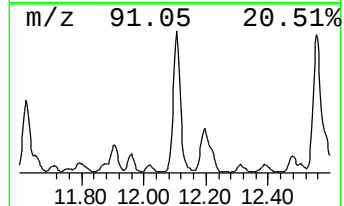
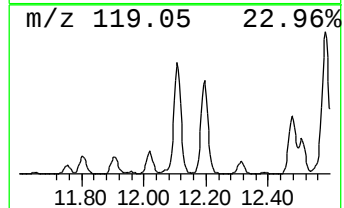
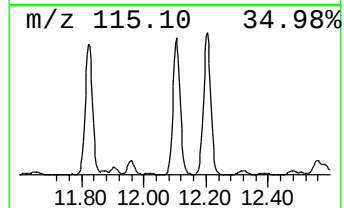
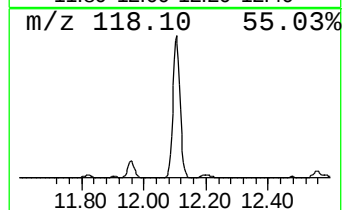
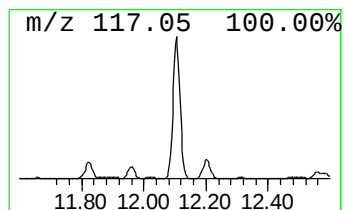
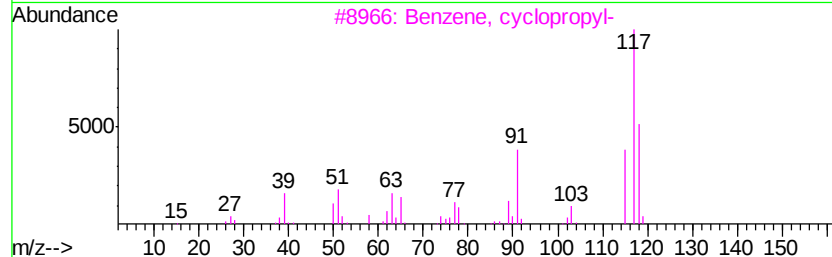
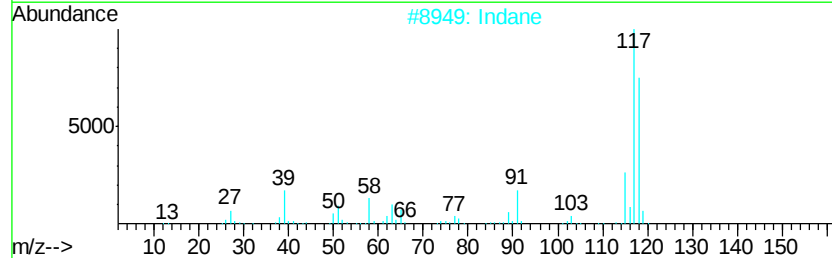
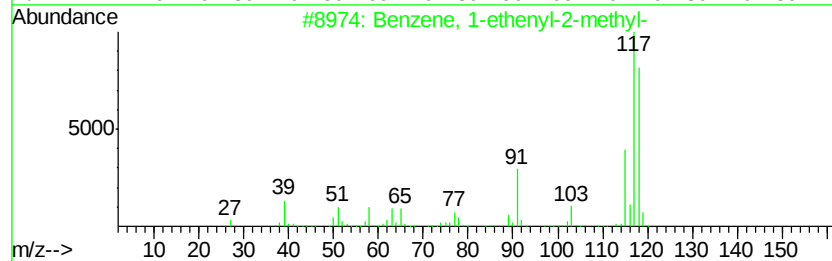
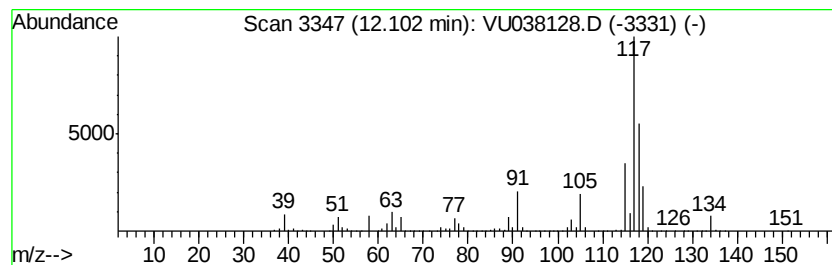
Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

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

\*\*\*\*\*  
Peak Number 20 Indane Concentration Rank 2

| R.T.    | EstConc   | Area                         | Relative to ISTD       | R.T.           |
|---------|-----------|------------------------------|------------------------|----------------|
| 12.10   | 5.06 ug/L | 1395310                      | 1,4-Dichlorobenzene-d4 | 11.82          |
| Hit# of | 5         | Tentative ID                 | MW MolForm             | CAS# Qual      |
| 1       |           | Benzene, 1-ethenyl-2-methyl- | 118 C9H10              | 000611-15-4 86 |
| 2       |           | Indane                       | 118 C9H10              | 000496-11-7 70 |
| 3       |           | Benzene, cyclopropyl-        | 118 C9H10              | 000873-49-4 64 |
| 4       |           | Benzene, cyclopropyl-        | 118 C9H10              | 000873-49-4 64 |
| 5       |           | Benzene, 2-propenyl-         | 118 C9H10              | 000300-57-2 64 |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

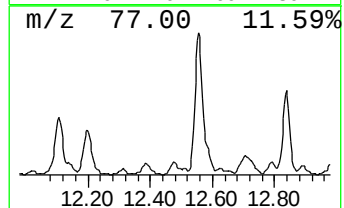
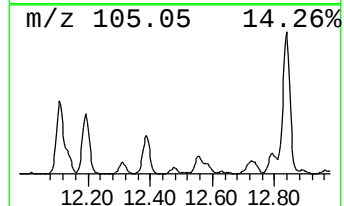
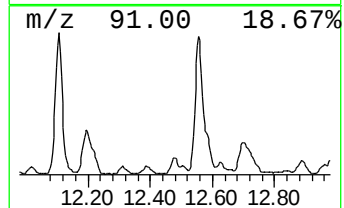
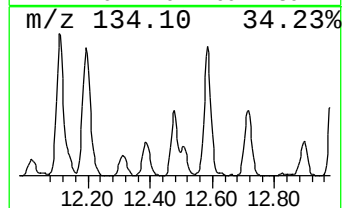
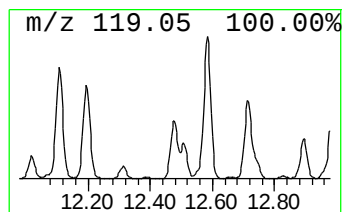
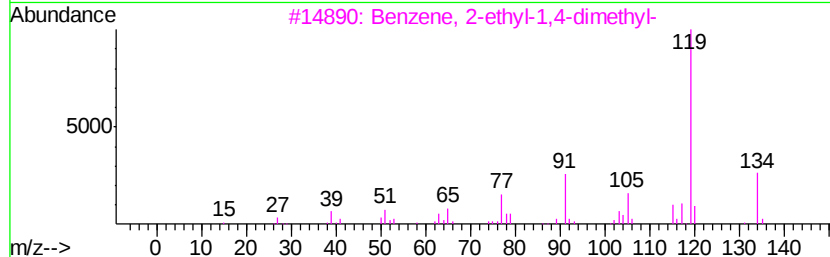
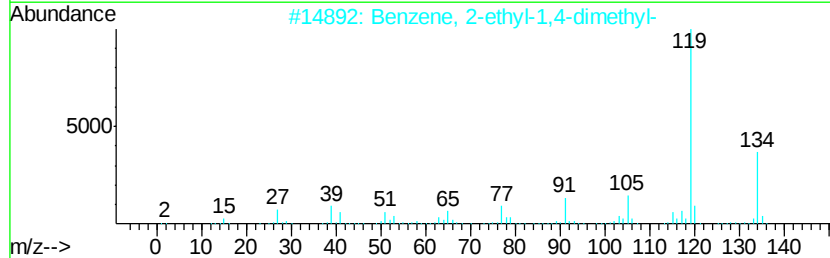
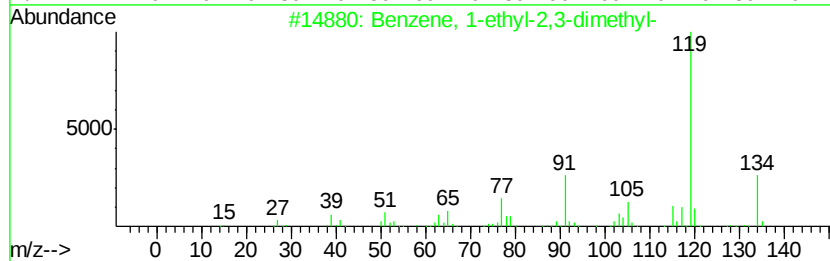
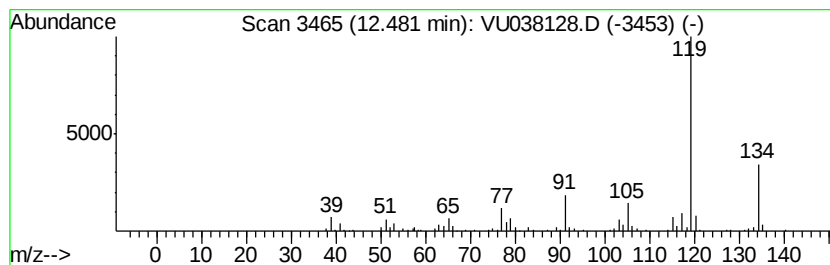
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 37

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.48 | 0.53 ug/L | 146874 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 96   |
| 2    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 95   |
| 3    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 95   |
| 4    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 94   |
| 5    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

