

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
 Data File : VV022511.D  
 Acq On : 02 Oct 2021 15:19  
 Operator : SY/MD  
 Sample : M3996-14  
 Misc : 5.0mL/MSVOA\_V/WATER  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 MSVOA\_V  
 ClientSampleId :  
 F4A29

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
 Title : VOC Analysis

Signal : TIC: VV022511.D\data.ms

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 1.214       | 49            | 54          | 60           | rBV      | 13752          | 12196         | 1.65%           | 0.149%        |
| 2         | 1.304       | 76            | 82          | 105          | rVB      | 138904         | 149325        | 20.25%          | 1.826%        |
| 3         | 1.568       | 156           | 164         | 174          | rBV      | 93665          | 105961        | 14.37%          | 1.296%        |
| 4         | 1.632       | 175           | 184         | 201          | rVB      | 151390         | 187163        | 25.39%          | 2.289%        |
| 5         | 1.806       | 230           | 238         | 249          | rBV3     | 17379          | 24619         | 3.34%           | 0.301%        |
| 6         | 1.947       | 280           | 282         | 291          | rVB7     | 1324           | 1268          | 0.17%           | 0.016%        |
| 7         | 2.024       | 296           | 306         | 321          | rVB      | 16644          | 23783         | 3.23%           | 0.291%        |
| 8         | 2.108       | 321           | 332         | 348          | rBV      | 202604         | 327812        | 44.46%          | 4.009%        |
| 9         | 2.497       | 441           | 453         | 458          | rBV      | 48245          | 89273         | 12.11%          | 1.092%        |
| 10        | 2.764       | 522           | 536         | 560          | rVB2     | 156468         | 331871        | 45.01%          | 4.059%        |
| 11        | 2.937       | 582           | 590         | 599          | rVB7     | 1357           | 2570          | 0.35%           | 0.031%        |
| 12        | 3.047       | 612           | 624         | 637          | rBV6     | 5694           | 12336         | 1.67%           | 0.151%        |
| 13        | 3.275       | 685           | 695         | 699          | rBV6     | 3641           | 6408          | 0.87%           | 0.078%        |
| 14        | 3.368       | 712           | 724         | 733          | rBV5     | 5320           | 11874         | 1.61%           | 0.145%        |
| 15        | 3.535       | 762           | 776         | 778          | rBV5     | 3103           | 5197          | 0.70%           | 0.064%        |
| 16        | 3.667       | 806           | 817         | 832          | rVB3     | 14787          | 35138         | 4.77%           | 0.430%        |
| 17        | 3.764       | 833           | 847         | 863          | rVV2     | 39274          | 94752         | 12.85%          | 1.159%        |
| 18        | 3.889       | 877           | 886         | 922          | rVB2     | 54183          | 165936        | 22.51%          | 2.029%        |
| 19        | 4.352       | 1013          | 1030        | 1059         | rBV2     | 100644         | 304237        | 41.27%          | 3.721%        |
| 20        | 4.677       | 1115          | 1131        | 1145         | rBV3     | 45863          | 116296        | 15.77%          | 1.422%        |
| 21        | 4.767       | 1146          | 1159        | 1180         | rVB2     | 49861          | 131013        | 17.77%          | 1.602%        |
| 22        | 4.915       | 1188          | 1205        | 1230         | rBV7     | 15544          | 53255         | 7.22%           | 0.651%        |
| 23        | 5.047       | 1230          | 1246        | 1258         | rBV2     | 289832         | 737271        | 100.00%         | 9.016%        |
| 24        | 5.494       | 1376          | 1385        | 1386         | rBV4     | 3690           | 4788          | 0.65%           | 0.059%        |
| 25        | 5.619       | 1411          | 1424        | 1461         | rBV      | 247124         | 538537        | 73.04%          | 6.586%        |
| 26        | 5.773       | 1464          | 1472        | 1485         | rVV8     | 3690           | 7328          | 0.99%           | 0.090%        |
| 27        | 5.847       | 1485          | 1495        | 1505         | rVV9     | 2950           | 5704          | 0.77%           | 0.070%        |
| 28        | 5.924       | 1506          | 1519        | 1537         | rVB9     | 17880          | 47036         | 6.38%           | 0.575%        |
| 29        | 6.072       | 1547          | 1565        | 1578         | rBV      | 159752         | 351348        | 47.66%          | 4.297%        |
| 30        | 6.204       | 1597          | 1606        | 1613         | rBV6     | 3905           | 6210          | 0.84%           | 0.076%        |
| 31        | 6.259       | 1617          | 1623        | 1632         | rVB6     | 6964           | 11433         | 1.55%           | 0.140%        |
| 32        | 6.368       | 1649          | 1657        | 1673         | rVB5     | 5650           | 12143         | 1.65%           | 0.149%        |
| 33        | 6.465       | 1675          | 1687        | 1699         | rVB8     | 6499           | 14141         | 1.92%           | 0.173%        |
| 34        | 6.545       | 1703          | 1712        | 1713         | rBV4     | 2633           | 3003          | 0.41%           | 0.037%        |
| 35        | 6.654       | 1733          | 1746        | 1761         | rVB      | 44768          | 92063         | 12.49%          | 1.126%        |
| 36        | 6.780       | 1770          | 1785        | 1797         | rBV2     | 58941          | 128927        | 17.49%          | 1.577%        |

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 MSVOA\_V  
 ClientSampleId :  
 F4A29

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
 Title : VOC Analysis

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 37 | 6.989  | 1835 | 1850 | 1869 | rBV  | 91175  | 176052 | 23.88% | 2.153% |
| 38 | 7.268  | 1926 | 1937 | 1941 | rBV7 | 10437  | 18437  | 2.50%  | 0.225% |
| 39 | 7.317  | 1942 | 1952 | 1970 | rVB  | 341863 | 597696 | 81.07% | 7.310% |
| 40 | 7.622  | 2039 | 2047 | 2068 | rVB  | 56134  | 108761 | 14.75% | 1.330% |
| 41 | 7.792  | 2082 | 2100 | 2109 | rBV5 | 8402   | 17850  | 2.42%  | 0.218% |
| 42 | 7.844  | 2109 | 2116 | 2126 | rBV3 | 11052  | 18552  | 2.52%  | 0.227% |
| 43 | 7.937  | 2137 | 2145 | 2153 | rBV3 | 6577   | 9960   | 1.35%  | 0.122% |
| 44 | 7.985  | 2153 | 2160 | 2169 | rVB7 | 1895   | 3254   | 0.44%  | 0.040% |
| 45 | 8.091  | 2183 | 2193 | 2206 | rBV  | 188467 | 331698 | 44.99% | 4.056% |
| 46 | 8.169  | 2210 | 2217 | 2228 | rVB  | 56256  | 96201  | 13.05% | 1.176% |
| 47 | 8.516  | 2313 | 2325 | 2332 | rBV7 | 4234   | 7647   | 1.04%  | 0.094% |
| 48 | 8.628  | 2353 | 2360 | 2372 | rVB5 | 8748   | 14970  | 2.03%  | 0.183% |
| 49 | 8.725  | 2383 | 2390 | 2402 | rBV3 | 5530   | 9383   | 1.27%  | 0.115% |
| 50 | 8.815  | 2408 | 2418 | 2419 | rBV6 | 3432   | 5434   | 0.74%  | 0.066% |
| 51 | 8.854  | 2420 | 2430 | 2457 | rVB  | 364834 | 585453 | 79.41% | 7.160% |
| 52 | 9.146  | 2516 | 2521 | 2529 | rVB4 | 6319   | 9255   | 1.26%  | 0.113% |
| 53 | 9.223  | 2538 | 2545 | 2555 | rBV2 | 4474   | 7177   | 0.97%  | 0.088% |
| 54 | 9.439  | 2602 | 2612 | 2616 | rBV3 | 1960   | 2815   | 0.38%  | 0.034% |
| 55 | 9.471  | 2620 | 2622 | 2633 | rVB8 | 2178   | 2652   | 0.36%  | 0.032% |
| 56 | 9.554  | 2640 | 2648 | 2656 | rBV6 | 2582   | 3884   | 0.53%  | 0.047% |
| 57 | 9.735  | 2693 | 2704 | 2717 | rVB4 | 8305   | 16469  | 2.23%  | 0.201% |
| 58 | 9.802  | 2718 | 2725 | 2732 | rBV7 | 2776   | 3966   | 0.54%  | 0.049% |
| 59 | 9.847  | 2734 | 2739 | 2741 | rBV3 | 1408   | 1347   | 0.18%  | 0.016% |
| 60 | 9.937  | 2756 | 2767 | 2775 | rVV  | 8318   | 13813  | 1.87%  | 0.169% |
| 61 | 9.998  | 2775 | 2786 | 2797 | rBV2 | 15626  | 26577  | 3.60%  | 0.325% |
| 62 | 10.091 | 2809 | 2815 | 2820 | rBV6 | 2506   | 3361   | 0.46%  | 0.041% |
| 63 | 10.136 | 2822 | 2829 | 2838 | rVB3 | 7723   | 12824  | 1.74%  | 0.157% |
| 64 | 10.217 | 2844 | 2854 | 2871 | rVB  | 226307 | 350438 | 47.53% | 4.286% |
| 65 | 10.407 | 2909 | 2913 | 2923 | rVB2 | 5563   | 8064   | 1.09%  | 0.099% |
| 66 | 10.538 | 2947 | 2954 | 2957 | rBV5 | 1644   | 2144   | 0.29%  | 0.026% |
| 67 | 10.741 | 3010 | 3017 | 3026 | rVB2 | 6261   | 9724   | 1.32%  | 0.119% |
| 68 | 10.853 | 3045 | 3052 | 3059 | rVB8 | 1328   | 2076   | 0.28%  | 0.025% |
| 69 | 10.921 | 3062 | 3073 | 3081 | rBV6 | 3775   | 5760   | 0.78%  | 0.070% |
| 70 | 11.056 | 3100 | 3115 | 3120 | rBV6 | 2782   | 4795   | 0.65%  | 0.059% |
| 71 | 11.185 | 3150 | 3155 | 3163 | rVB6 | 2722   | 3897   | 0.53%  | 0.048% |
| 72 | 11.249 | 3164 | 3175 | 3191 | rVV  | 386956 | 590583 | 80.10% | 7.223% |
| 73 | 11.336 | 3195 | 3202 | 3213 | rVV  | 27457  | 44660  | 6.06%  | 0.546% |
| 74 | 11.394 | 3213 | 3220 | 3224 | rVB7 | 2333   | 3146   | 0.43%  | 0.038% |
| 75 | 11.548 | 3256 | 3268 | 3278 | rBV8 | 9947   | 18440  | 2.50%  | 0.226% |
| 76 | 11.625 | 3281 | 3292 | 3310 | rVB  | 350294 | 544740 | 73.89% | 6.662% |