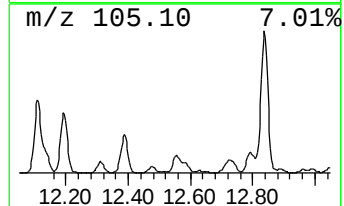
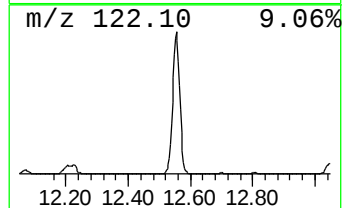
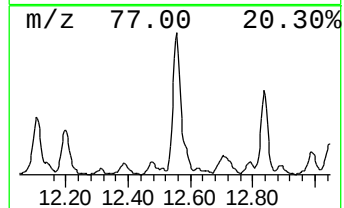
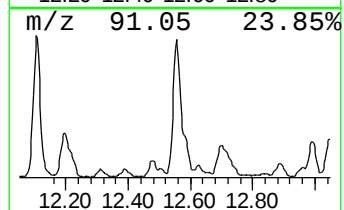
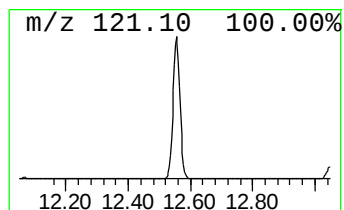
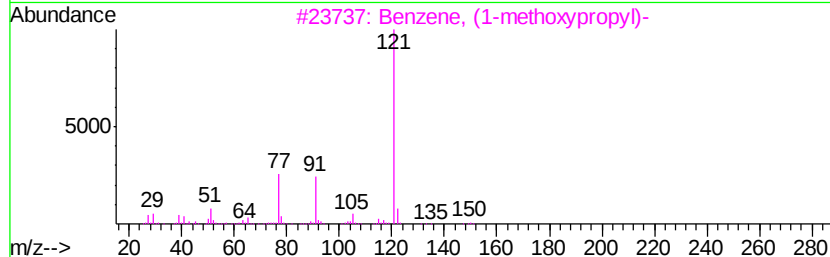
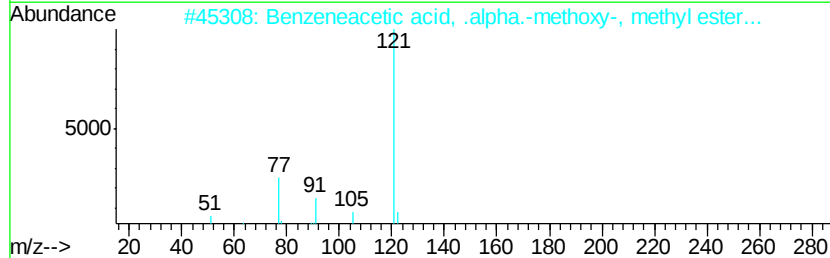
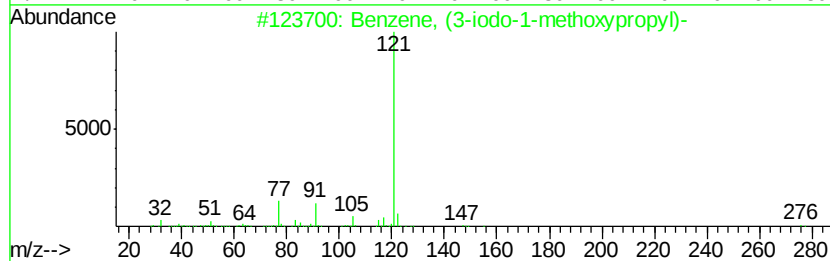
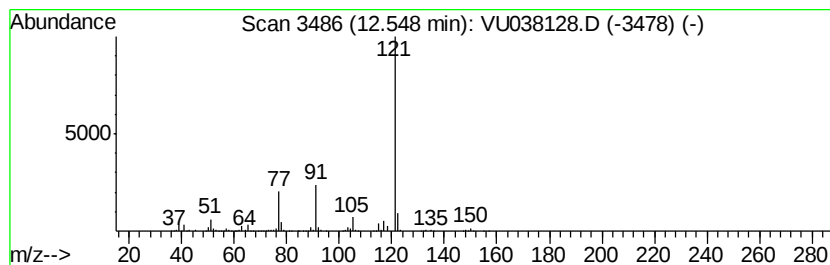
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 Benzene, (3-iodo-1-methoxyp... Concentration Rank 7

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.55 | 2.44 ug/L | 673023 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Benzene, (3-iodo-1-methoxypropyl)-  | 276 | C10H13IO | 1000327-31-9 | 72   |
| 2    |    | Benzeneacetic acid, .alpha.-meth... | 180 | C10H12O3 | 056143-21-6  | 72   |
| 3    |    | Benzene, (1-methoxypropyl)-         | 150 | C10H14O  | 059588-12-4  | 68   |
| 4    |    | Benzene, (1,2-dimethoxyethyl)-      | 166 | C10H14O2 | 004013-37-0  | 64   |
| 5    |    | Benzene, (3-iodo-1-methoxybutyl)-   | 290 | C11H15IO | 1000327-38-8 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

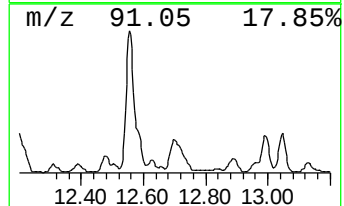
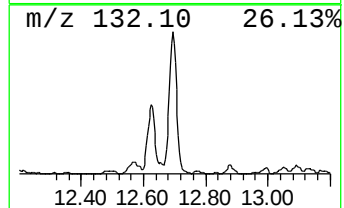
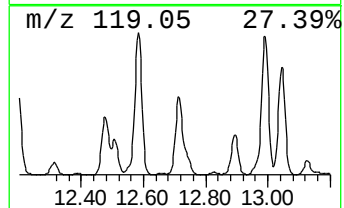
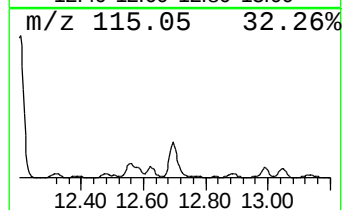
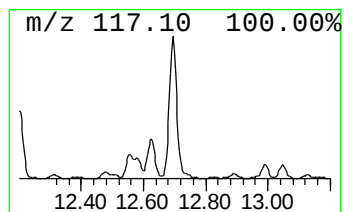
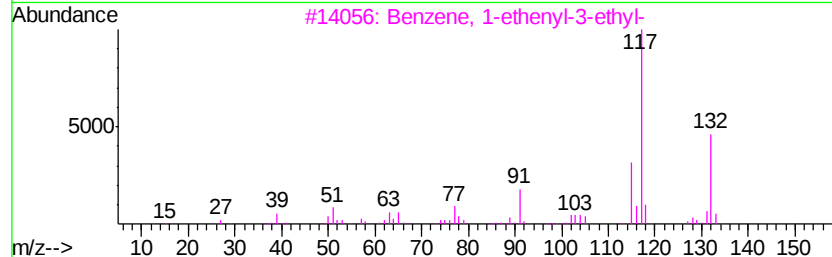
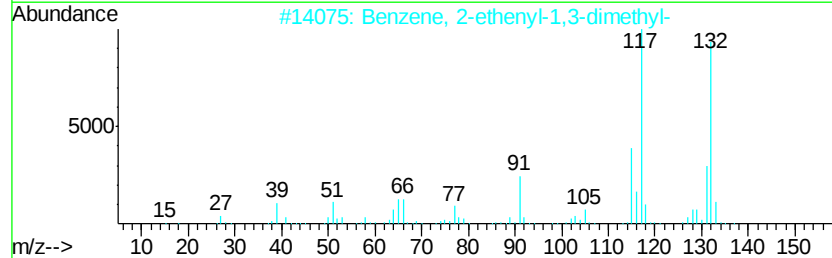
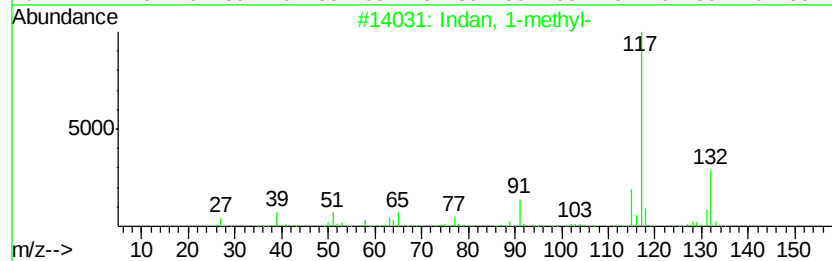
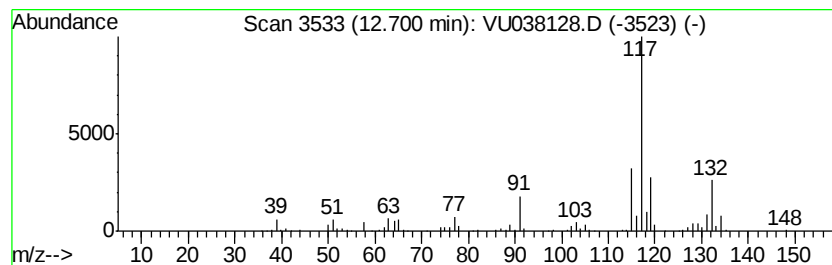
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 Indan, 1-methyl- Concentration Rank 12

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.70 | 1.60 ug/L | 442179 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Indan, 1-methyl-                    | 132 | C10H12  | 000767-58-8 | 93   |
| 2    |    | Benzene, 2-ethenyl-1,3-dimethyl-    | 132 | C10H12  | 002039-90-9 | 87   |
| 3    |    | Benzene, 1-ethenyl-3-ethyl-         | 132 | C10H12  | 007525-62-4 | 83   |
| 4    |    | Benzene, 1-ethenyl-4-ethyl-         | 132 | C10H12  | 003454-07-7 | 83   |
| 5    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 80   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

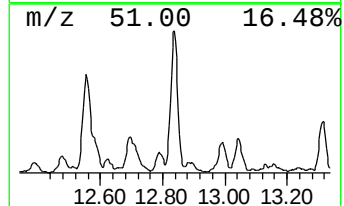
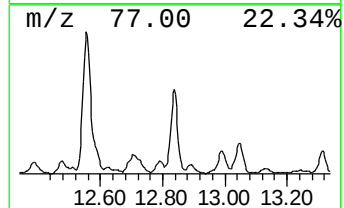
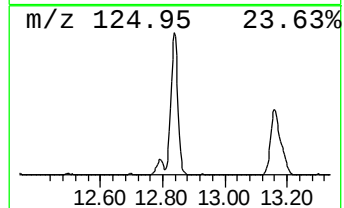
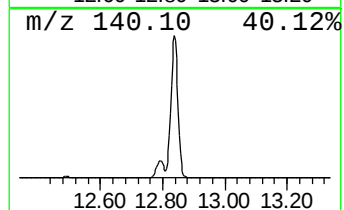
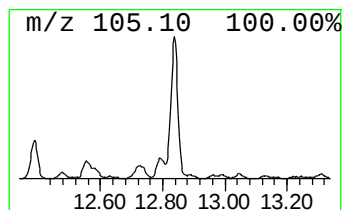
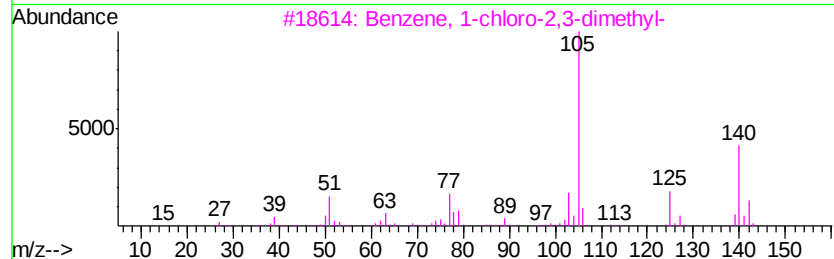
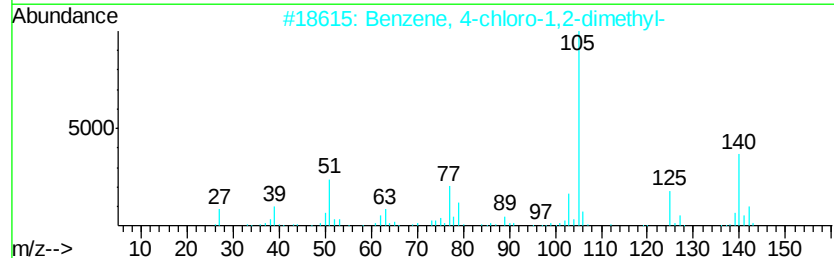
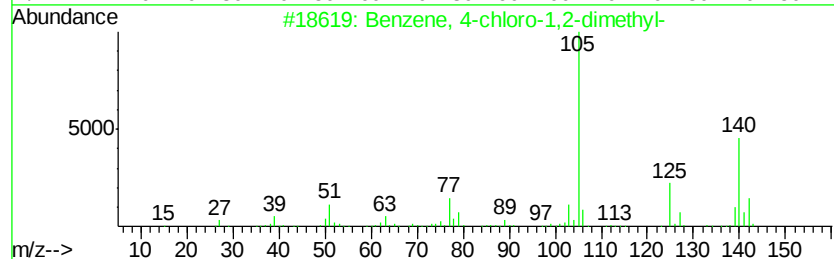
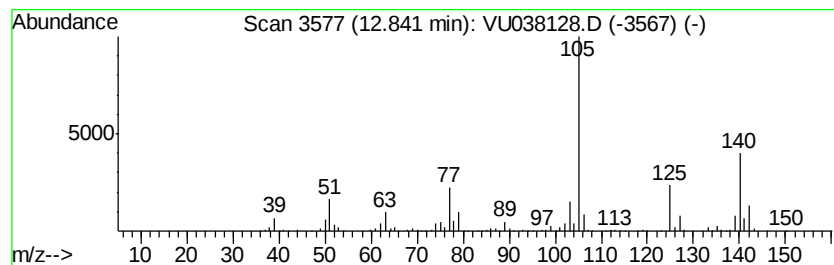
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 Benzene, 4-chloro-1,2-dimet... Concentration Rank 9