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Instrument :  
 MSVOA\_V  
 ClientSampleId :  
 F4A29

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
 Title : VOC Analysis

|     |        |      |      |      |      |       |       |        |        |
|-----|--------|------|------|------|------|-------|-------|--------|--------|
| 77  | 11.747 | 3324 | 3330 | 3338 | rVV6 | 3719  | 4860  | 0.66%  | 0.059% |
| 78  | 12.021 | 3407 | 3415 | 3424 | rVB2 | 4581  | 6370  | 0.86%  | 0.078% |
| 79  | 12.117 | 3435 | 3445 | 3462 | rVB2 | 30159 | 49846 | 6.76%  | 0.610% |
| 80  | 12.287 | 3488 | 3498 | 3503 | rBV6 | 1502  | 2702  | 0.37%  | 0.033% |
| 81  | 12.320 | 3504 | 3508 | 3516 | rVV7 | 2271  | 3010  | 0.41%  | 0.037% |
| 82  | 12.416 | 3526 | 3538 | 3549 | rBV  | 13227 | 20660 | 2.80%  | 0.253% |
| 83  | 12.468 | 3552 | 3554 | 3559 | rVV4 | 1544  | 1242  | 0.17%  | 0.015% |
| 84  | 12.725 | 3630 | 3634 | 3646 | rVB5 | 1911  | 2728  | 0.37%  | 0.033% |
| 85  | 12.886 | 3674 | 3684 | 3695 | rBV2 | 30460 | 48132 | 6.53%  | 0.589% |
| 86  | 12.950 | 3695 | 3704 | 3713 | rVV4 | 9151  | 14234 | 1.93%  | 0.174% |
| 87  | 13.053 | 3728 | 3736 | 3749 | rVB7 | 3137  | 6365  | 0.86%  | 0.078% |
| 88  | 13.194 | 3772 | 3780 | 3782 | rBV2 | 1940  | 2102  | 0.29%  | 0.026% |
| 89  | 13.307 | 3808 | 3815 | 3817 | rVV4 | 1592  | 1600  | 0.22%  | 0.020% |
| 90  | 13.336 | 3820 | 3824 | 3830 | rVV6 | 2987  | 4375  | 0.59%  | 0.054% |
| 91  | 13.371 | 3831 | 3835 | 3841 | rVB4 | 2529  | 3067  | 0.42%  | 0.038% |
| 92  | 13.503 | 3865 | 3876 | 3890 | rBV  | 50522 | 80171 | 10.87% | 0.980% |
| 93  | 13.628 | 3906 | 3915 | 3925 | rBV  | 10110 | 15016 | 2.04%  | 0.184% |
| 94  | 13.767 | 3952 | 3958 | 3967 | rVB3 | 1187  | 1616  | 0.22%  | 0.020% |
| 95  | 13.911 | 3998 | 4003 | 4010 | rVB4 | 1424  | 1581  | 0.21%  | 0.019% |
| 96  | 14.152 | 4070 | 4078 | 4085 | rBV4 | 2481  | 3772  | 0.51%  | 0.046% |
| 97  | 14.236 | 4095 | 4104 | 4105 | rBV5 | 1552  | 1671  | 0.23%  | 0.020% |
| 98  | 14.355 | 4135 | 4141 | 4148 | rBV5 | 2015  | 2331  | 0.32%  | 0.029% |
| 99  | 14.577 | 4201 | 4210 | 4226 | rVB3 | 5881  | 9514  | 1.29%  | 0.116% |
| 100 | 14.818 | 4274 | 4285 | 4303 | rBV2 | 26381 | 43824 | 5.94%  | 0.536% |

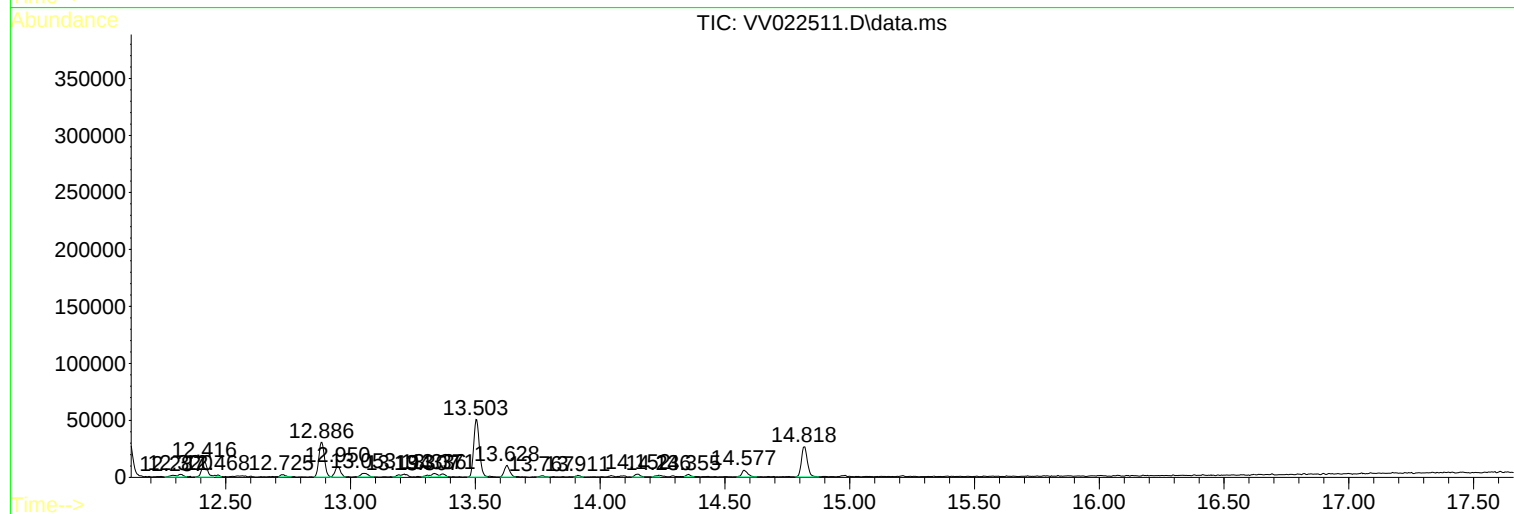
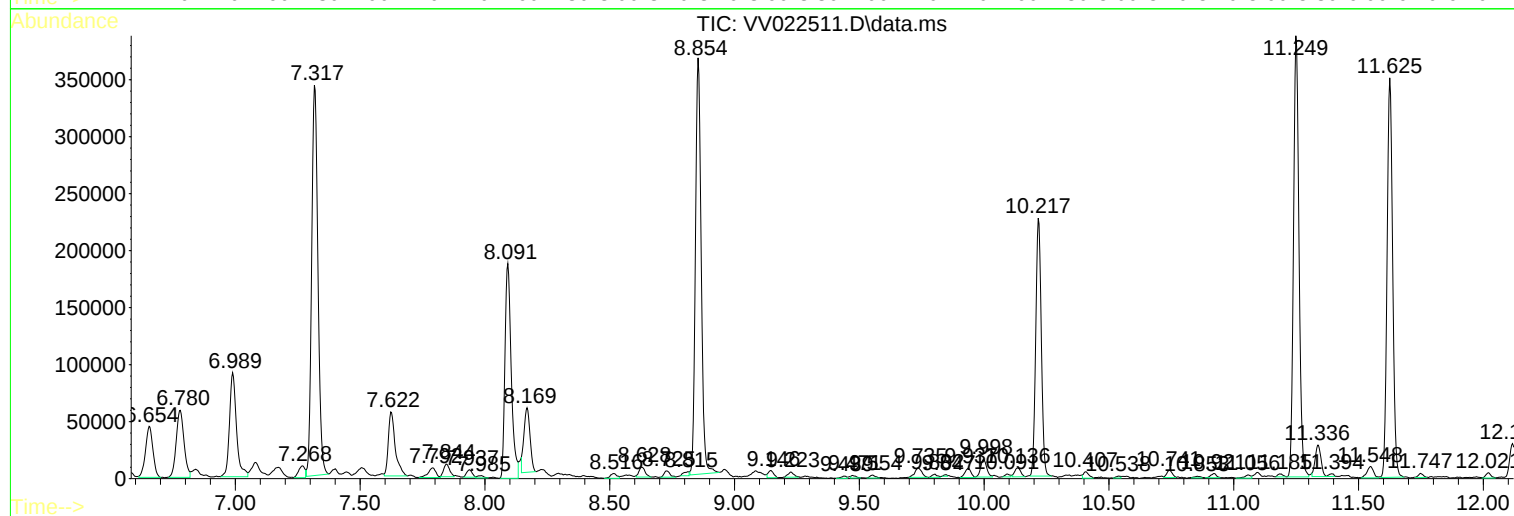
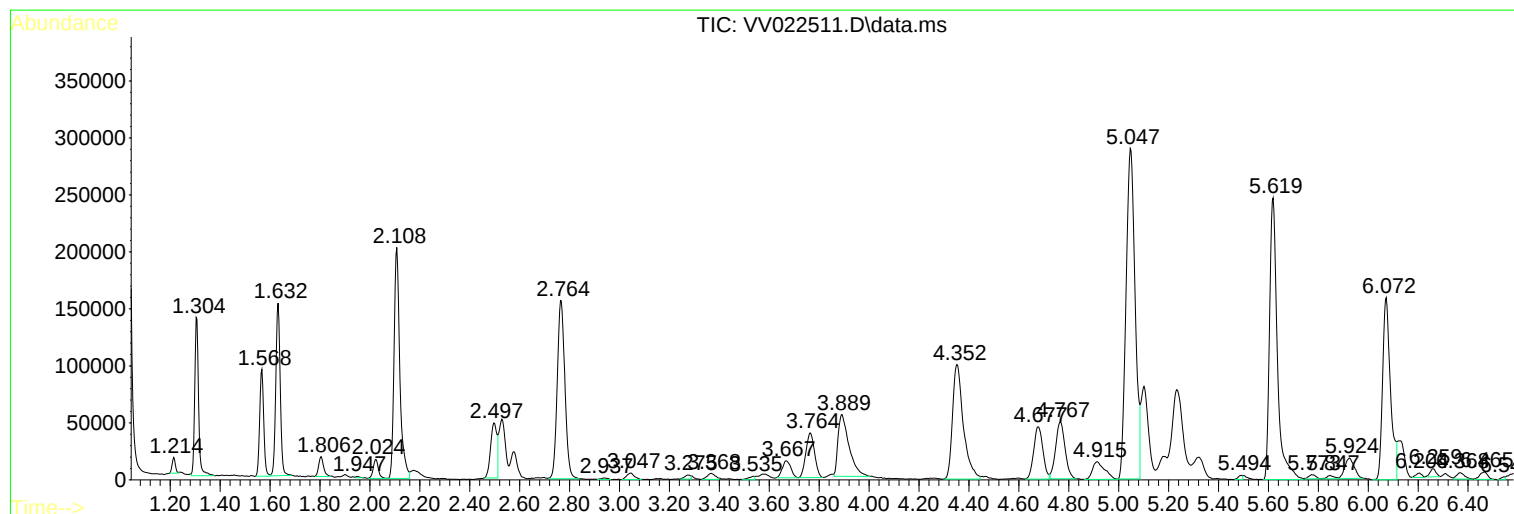
Sum of corrected areas: 8176959