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.84 | 1.94 ug/L | 536406 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-chloro-1,2-dimethyl- | 140 | C8H9Cl  | 000615-60-1 | 95   |
| 2    |    | Benzene, 4-chloro-1,2-dimethyl- | 140 | C8H9Cl  | 000615-60-1 | 95   |
| 3    |    | Benzene, 1-chloro-2,3-dimethyl- | 140 | C8H9Cl  | 000608-23-1 | 94   |
| 4    |    | Benzene, 2-chloro-1,4-dimethyl- | 140 | C8H9Cl  | 000095-72-7 | 94   |
| 5    |    | Benzene, 4-chloro-1,2-dimethyl- | 140 | C8H9Cl  | 000615-60-1 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

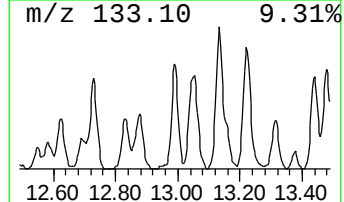
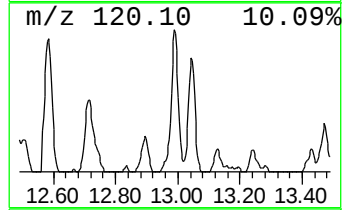
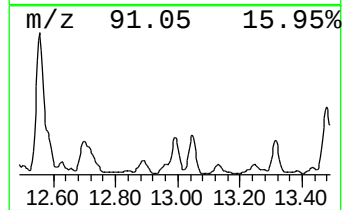
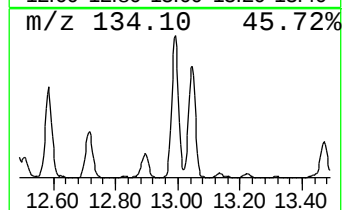
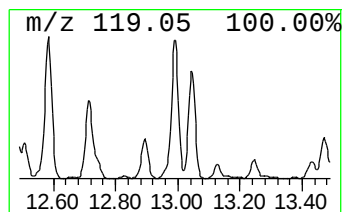
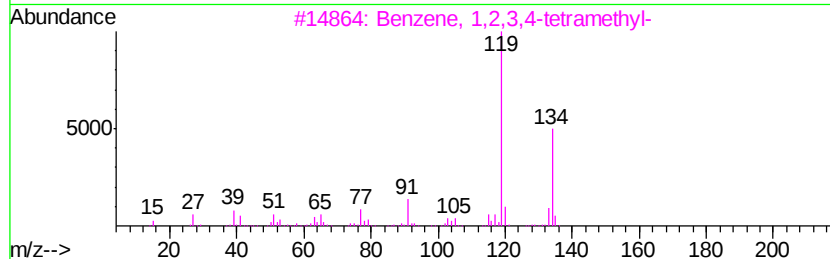
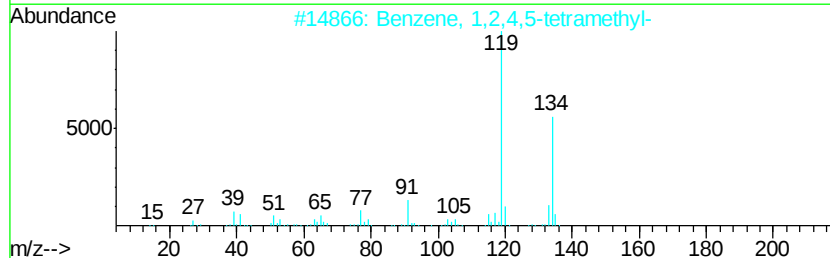
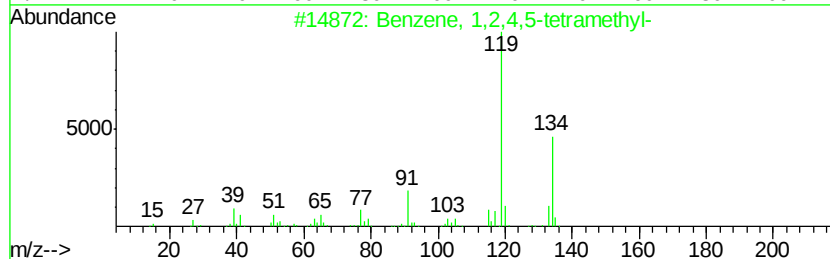
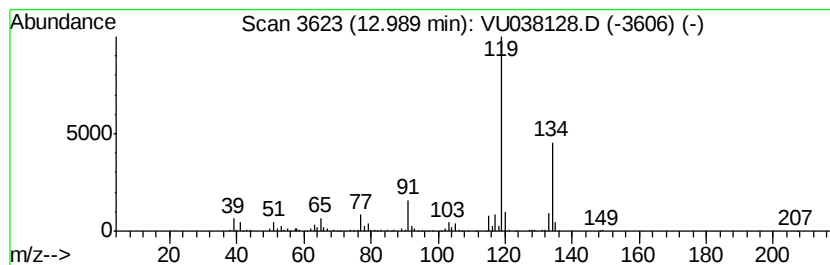
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 20

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.99 | 1.26 ug/L | 347564 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 97   |
| 2    |    | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 95   |
| 3    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 95   |
| 4    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 95   |
| 5    |    | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 94   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

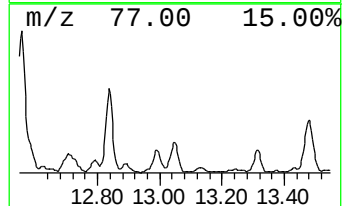
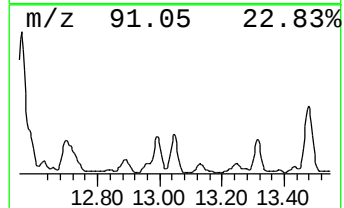
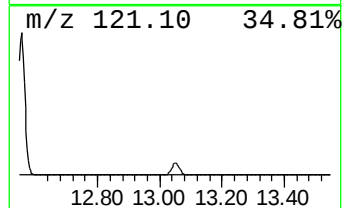
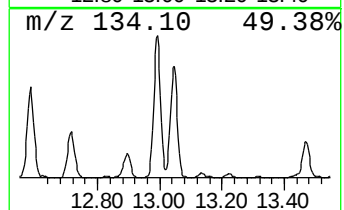
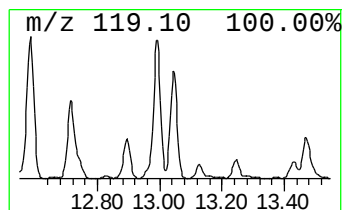
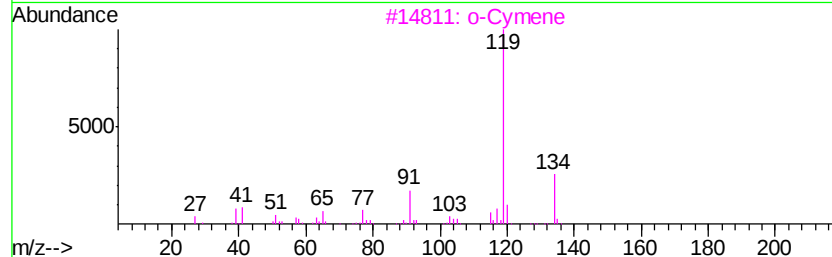
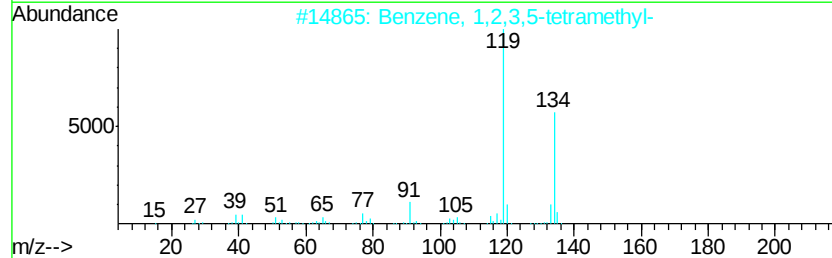
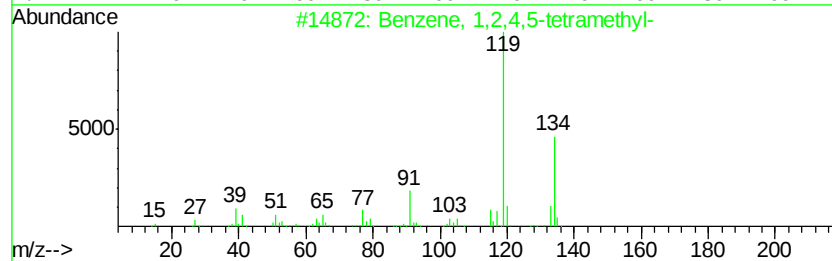
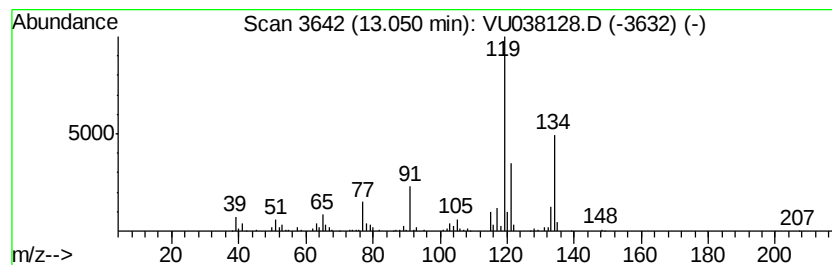
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 23

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 13.05 | 1.12 ug/L | 309281 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 95   |
| 2    |    | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 94   |
| 3    |    | o-Cymene                      | 134 | C10H14  | 000527-84-4 | 94   |
| 4    |    | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 94   |
| 5    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 93   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