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Instrument :  
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F4A29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
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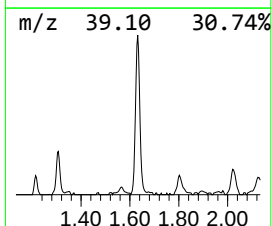
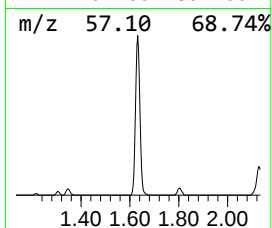
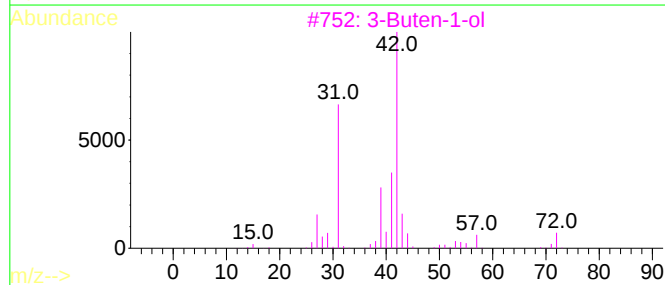
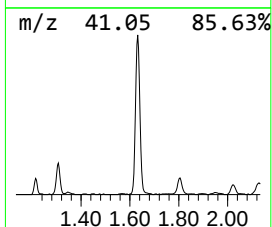
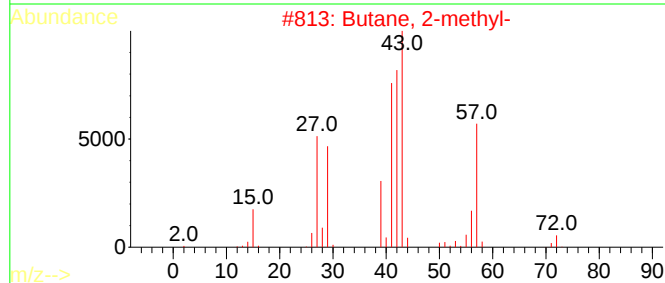
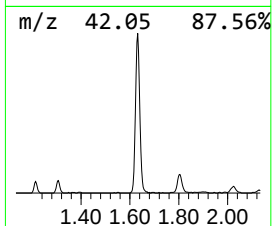
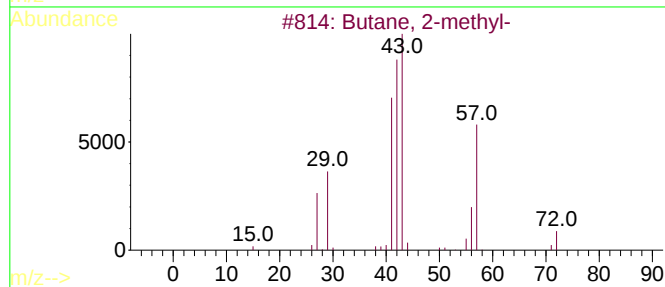
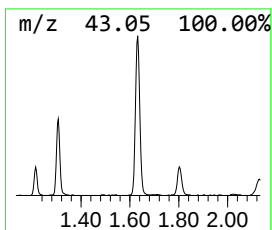
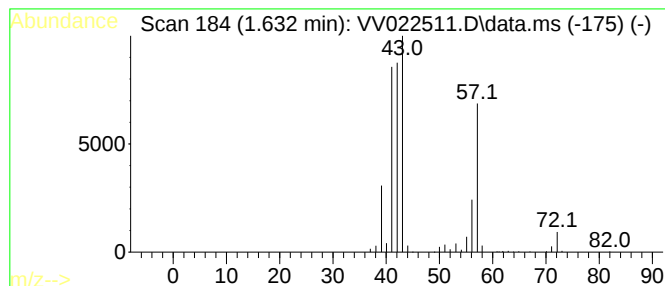
Instrument :  
MSVOA\_V  
ClientSampleId :  
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
Quant Title : VOC Analysis

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

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

| R.T.      | EstConc           | Area   | Relative to ISTD    | R.T.        |      |
|-----------|-------------------|--------|---------------------|-------------|------|
| 1.632     | 17.38 ug/L        | 187163 | 1,4-Difluorobenzene | 5.619       |      |
| Hit# of 5 | Tentative ID      | MW     | MolForm             | CAS#        | Qual |
| 1         | Butane, 2-methyl- | 72     | C5H12               | 000078-78-4 | 87   |
| 2         | Butane, 2-methyl- | 72     | C5H12               | 000078-78-4 | 86   |
| 3         | 3-Buten-1-ol      | 72     | C4H8O               | 000627-27-0 | 47   |
| 4         | Pentane           | 72     | C5H12               | 000109-66-0 | 37   |
| 5         | 1-Butene          | 56     | C4H8                | 000106-98-9 | 14   |



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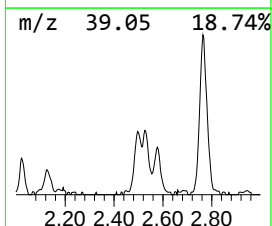
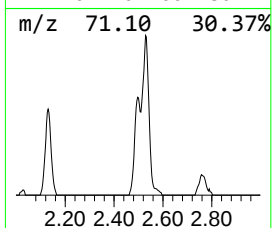
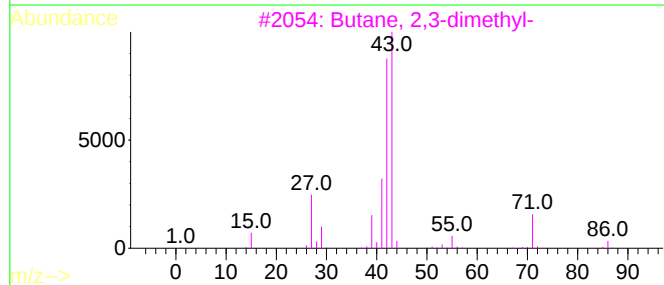
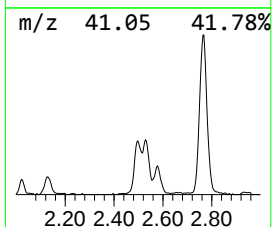
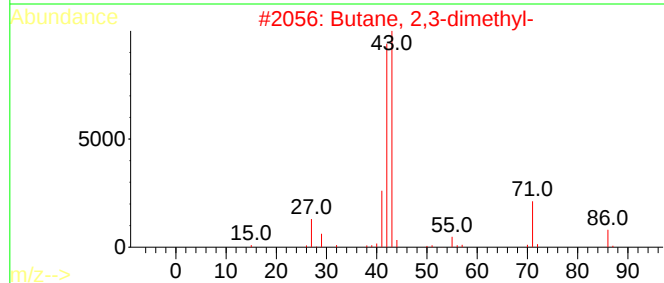
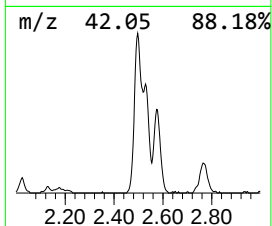
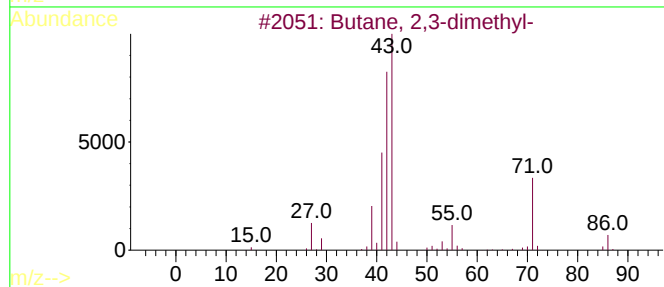
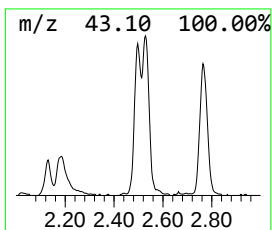
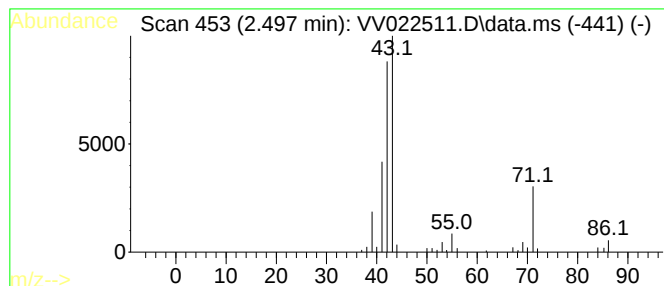
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc   | Area  | Relative to ISTD    | R.T.  |
|-------|-----------|-------|---------------------|-------|
| 2.497 | 8.29 ug/L | 89273 | 1,4-Difluorobenzene | 5.619 |

| Hit# | of 5 | Tentative ID          | MW | MolForm | CAS#        | Qual |
|------|------|-----------------------|----|---------|-------------|------|
| 1    |      | Butane, 2,3-dimethyl- | 86 | C6H14   | 000079-29-8 | 91   |
| 2    |      | Butane, 2,3-dimethyl- | 86 | C6H14   | 000079-29-8 | 86   |
| 3    |      | Butane, 2,3-dimethyl- | 86 | C6H14   | 000079-29-8 | 80   |
| 4    |      | Butane, 2,3-dimethyl- | 86 | C6H14   | 000079-29-8 | 80   |
| 5    |      | Butane, 2,3-dimethyl- | 86 | C6H14   | 000079-29-8 | 78   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
Quant Title : VOC Analysis

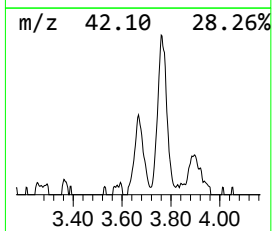
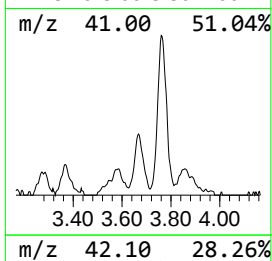
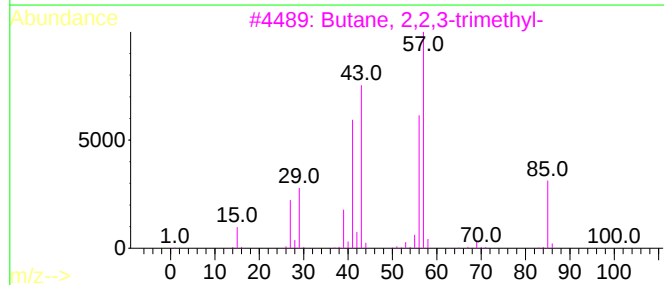
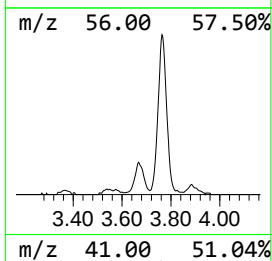
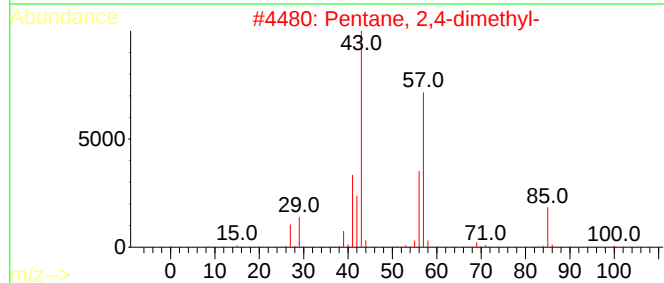
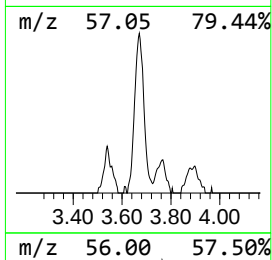
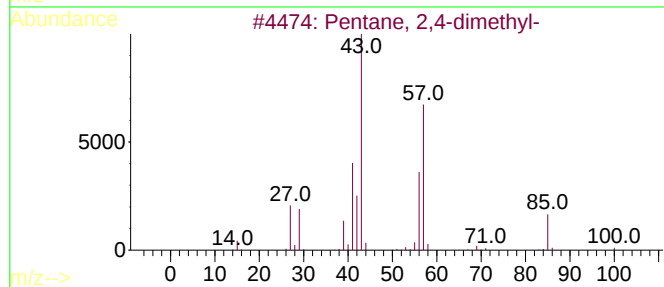
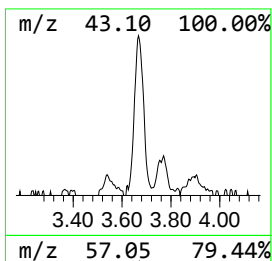
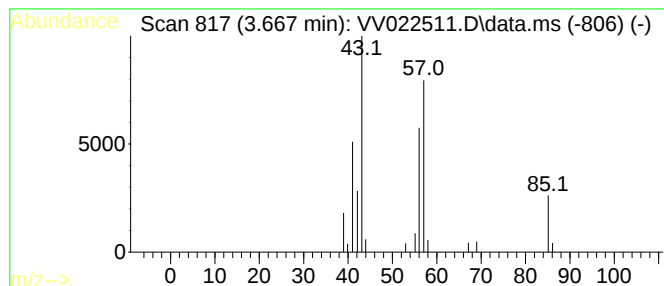
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc   | Area  | Relative to ISTD    | R.T.  |
|-------|-----------|-------|---------------------|-------|
| 3.667 | 3.26 ug/L | 35138 | 1,4-Difluorobenzene | 5.619 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentane, 2,4-dimethyl-   | 100 | C7H16   | 000108-08-7 | 74   |
| 2    |    |   | Pentane, 2,4-dimethyl-   | 100 | C7H16   | 000108-08-7 | 64   |
| 3    |    |   | Butane, 2,2,3-trimethyl- | 100 | C7H16   | 000464-06-2 | 59   |
| 4    |    |   | Pentane, 2,2-dimethyl-   | 100 | C7H16   | 000590-35-2 | 50   |
| 5    |    |   | Pentane, 2,2-dimethyl-   | 100 | C7H16   | 000590-35-2 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
Quant Title : VOC Analysis

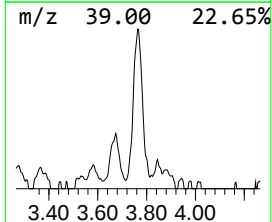
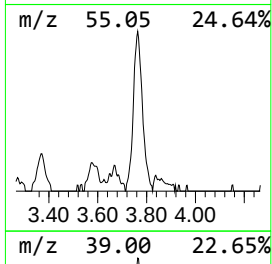
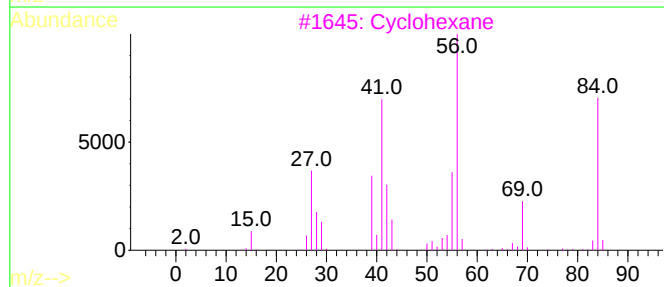
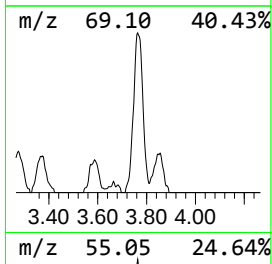
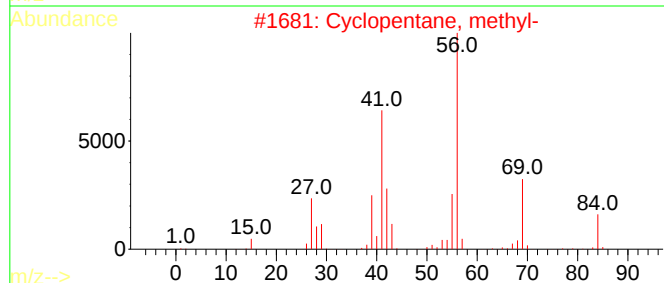
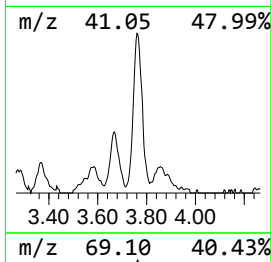
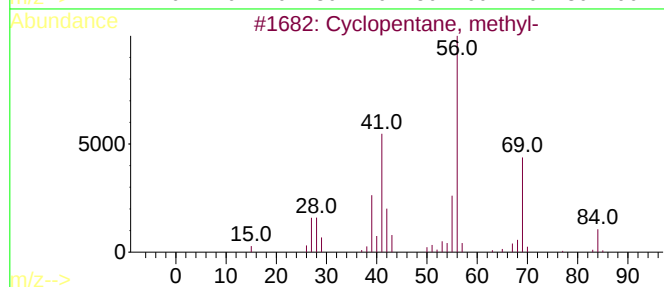
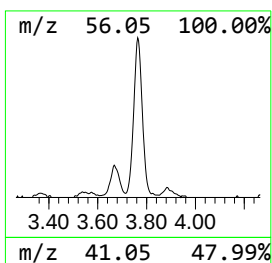
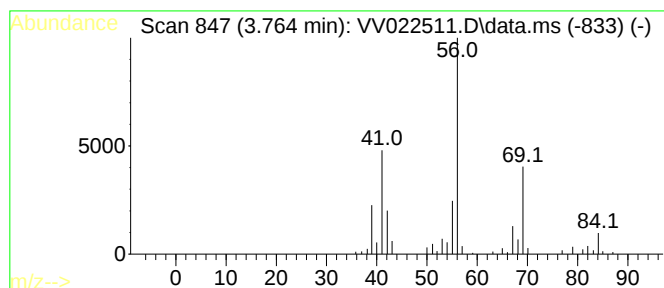
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 (DEL) Alkane: Cyclic3.764 Concentration Rank 5

| R.T.  | EstConc   | Area  | Relative to ISTD    | R.T.  |
|-------|-----------|-------|---------------------|-------|
| 3.764 | 8.80 ug/L | 94752 | 1,4-Difluorobenzene | 5.619 |

| Hit# | of 5 | Tentative ID            | MW | MolForm | CAS#        | Qual |
|------|------|-------------------------|----|---------|-------------|------|
| 1    |      | Cyclopentane, methyl-   | 84 | C6H12   | 000096-37-7 | 90   |
| 2    |      | Cyclopentane, methyl-   | 84 | C6H12   | 000096-37-7 | 83   |
| 3    |      | Cyclohexane             | 84 | C6H12   | 000110-82-7 | 80   |
| 4    |      | 1H-Tetrazole, 5-methyl- | 84 | C2H4N4  | 004076-36-2 | 72   |
| 5    |      | Cyclohexane             | 84 | C6H12   | 000110-82-7 | 72   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