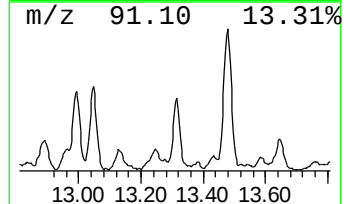
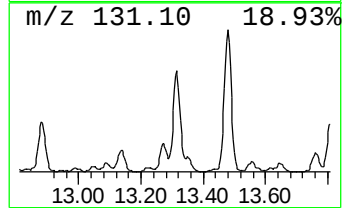
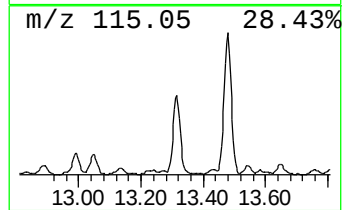
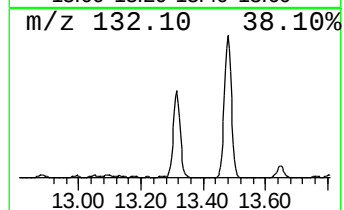
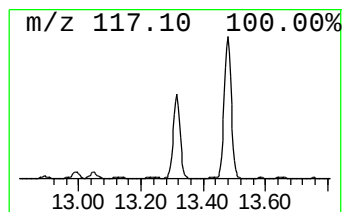
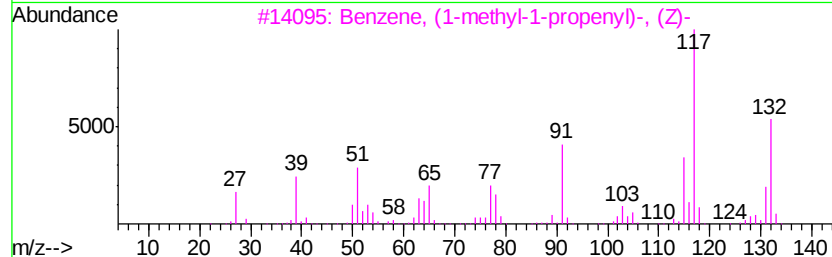
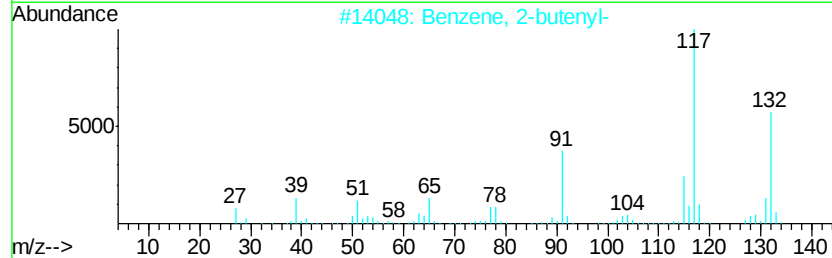
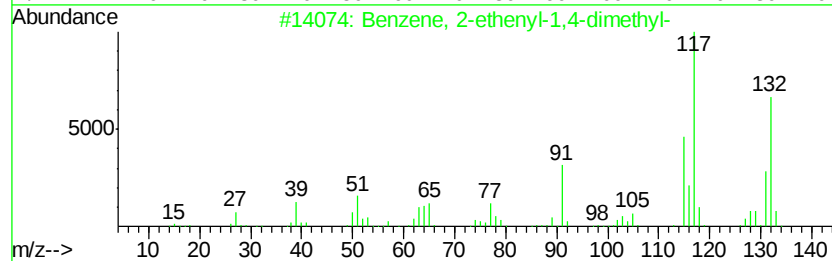
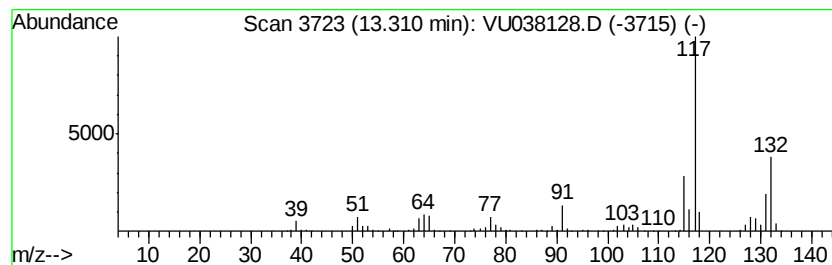
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 18

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 13.31 | 1.32 ug/L | 365162 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethenyl-1,4-dimethyl-    | 132 | C10H12  | 002039-89-6 | 92   |
| 2    |    |   | Benzene, 2-butenyl-                 | 132 | C10H12  | 001560-06-1 | 91   |
| 3    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000767-99-7 | 90   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-4-methyl-    | 132 | C10H12  | 000824-22-6 | 87   |
| 5    |    |   | Benzene, 1-methyl-2-(2-propenyl)-   | 132 | C10H12  | 001587-04-8 | 87   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

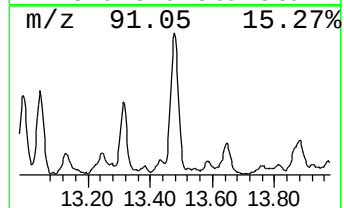
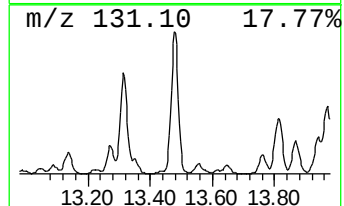
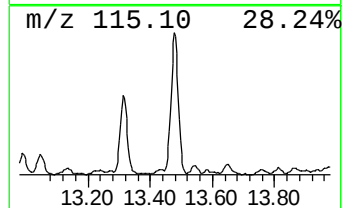
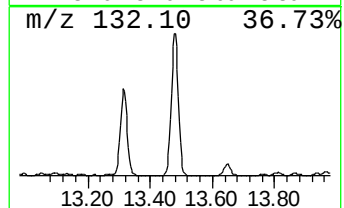
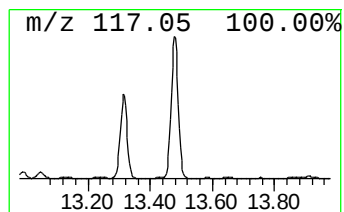
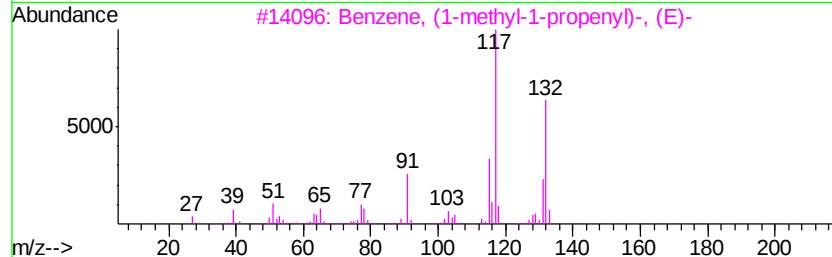
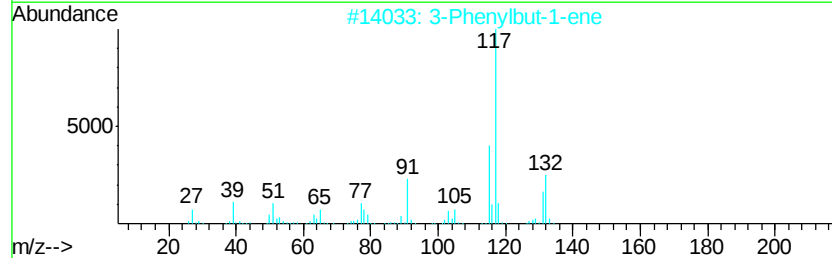
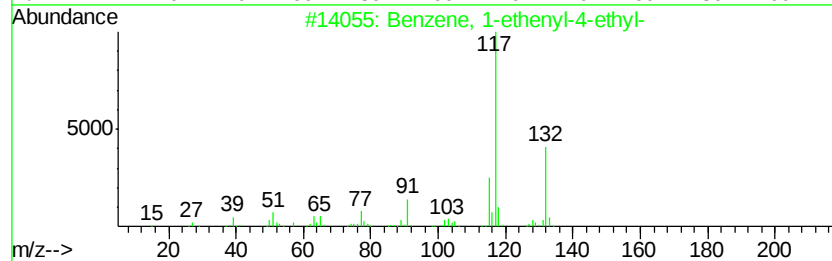
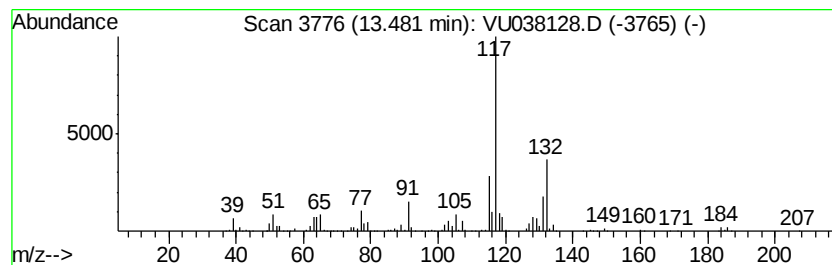
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 5

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 13.48 | 3.04 ug/L | 839678 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethenyl-4-ethyl-          | 132 | C10H12  | 003454-07-7 | 90   |
| 2    |    | 3-Phenylbut-1-ene                    | 132 | C10H12  | 000934-10-1 | 87   |
| 3    |    | Benzene, (1-methyl-1-propenyl)-, ... | 132 | C10H12  | 000768-00-3 | 83   |
| 4    |    | Benzene, 1-methyl-4-(2-propenyl)-    | 132 | C10H12  | 003333-13-9 | 83   |
| 5    |    | Benzene, 2-ethenyl-1,4-dimethyl-     | 132 | C10H12  | 002039-89-6 | 83   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleID :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

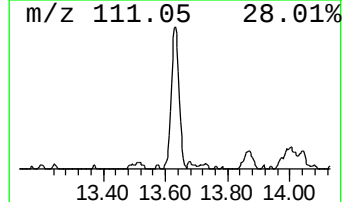
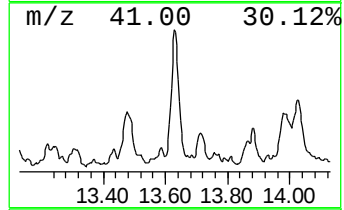
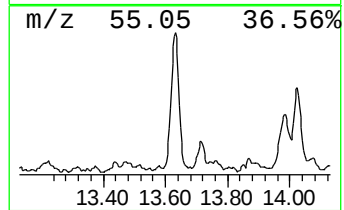
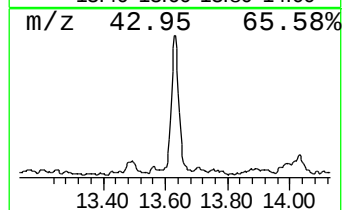
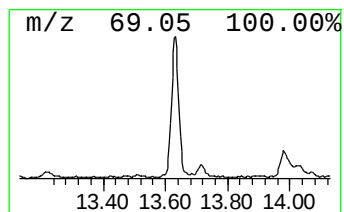
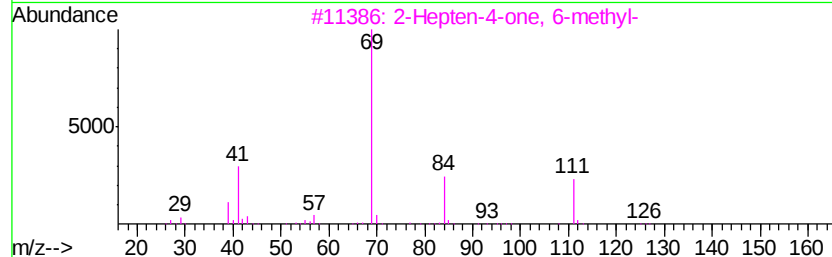
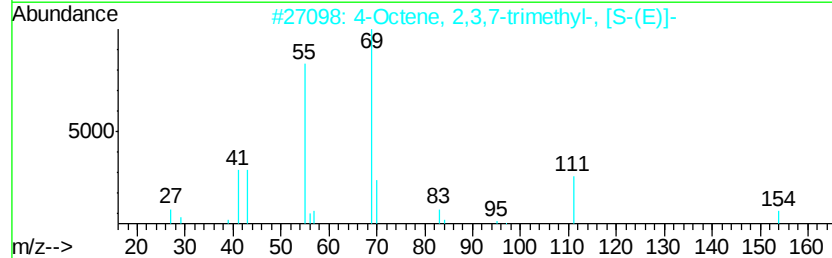
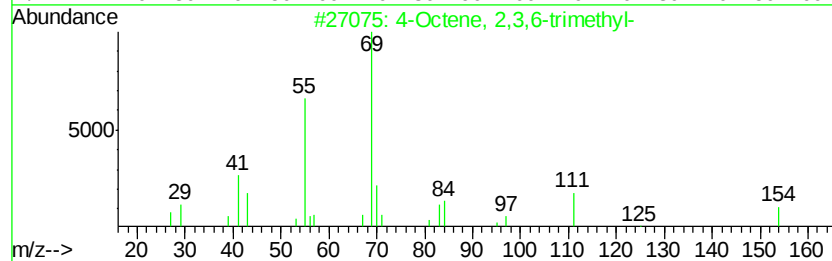
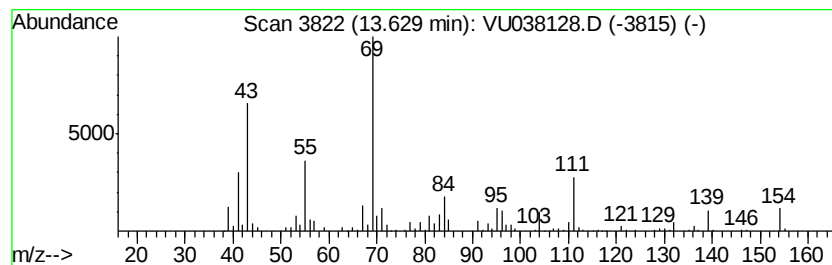
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 26