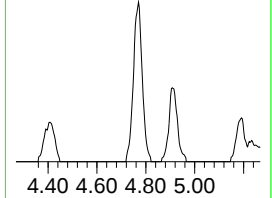
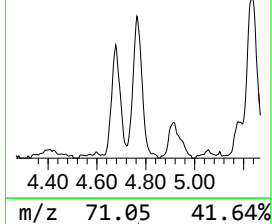
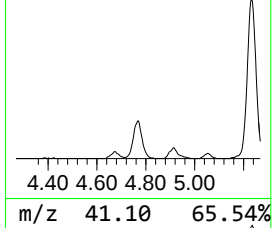
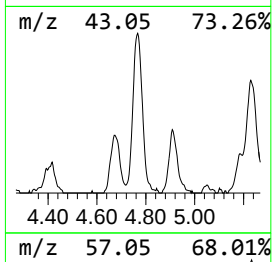
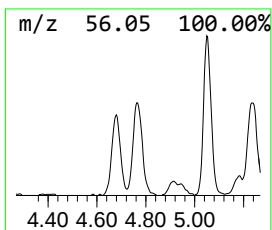
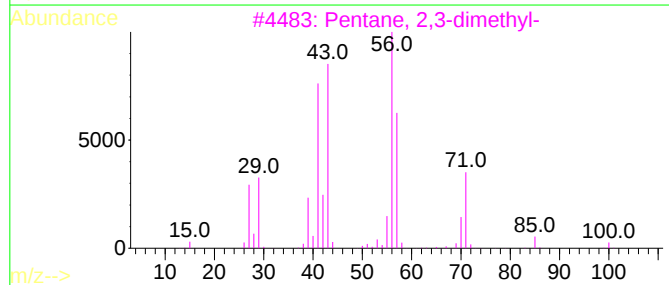
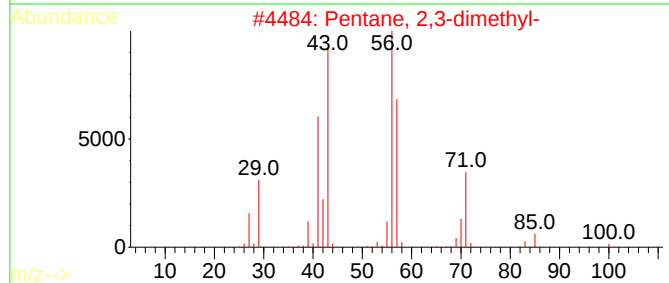
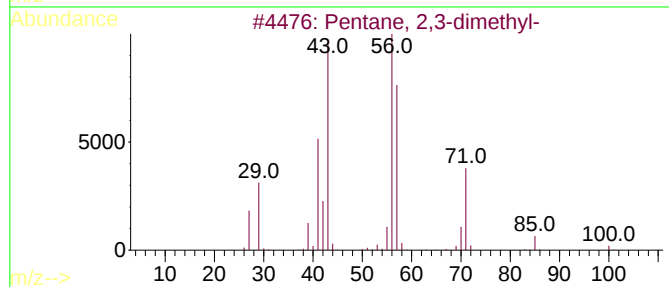
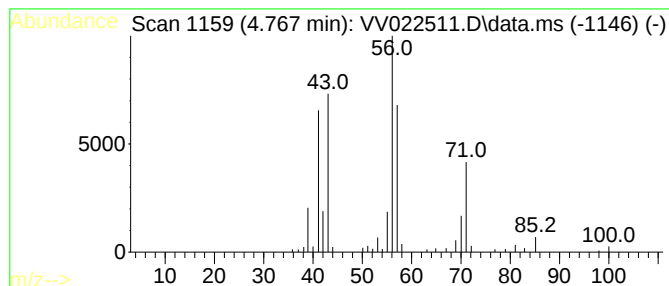
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Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 4.767 | 12.16 ug/L | 131013 | 1,4-Difluorobenzene | 5.619 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentane, 2,3-dimethyl-       | 100 | C7H16   | 000565-59-3 | 91   |
| 2    |    |   | Pentane, 2,3-dimethyl-       | 100 | C7H16   | 000565-59-3 | 90   |
| 3    |    |   | Pentane, 2,3-dimethyl-       | 100 | C7H16   | 000565-59-3 | 83   |
| 4    |    |   | Pentane, 2,3-dimethyl-       | 100 | C7H16   | 000565-59-3 | 80   |
| 5    |    |   | Hexane, 2,2,5,5-tetramethyl- | 142 | C10H22  | 001071-81-4 | 72   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

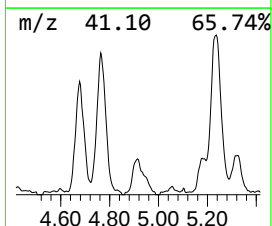
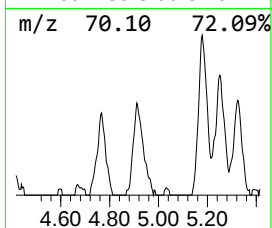
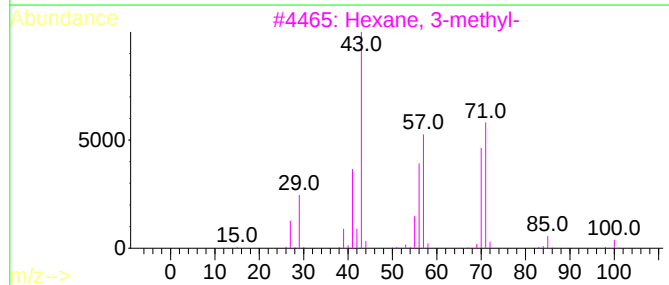
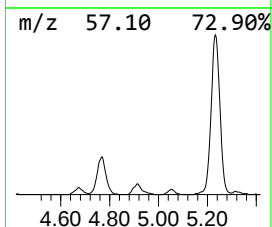
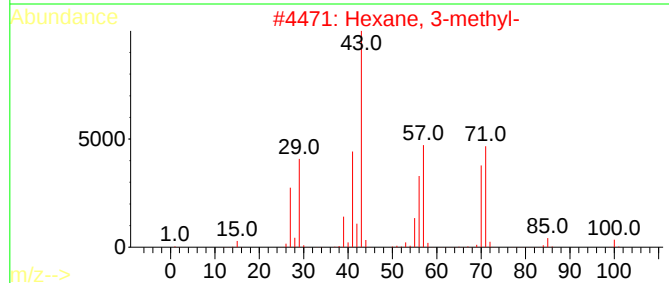
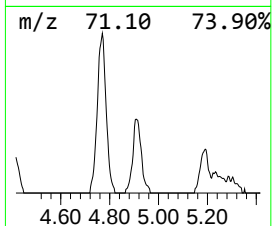
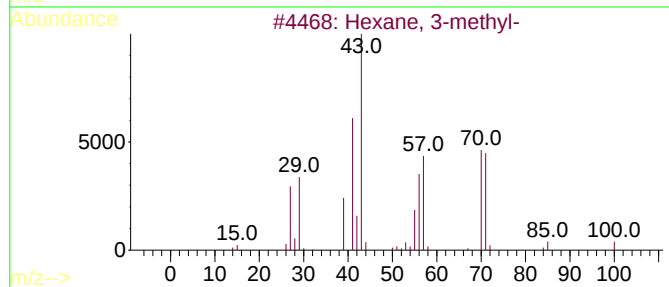
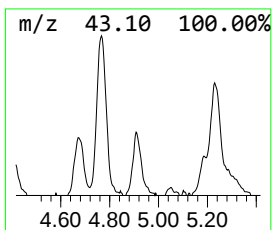
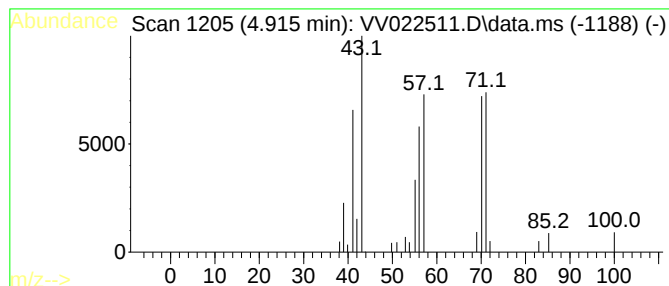
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Quant Title : VOC Analysis