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 13.63 | 0.90 ug/L | 247266 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 4-Octene, 2,3,6-trimethyl-          | 154 | C11H22  | 063830-65-9 | 64   |
| 2    |    |   | 4-Octene, 2,3,7-trimethyl-, [S-(... | 154 | C11H22  | 052763-13-0 | 50   |
| 3    |    |   | 2-Hepten-4-one, 6-methyl-           | 126 | C8H14O  | 049852-35-9 | 50   |
| 4    |    |   | 3-Ethyl-4-octene                    | 140 | C10H20  | 019781-32-9 | 47   |
| 5    |    |   | Cyclohexane, 1-ethyl-1,3-dimethy... | 140 | C10H20  | 062238-31-7 | 47   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleID :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

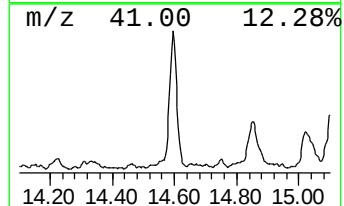
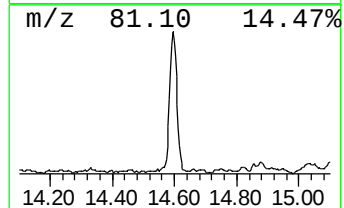
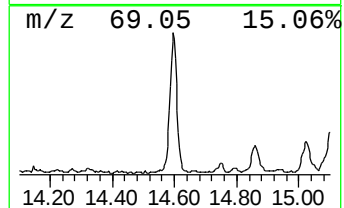
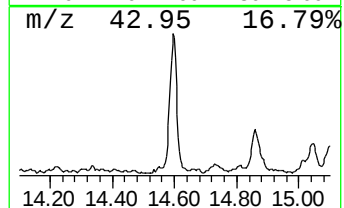
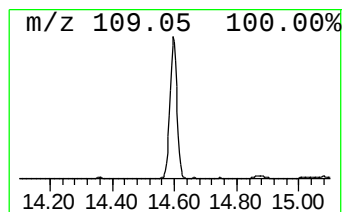
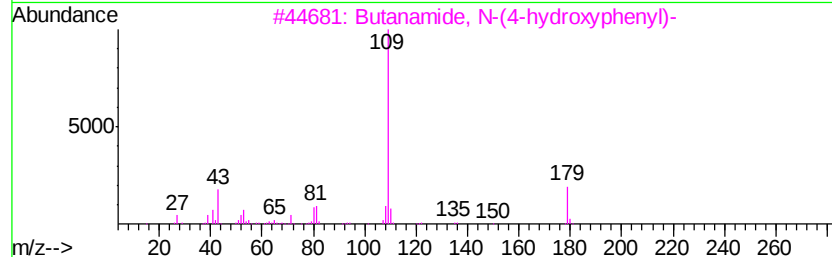
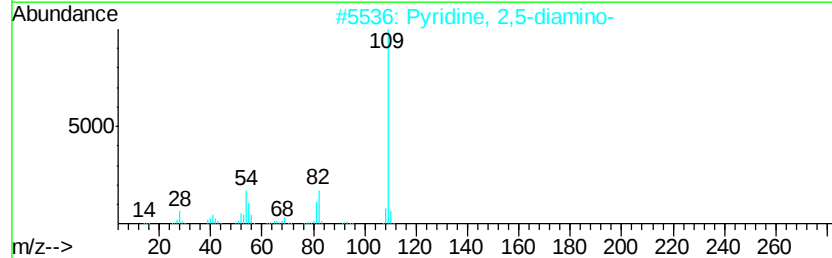
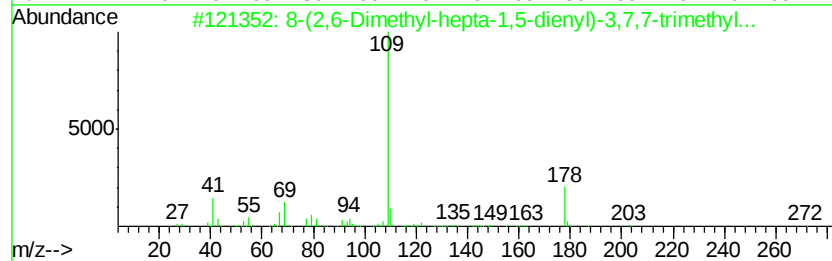
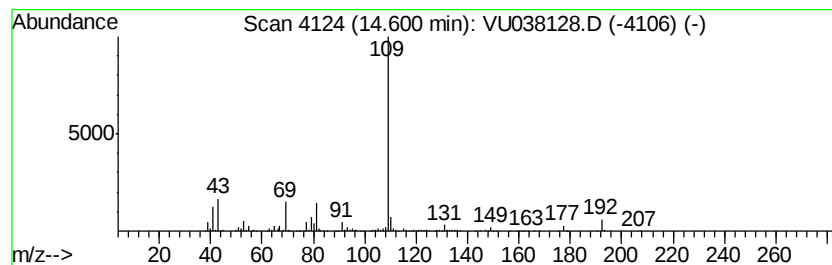
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 8-(2,6-Dimethyl-hepta-1,5-d... Concentration Rank 16

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 14.60 | 1.41 ug/L | 388899 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 8-(2,6-Dimethyl-hepta-1,5-dienyl... | 272 | C20H32    | 113725-56-7  | 64   |
| 2    |    | Pyridine, 2,5-diamino-              | 109 | C5H7N3    | 004318-76-7  | 64   |
| 3    |    | Butanamide, N-(4-hydroxyphenyl)-    | 179 | C10H13NO2 | 000101-91-7  | 59   |
| 4    |    | 1,2-Benzenediol, O-(ethoxycarbon... | 264 | C13H12O6  | 1000329-71-6 | 56   |
| 5    |    | Pyridine, 3-methoxy-                | 109 | C6H7NO    | 007295-76-3  | 50   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

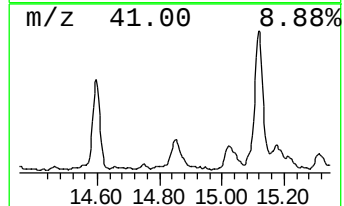
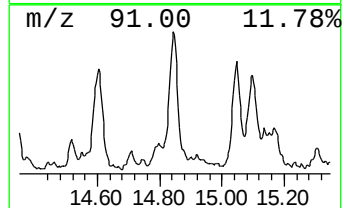
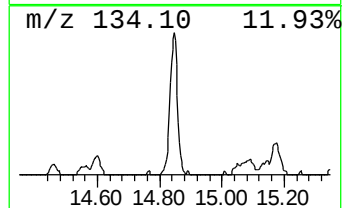
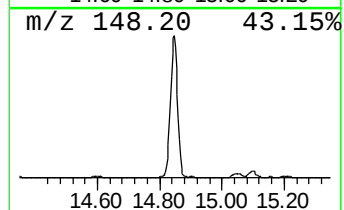
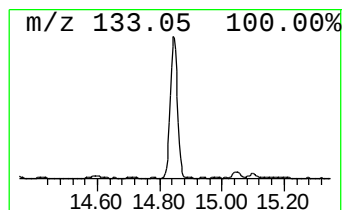
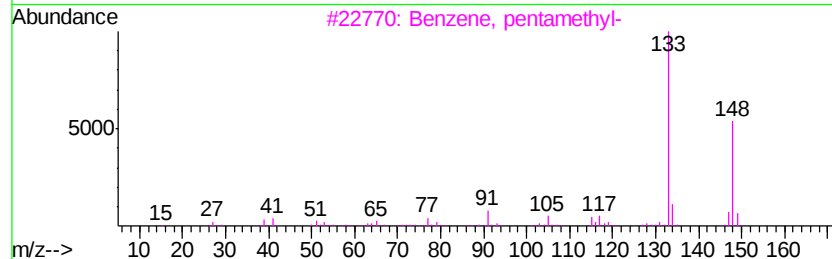
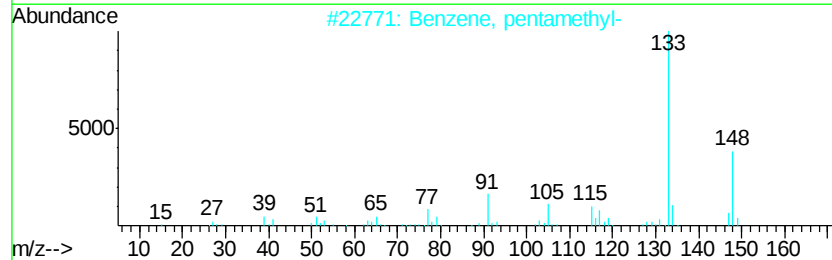
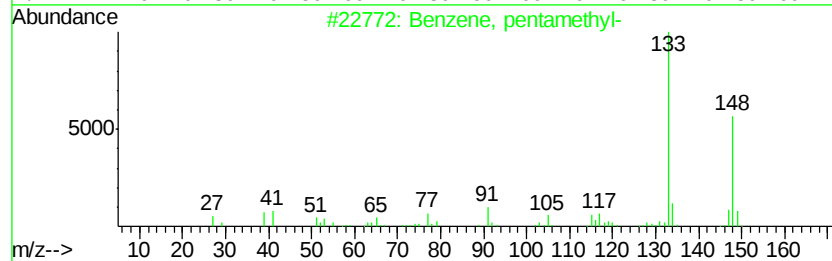
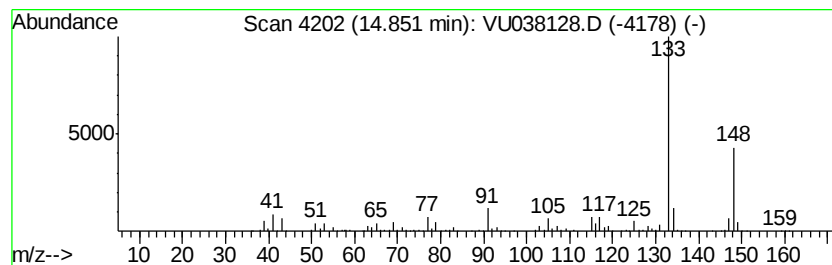
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 31 Benzene, pentamethyl- Concentration Rank 22

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 14.85 | 1.13 ug/L | 312449 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9  | 95   |
| 2    |    | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9  | 95   |
| 3    |    | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9  | 93   |
| 4    |    | 3,4-Dimethylcumene                  | 148 | C11H16  | 1000370-34-1 | 91   |
| 5    |    | Benzene, 2,4-dimethyl-1-(1-methy... | 148 | C11H16  | 004706-89-2  | 90   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