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

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Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc   | Area  | Relative to ISTD    | R.T.  |
|-------|-----------|-------|---------------------|-------|
| 4.915 | 4.94 ug/L | 53255 | 1,4-Difluorobenzene | 5.619 |

| Hit# | of 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------|-----|---------|-------------|------|
| 1    |      | Hexane, 3-methyl-        | 100 | C7H16   | 000589-34-4 | 86   |
| 2    |      | Hexane, 3-methyl-        | 100 | C7H16   | 000589-34-4 | 83   |
| 3    |      | Hexane, 3-methyl-        | 100 | C7H16   | 000589-34-4 | 80   |
| 4    |      | Decane, 3,3,4-trimethyl- | 184 | C13H28  | 049622-18-6 | 59   |
| 5    |      | Hexane, 3,3,4-trimethyl- | 128 | C9H20   | 016747-31-2 | 53   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
Quant Title : VOC Analysis

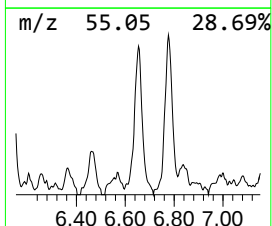
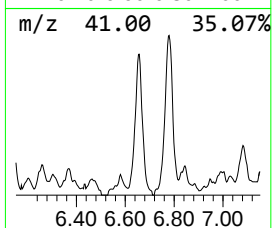
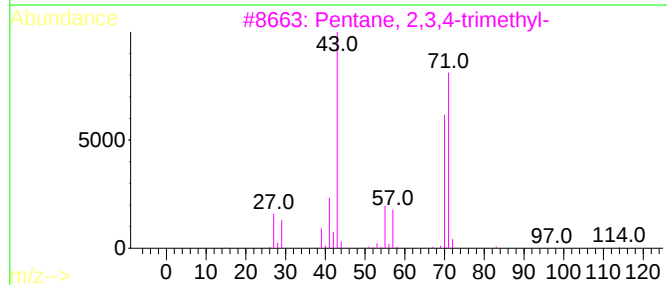
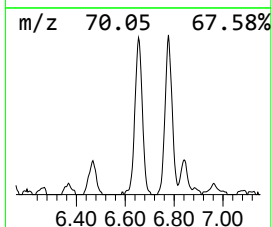
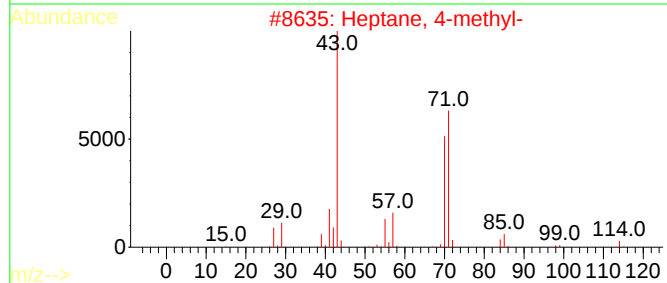
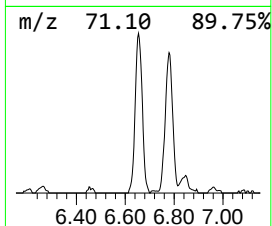
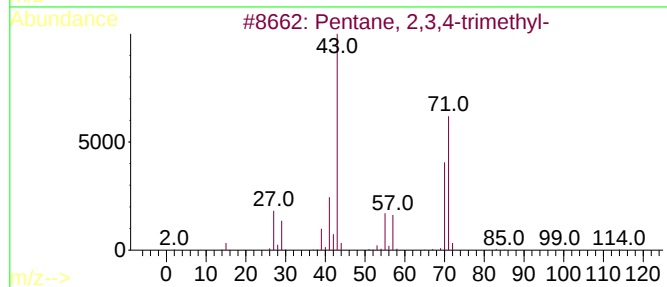
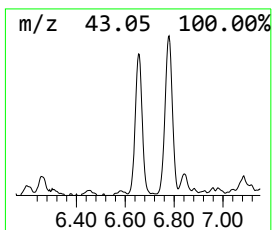
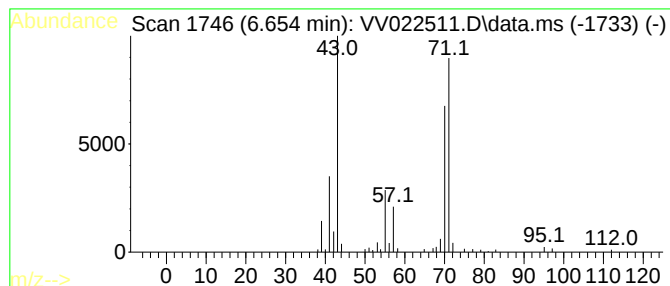
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc   | Area  | Relative to ISTD    | R.T.  |
|-------|-----------|-------|---------------------|-------|
| 6.654 | 8.55 ug/L | 92063 | 1,4-Difluorobenzene | 5.619 |

| Hit# | of 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------------|-----|---------|-------------|------|
| 1    |      | Pentane, 2,3,4-trimethyl- | 114 | C8H18   | 000565-75-3 | 86   |
| 2    |      | Heptane, 4-methyl-        | 114 | C8H18   | 000589-53-7 | 83   |
| 3    |      | Pentane, 2,3,4-trimethyl- | 114 | C8H18   | 000565-75-3 | 83   |
| 4    |      | Pentane, 2,3,4-trimethyl- | 114 | C8H18   | 000565-75-3 | 78   |
| 5    |      | Pentane, 3-ethyl-         | 100 | C7H16   | 000617-78-7 | 78   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
Quant Title : VOC Analysis

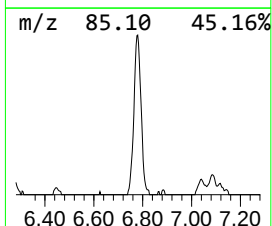
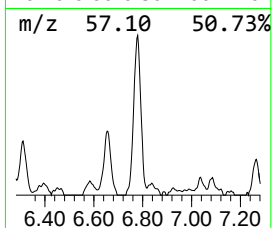
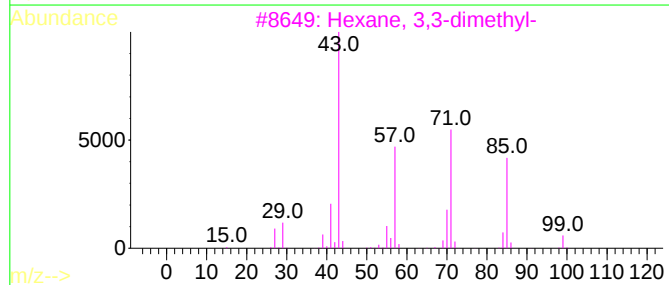
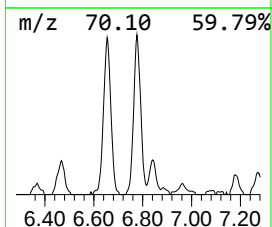
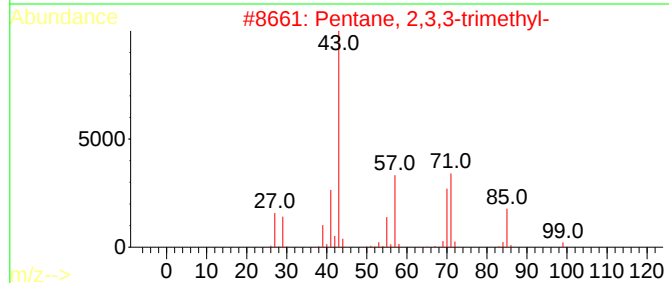
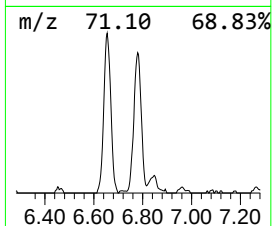
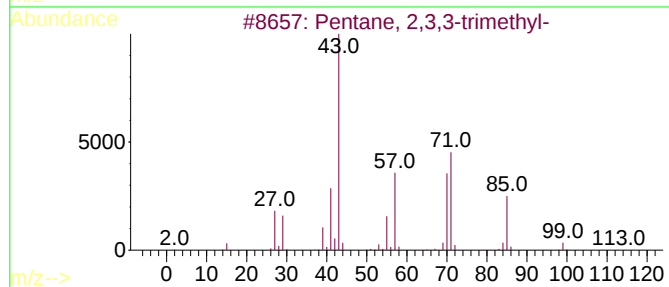
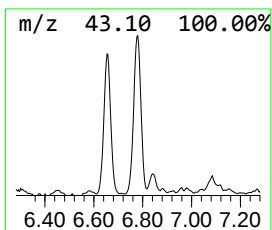
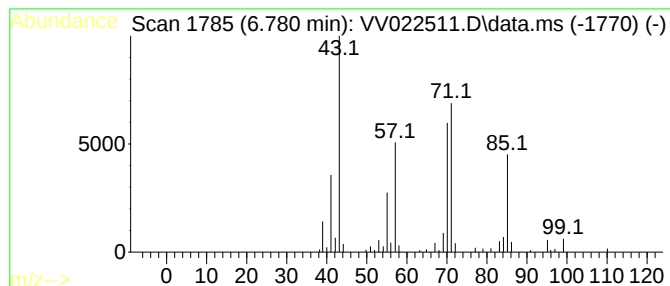
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD    | R.T.  |
|-------|------------|--------|---------------------|-------|
| 6.780 | 11.97 ug/L | 128927 | 1,4-Difluorobenzene | 5.619 |

| Hit# | of 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------------|-----|---------|-------------|------|
| 1    |      | Pentane, 2,3,3-trimethyl- | 114 | C8H18   | 000560-21-4 | 90   |
| 2    |      | Pentane, 2,3,3-trimethyl- | 114 | C8H18   | 000560-21-4 | 83   |
| 3    |      | Hexane, 3,3-dimethyl-     | 114 | C8H18   | 000563-16-6 | 64   |
| 4    |      | Heptane, 4-methyl-        | 114 | C8H18   | 000589-53-7 | 53   |
| 5    |      | Hexane, 3-methyl-         | 100 | C7H16   | 000589-34-4 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