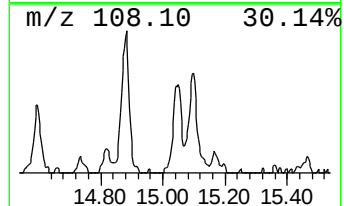
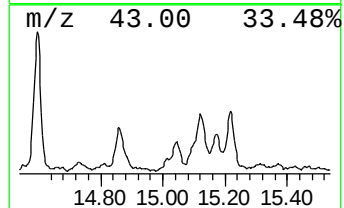
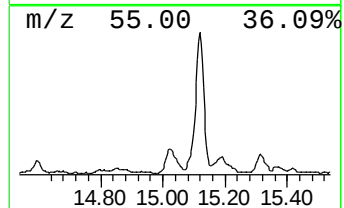
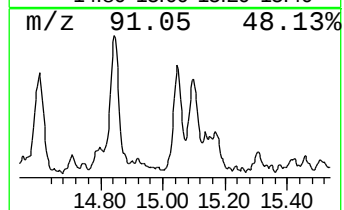
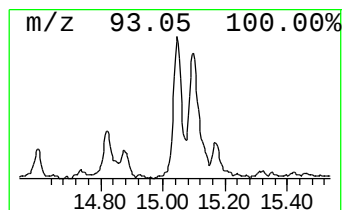
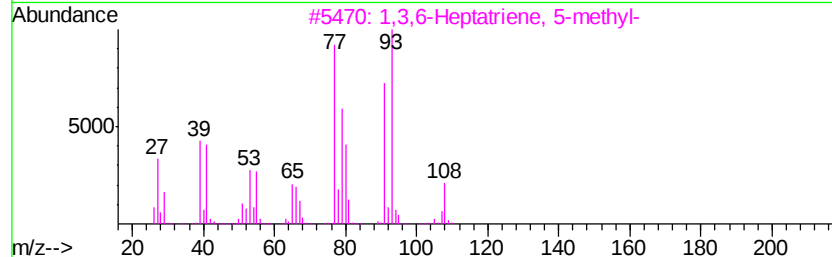
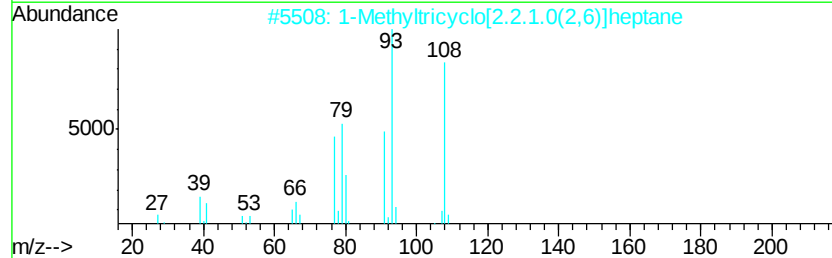
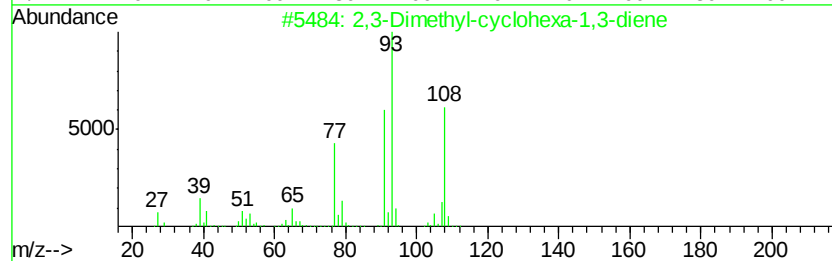
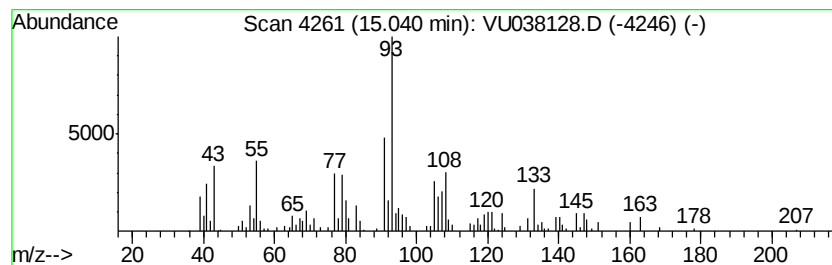
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 unknown-02 Concentration Rank 32

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 15.04 | 0.60 ug/L | 164518 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2,3-Dimethyl-cyclohexa-1,3-diene    | 108 | C8H12        | 004430-91-5 | 46      |      |      |
| 2    | 1-Methyltricyclo[2.2.1.0(2,6)]he... | 108 | C8H12        | 004601-85-8 | 43      |      |      |
| 3    | 1,3,6-Heptatriene, 5-methyl-        | 108 | C8H12        | 000925-52-0 | 43      |      |      |
| 4    | 1,3-Cyclohexadiene, 5,6-dimethyl-   | 108 | C8H12        | 005715-27-5 | 43      |      |      |
| 5    | 1,3,6-Heptatriene, 5-methyl-, (E)-  | 108 | C8H12        | 000818-48-4 | 43      |      |      |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampled :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

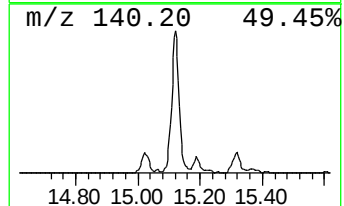
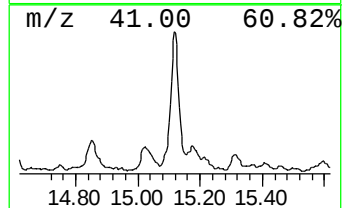
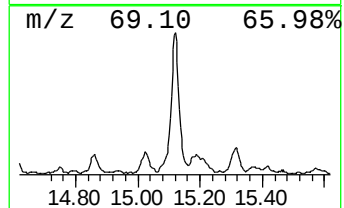
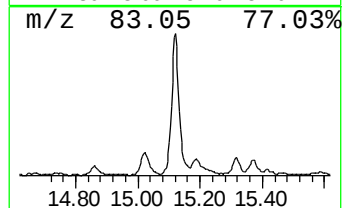
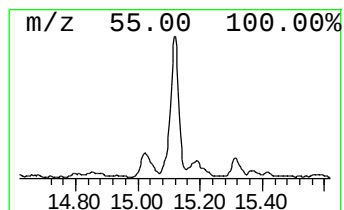
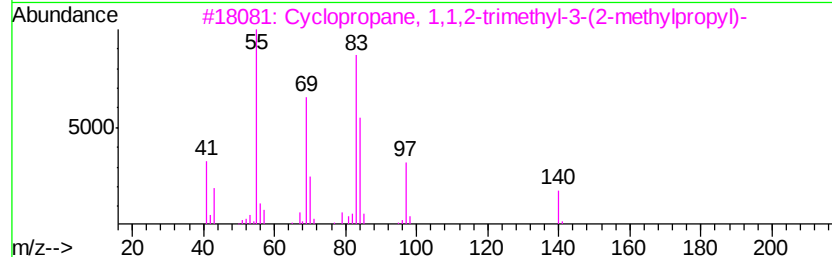
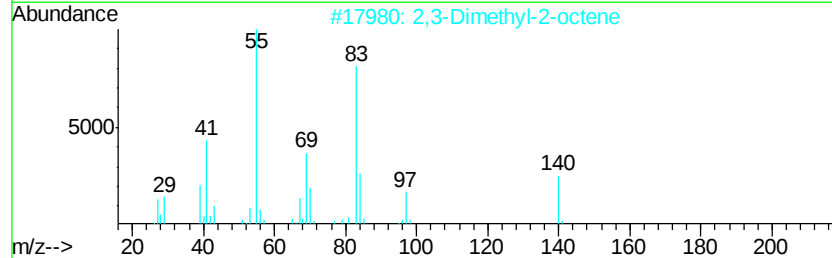
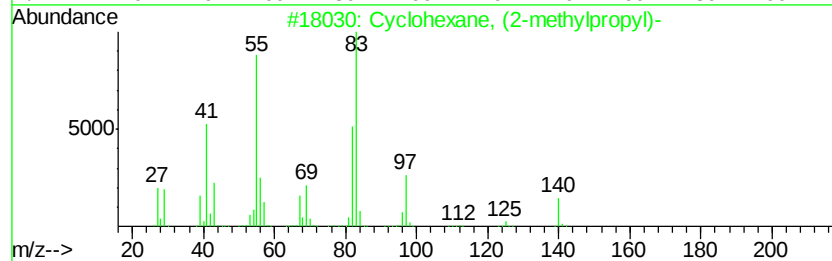
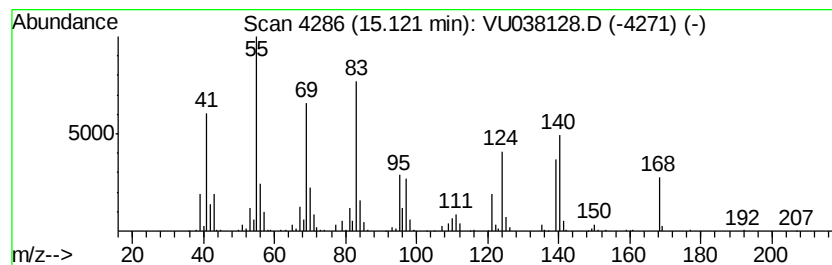
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 (DEL) Alkane: Cyclic15.12 Concentration Rank 13

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 15.12 | 1.57 ug/L | 433398 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, (2-methylpropyl)-      | 140 | C10H20  | 001678-98-4 | 55   |
| 2    |    |   | 2,3-Dimethyl-2-octene               | 140 | C10H20  | 019781-18-1 | 55   |
| 3    |    |   | Cyclopropane, 1,1,2-trimethyl-3-... | 140 | C10H20  | 041977-43-9 | 49   |
| 4    |    |   | Cyclohexane, (2-methylpropyl)-      | 140 | C10H20  | 001678-98-4 | 46   |
| 5    |    |   | 2-Methyl-3-ethyl-2-heptene          | 140 | C10H20  | 019780-61-1 | 43   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

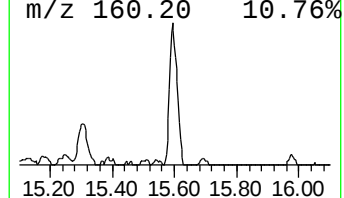
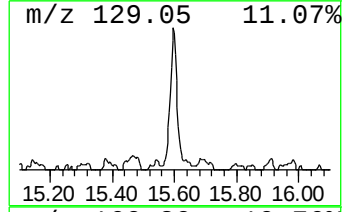
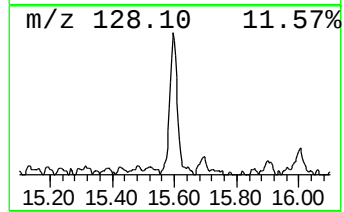
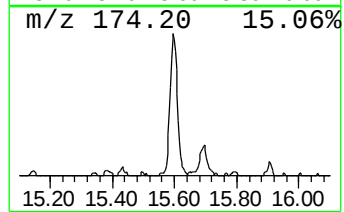
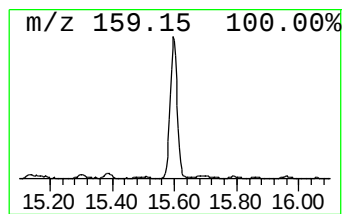
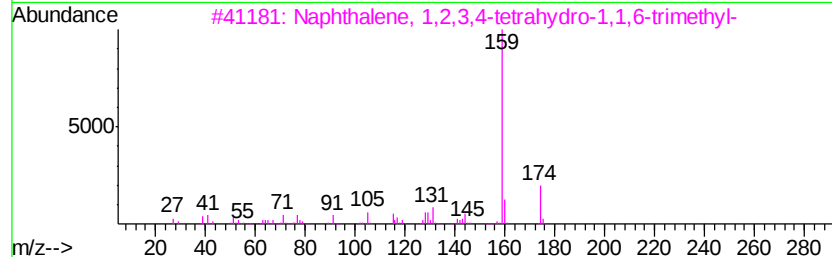
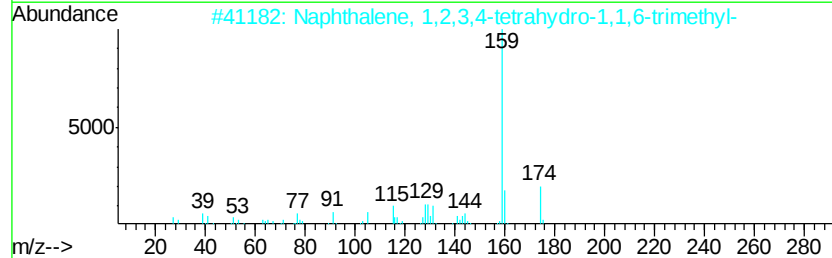
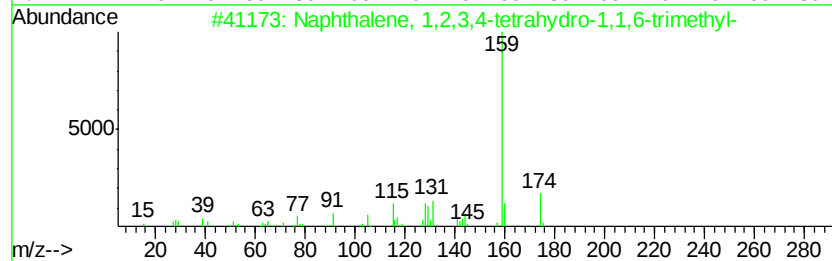
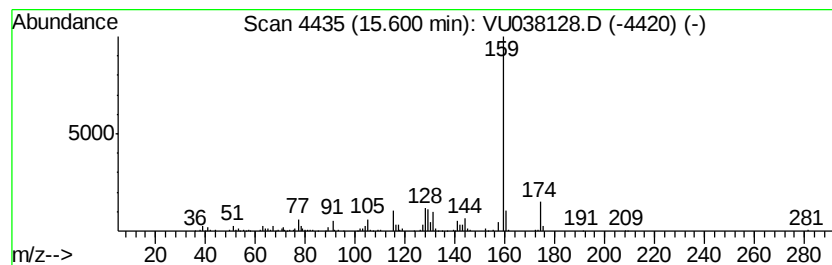
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 34 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 30

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 15.60 | 0.71 ug/L | 194562 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro-... | 174 | C13H18  | 000475-03-6 | 95   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 174 | C13H18  | 000475-03-6 | 95   |
| 3    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 174 | C13H18  | 000475-03-6 | 87   |
| 4    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 174 | C13H18  | 030316-36-0 | 87   |
| 5    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 202 | C15H22  | 000483-77-2 | 78   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

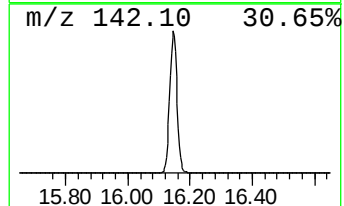
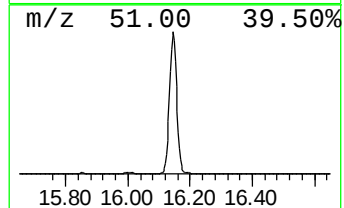
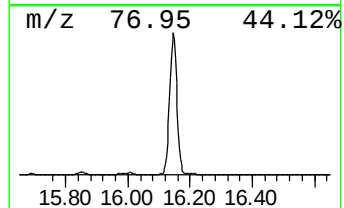
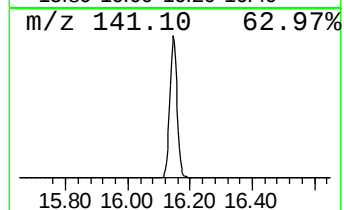
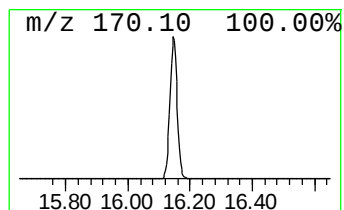
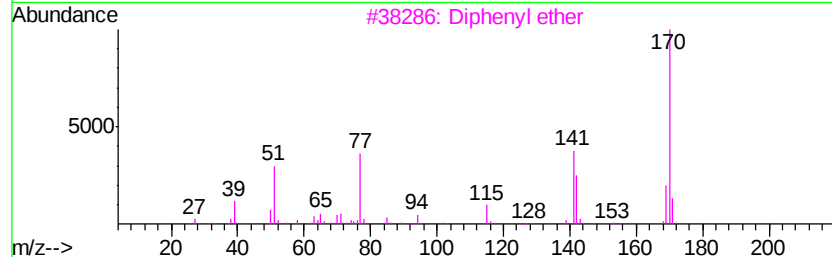
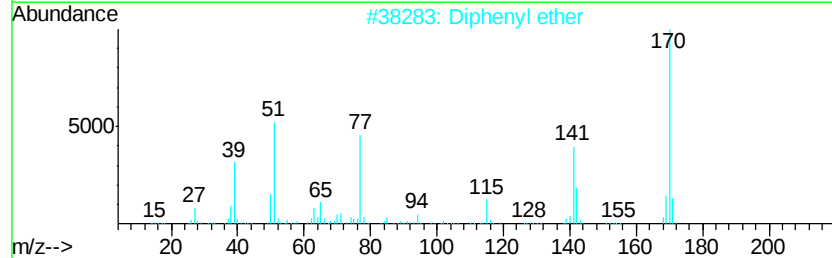
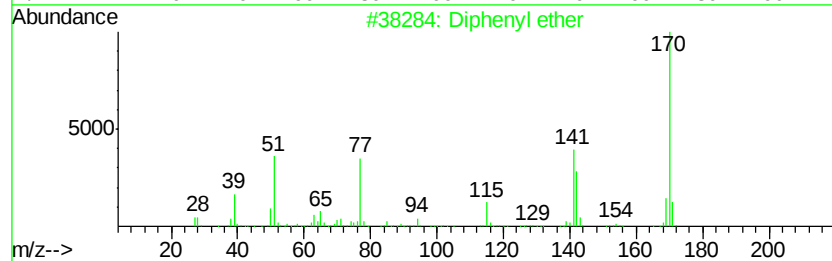
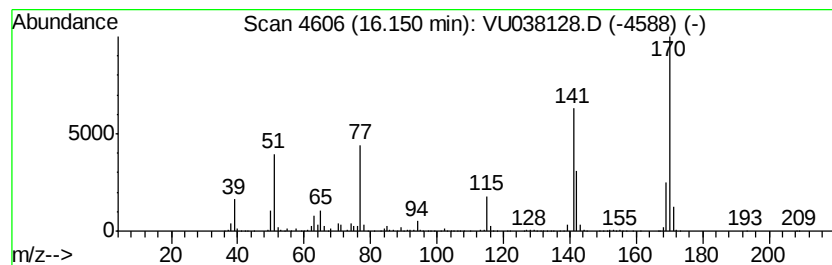
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 Diphenyl ether Concentration Rank 1

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 16.15 | 7.62 ug/L | 2102960 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Diphenyl ether                      | 170 | C12H10O | 000101-84-8 | 91   |
| 2    |    | Diphenyl ether                      | 170 | C12H10O | 000101-84-8 | 90   |
| 3    |    | Diphenyl ether                      | 170 | C12H10O | 000101-84-8 | 87   |
| 4    |    | Diphenyl ether                      | 170 | C12H10O | 000101-84-8 | 74   |
| 5    |    | Spiro[cyclopropane-1,1'(4'H)-nap... | 170 | C12H10O | 033498-24-7 | 58   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

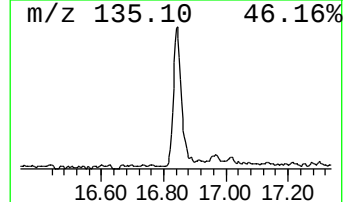
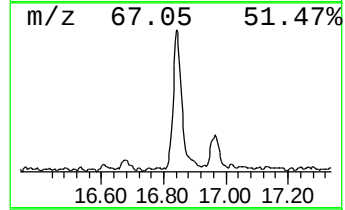
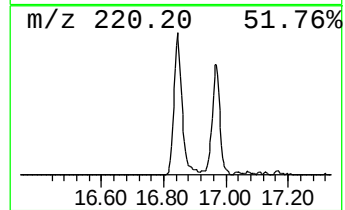
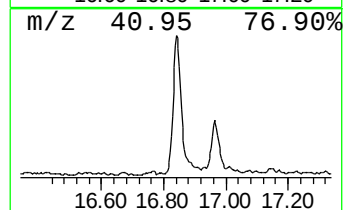
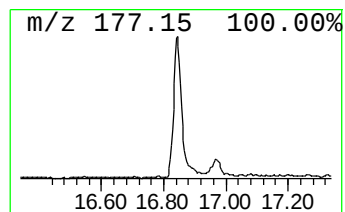
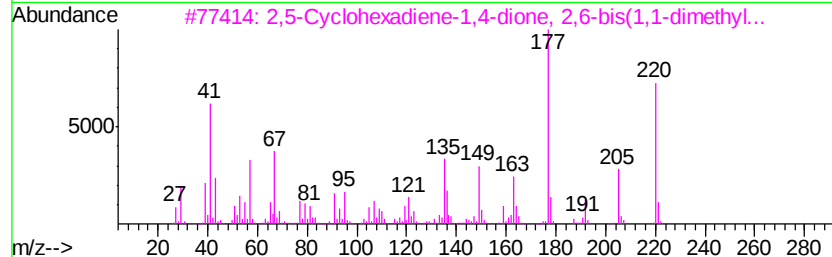
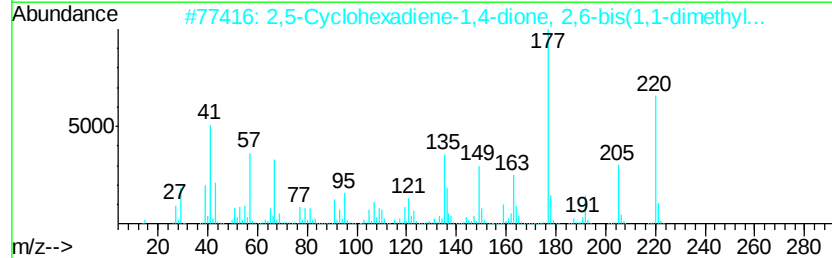
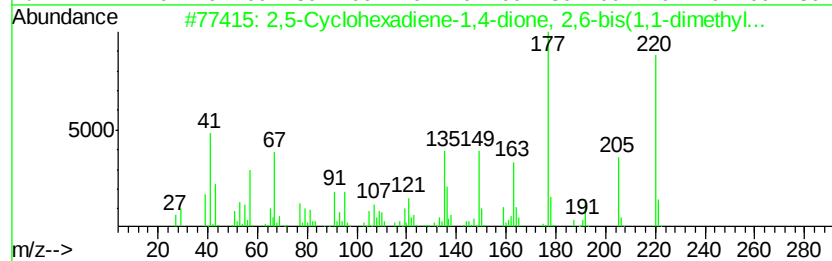
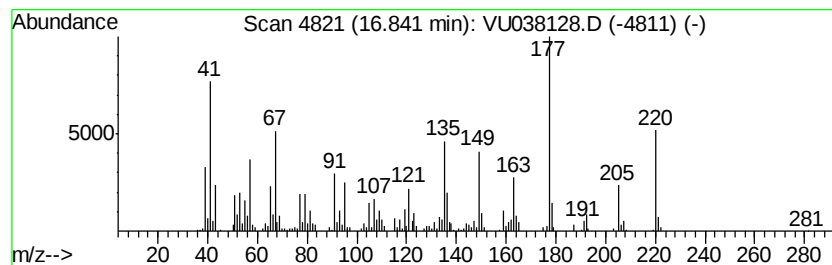
Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