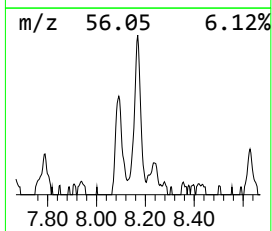
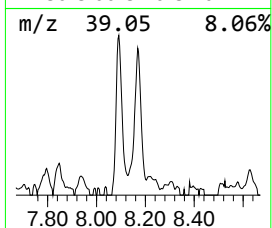
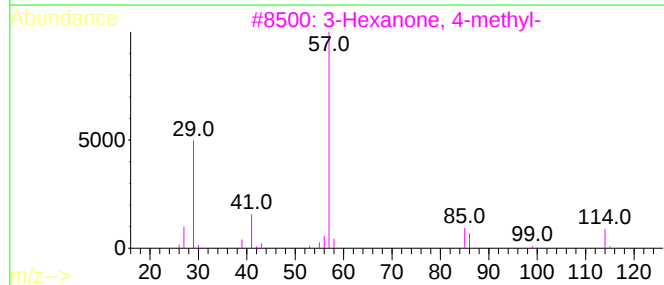
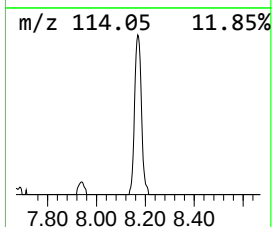
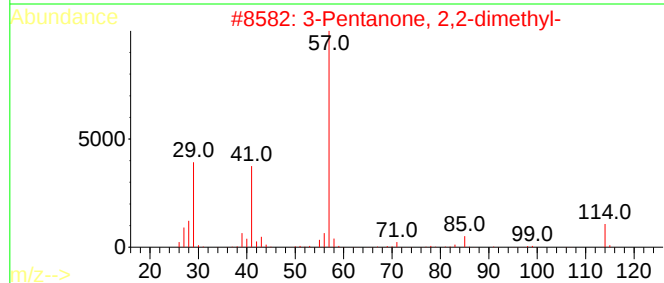
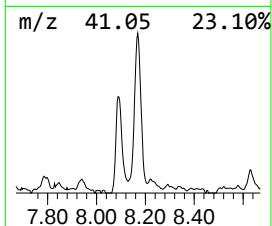
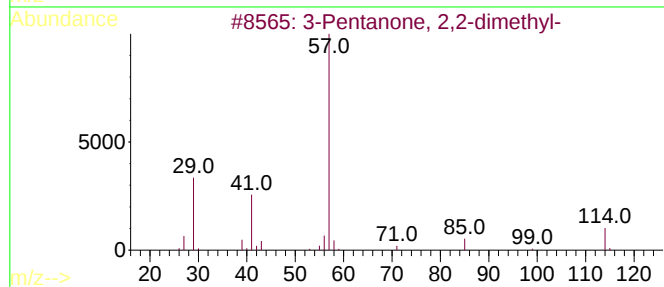
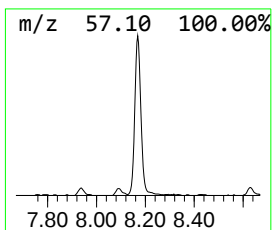
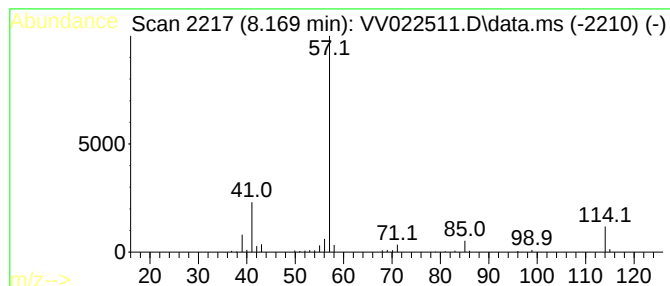
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Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 10 3-Pentanone, 2,2-dimethyl- Concentration Rank 8

| R.T.  | EstConc   | Area  | Relative to ISTD | R.T.  |
|-------|-----------|-------|------------------|-------|
| 8.169 | 8.22 ug/L | 96201 | Chlorobenzene-d5 | 8.854 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | 3-Pentanone, 2,2-dimethyl- | 114 | C7H14O  | 000564-04-5 | 64   |
| 2    |    |   | 3-Pentanone, 2,2-dimethyl- | 114 | C7H14O  | 000564-04-5 | 59   |
| 3    |    |   | 3-Hexanone, 4-methyl-      | 114 | C7H14O  | 017042-16-9 | 50   |
| 4    |    |   | 3-Heptanone                | 114 | C7H14O  | 000106-35-4 | 47   |
| 5    |    |   | Aluminum, triethyl-        | 114 | C6H15Al | 000097-93-8 | 40   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

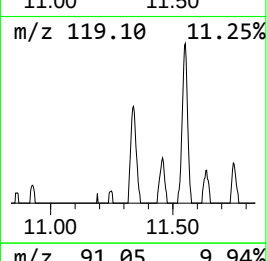
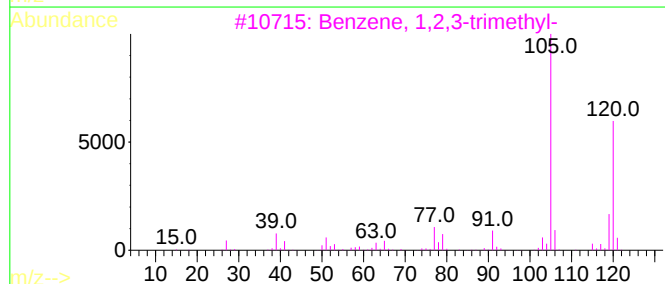
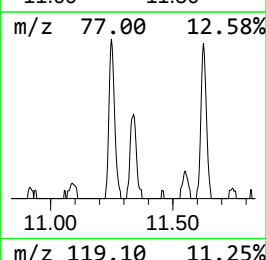
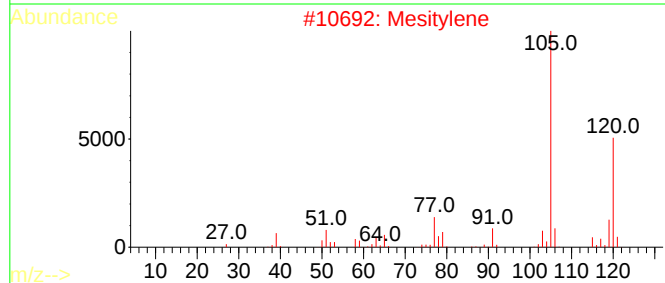
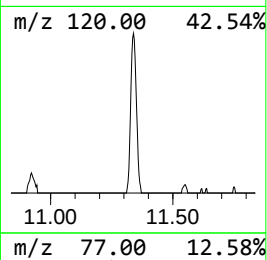
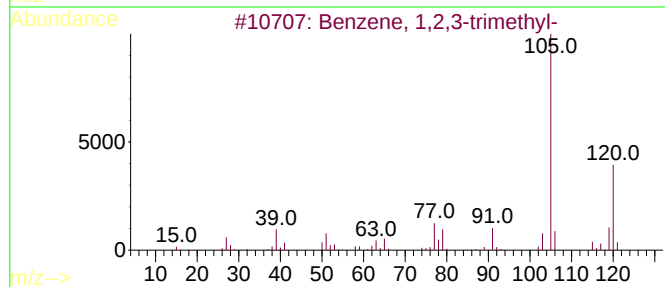
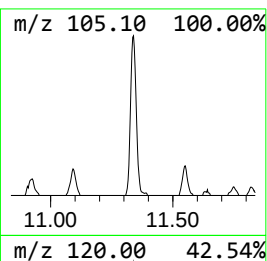
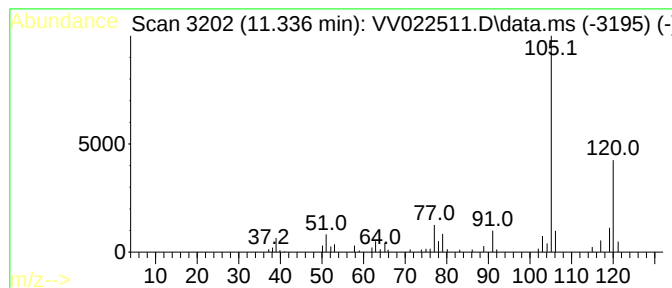
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Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 11 Mesitylene Concentration Rank 13

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 11.336 | 3.78 ug/L | 44660 | 1,4-Dichlorobenzene-d4 | 11.249 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 95   |
| 3    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 94   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 5    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
Quant Title : VOC Analysis