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

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Peak Number 36 2,5-Cyclohexadiene-1,4-dione... Concentration Rank 14

| R.T.      | EstConc                             | Area   | Relative to ISTD       | R.T.         |      |
|-----------|-------------------------------------|--------|------------------------|--------------|------|
| 16.84     | 1.52 ug/L                           | 419398 | 1,4-Dichlorobenzene-d4 | 11.82        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm                | CAS#         | Qual |
| 1         | 2,5-Cyclohexadiene-1,4-dione, 2,... | 220    | C14H20O2               | 000719-22-2  | 98   |
| 2         | 2,5-Cyclohexadiene-1,4-dione, 2,... | 220    | C14H20O2               | 000719-22-2  | 98   |
| 3         | 2,5-Cyclohexadiene-1,4-dione, 2,... | 220    | C14H20O2               | 000719-22-2  | 98   |
| 4         | Thiocoumarin, 4,4,6,8-tetramethyl-  | 220    | C13H16OS               | 1000128-90-3 | 38   |
| 5         | Propanamine, 2-(1-adamantyl)-2-m... | 207    | C14H25N                | 079594-24-4  | 30   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
Data File : VU038128.D  
Acq On : 12 May 2020 16:45  
Operator : JC/MD  
Sample : L2570-07  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 1 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
BFSP5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
Quant Title : TRACE VOA SOM01.0

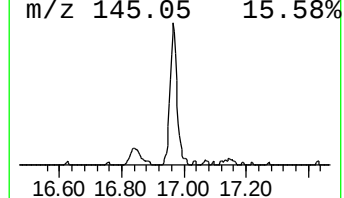
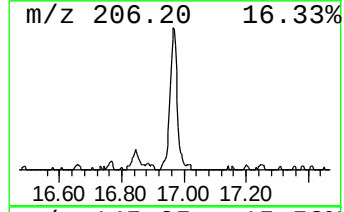
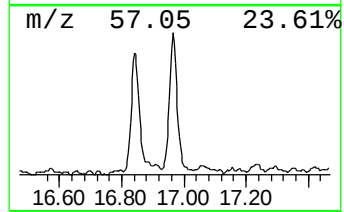
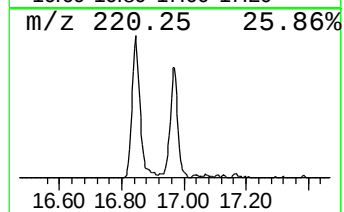
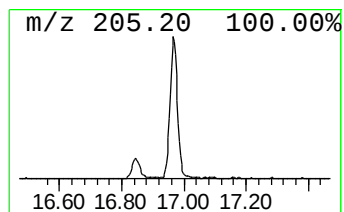
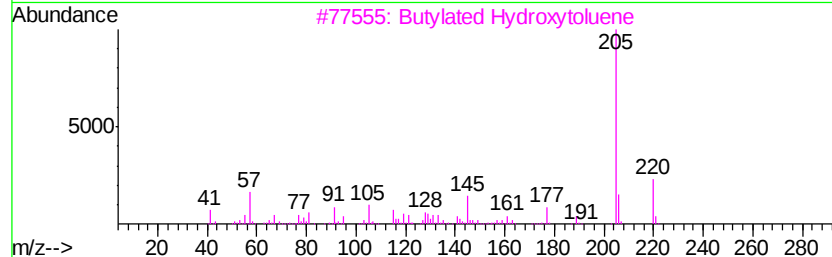
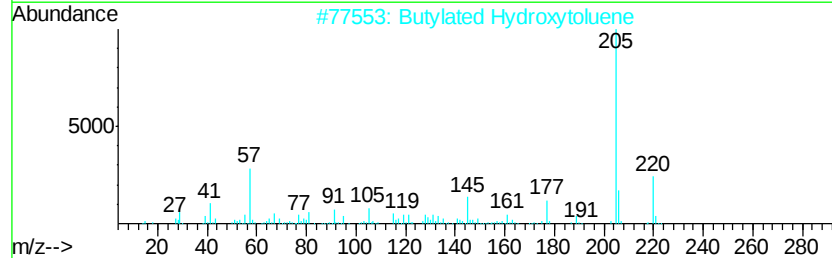
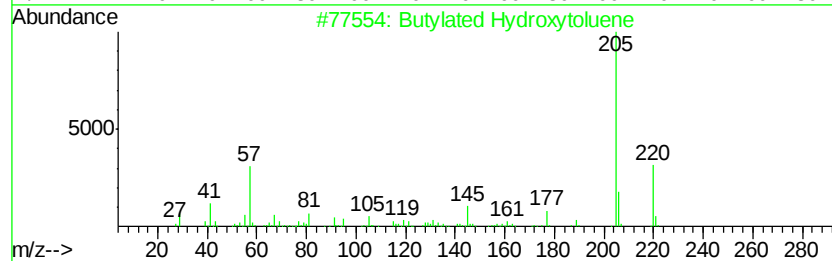
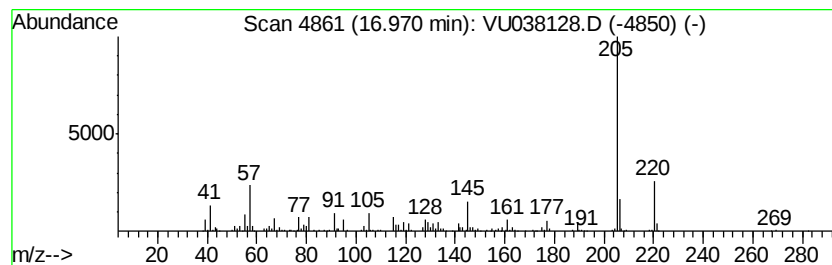
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 37 Butylated Hydroxytoluene Concentration Rank 28

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 16.97 | 0.81 ug/L | 222662 | 1,4-Dichlorobenzene-d4 | 11.82 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Butylated Hydroxytoluene            | 220 | C15H24O | 000128-37-0 | 97   |
| 2    |    | Butylated Hydroxytoluene            | 220 | C15H24O | 000128-37-0 | 96   |
| 3    |    | Butylated Hydroxytoluene            | 220 | C15H24O | 000128-37-0 | 96   |
| 4    |    | Butylated Hydroxytoluene            | 220 | C15H24O | 000128-37-0 | 94   |
| 5    |    | Phenol, 4,6-di(1,1-dimethylethyl... | 220 | C15H24O | 000616-55-7 | 81   |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_U\DATA\VU051220\  
 Data File : VU038128.D  
 Acq On : 12 May 2020 16:45  
 Operator : JC/MD  
 Sample : L2570-07  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 1 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 BFSP5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_U\METHOD\SOMUTR050720WMA.M  
 Quant Title : TRACE VOA SOM01.0

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| unknown-01           | 2.40  | 1.5     | ug/L  | 263919   | 1                     | 6.28  | 877343  | 5.0  |
| (DEL) Alkane: Str... | 4.33  | 0.5     | ug/L  | 93446    | 1                     | 6.28  | 877343  | 5.0  |
| (DEL) Alkane: Str... | 4.47  | 0.5     | ug/L  | 94152    | 1                     | 6.28  | 877343  | 5.0  |
| (DEL) Alkane: Cyc... | 4.57  | 1.4     | ug/L  | 236638   | 1                     | 6.28  | 877343  | 5.0  |
| (DEL) Alkane: Str... | 5.17  | 1.1     | ug/L  | 187208   | 1                     | 6.28  | 877343  | 5.0  |
| (DEL) Alkane: Str... | 5.49  | 2.5     | ug/L  | 447982   | 1                     | 6.28  | 877343  | 5.0  |
| (DEL) Alkane: Str... | 5.63  | 0.8     | ug/L  | 138915   | 1                     | 6.28  | 877343  | 5.0  |
| (DEL) Alkane: Cyc... | 5.67  | 0.6     | ug/L  | 99329    | 1                     | 6.28  | 877343  | 5.0  |
| 1-Butanol, 2-ethyl-  | 5.88  | 1.8     | ug/L  | 321673   | 1                     | 6.28  | 877343  | 5.0  |
| (DEL) Alkane: Cyc... | 5.95  | 0.9     | ug/L  | 163870   | 1                     | 6.28  | 877343  | 5.0  |
| (DEL) Alkane: Str... | 6.02  | 0.6     | ug/L  | 103789   | 1                     | 6.28  | 877343  | 5.0  |
| 1,5-Cyclooctadiene   | 10.58 | 2.1     | ug/L  | 1427720  | 2                     | 9.44  | 3471490 | 5.0  |
| Benzene, propyl-     | 10.92 | 4.6     | ug/L  | 1264390  | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, 1-ethyl-... | 11.01 | 0.8     | ug/L  | 233201   | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, 1,2,3-tr... | 11.10 | 0.7     | ug/L  | 182810   | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, 1-ethyl-... | 11.30 | 1.2     | ug/L  | 338013   | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, 1,2,4-tr... | 11.48 | 3.7     | ug/L  | 1029480  | 3                     | 11.82 | 1379510 | 5.0  |
| Mesitylene           | 11.91 | 1.3     | ug/L  | 352900   | 3                     | 11.82 | 1379510 | 5.0  |
| Indane               | 12.10 | 5.1     | ug/L  | 1395310  | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, 1-ethyl-... | 12.48 | 0.5     | ug/L  | 146874   | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, (3-iodo-... | 12.55 | 2.4     | ug/L  | 673023   | 3                     | 11.82 | 1379510 | 5.0  |
| Indan, 1-methyl-     | 12.70 | 1.6     | ug/L  | 442179   | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, 4-chloro... | 12.84 | 1.9     | ug/L  | 536406   | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, 1,2,4,5-... | 12.99 | 1.3     | ug/L  | 347564   | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, 1,2,3,5-... | 13.05 | 1.1     | ug/L  | 309281   | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, 2-etheny... | 13.31 | 1.3     | ug/L  | 365162   | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, 1-etheny... | 13.48 | 3.0     | ug/L  | 839678   | 3                     | 11.82 | 1379510 | 5.0  |
| (DEL) Alkane: Str... | 13.63 | 0.9     | ug/L  | 247266   | 3                     | 11.82 | 1379510 | 5.0  |
| 8-(2,6-Dimethyl-h... | 14.60 | 1.4     | ug/L  | 388899   | 3                     | 11.82 | 1379510 | 5.0  |
| Benzene, pentamet... | 14.85 | 1.1     | ug/L  | 312449   | 3                     | 11.82 | 1379510 | 5.0  |
| unknown-02           | 15.04 | 0.6     | ug/L  | 164518   | 3                     | 11.82 | 1379510 | 5.0  |
| (DEL) Alkane: Cyc... | 15.12 | 1.6     | ug/L  | 433398   | 3                     | 11.82 | 1379510 | 5.0  |
| Naphthalene, 1,2,... | 15.60 | 0.7     | ug/L  | 194562   | 3                     | 11.82 | 1379510 | 5.0  |
| Diphenyl ether       | 16.15 | 7.6     | ug/L  | 2102960  | 3                     | 11.82 | 1379510 | 5.0  |
| 2,5-Cyclohexadien... | 16.84 | 1.5     | ug/L  | 419398   | 3                     | 11.82 | 1379510 | 5.0  |
| Butylated Hydroxy... | 16.97 | 0.8     | ug/L  | 222662   | 3                     | 11.82 | 1379510 | 5.0  |