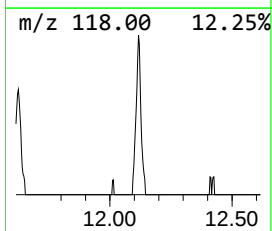
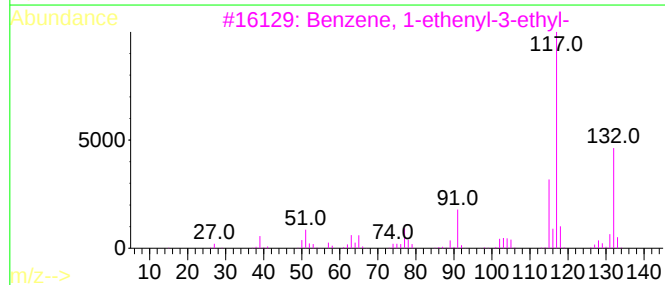
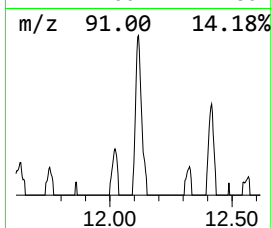
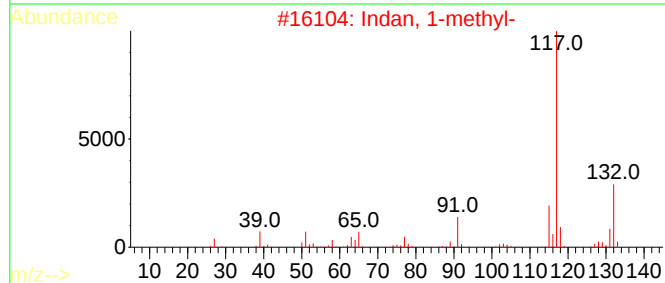
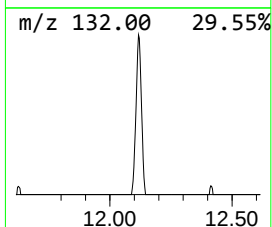
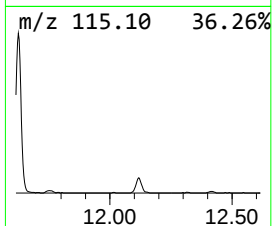
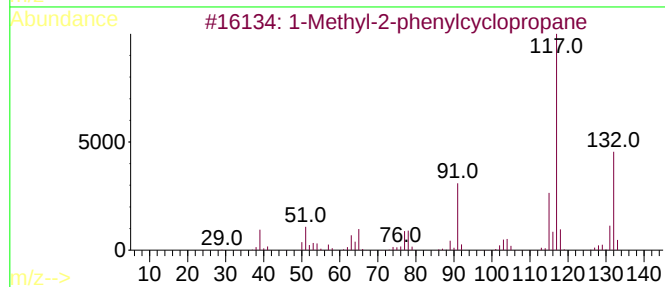
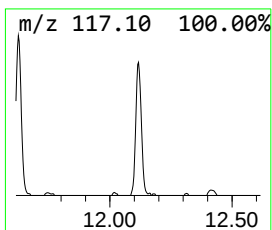
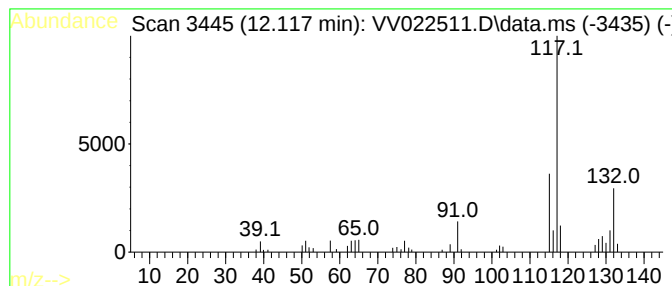
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 1-Methyl-2-phenylcyclopropane Concentration Rank 11

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 12.117 | 4.22 ug/L | 49846 | 1,4-Dichlorobenzene-d4 | 11.249 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Methyl-2-phenylcyclopropane       | 132 | C10H12  | 003145-76-4 | 90   |
| 2    |    |   | Indan, 1-methyl-                    | 132 | C10H12  | 000767-58-8 | 87   |
| 3    |    |   | Benzene, 1-ethenyl-3-ethyl-         | 132 | C10H12  | 007525-62-4 | 86   |
| 4    |    |   | Benzene, 1-methyl-2-(2-propenyl)-   | 132 | C10H12  | 001587-04-8 | 86   |
| 5    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 86   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

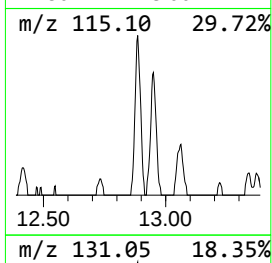
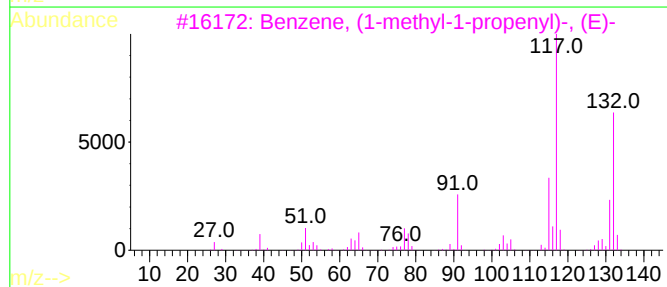
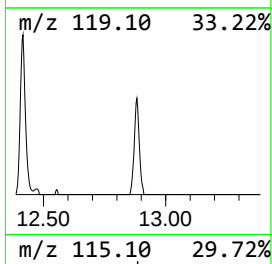
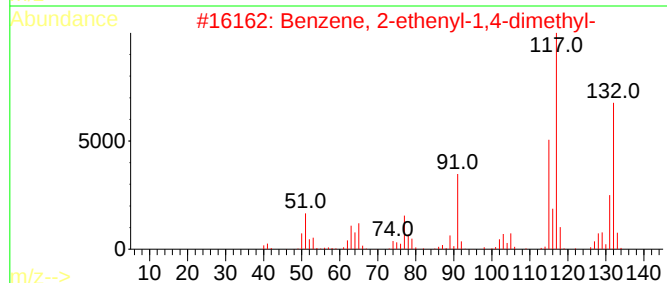
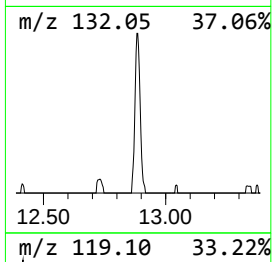
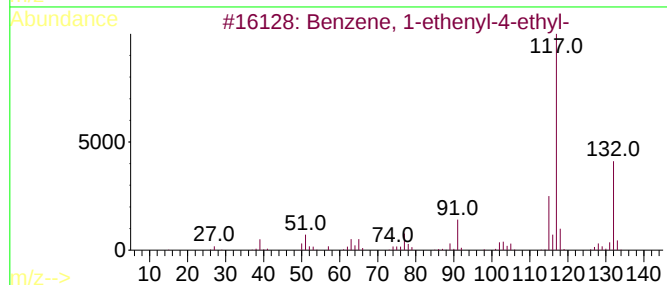
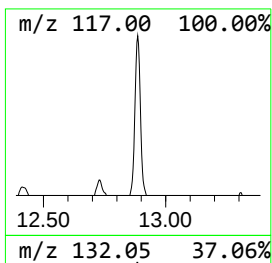
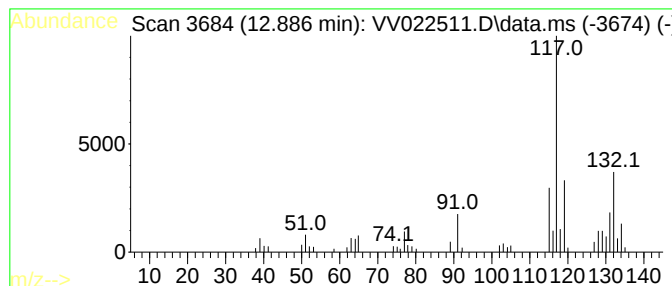
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Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 13 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 12

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 12.886 | 4.07 ug/L | 48132 | 1,4-Dichlorobenzene-d4 | 11.249 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethenyl-4-ethyl-         | 132 | C10H12  | 003454-07-7 | 93   |
| 2    |    |   | Benzene, 2-ethenyl-1,4-dimethyl-    | 132 | C10H12  | 002039-89-6 | 87   |
| 3    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 76   |
| 4    |    |   | Benzene, 2-butenyl-                 | 132 | C10H12  | 001560-06-1 | 76   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-2-methyl-    | 132 | C10H12  | 000824-63-5 | 76   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

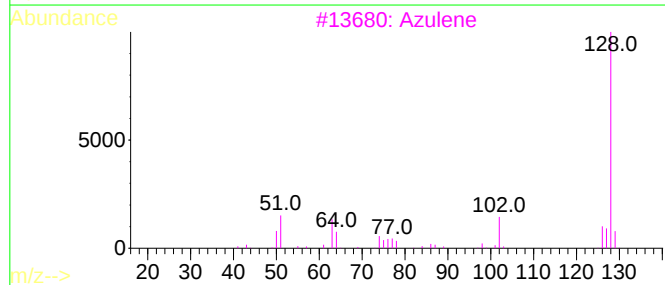
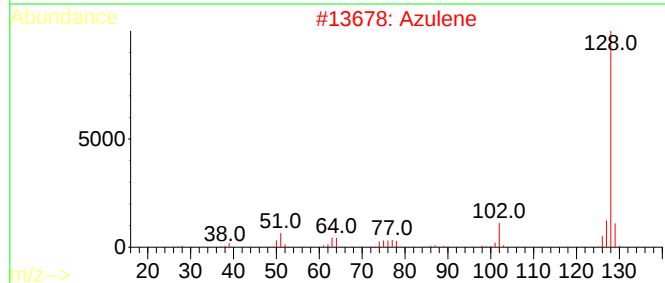
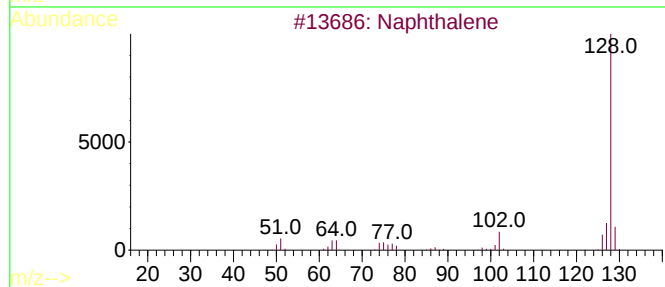
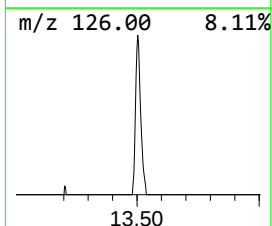
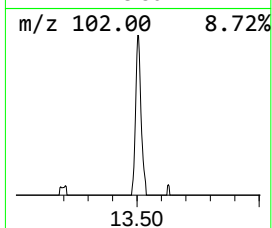
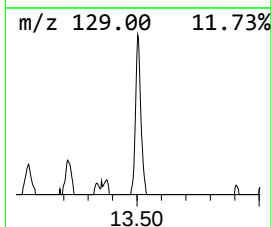
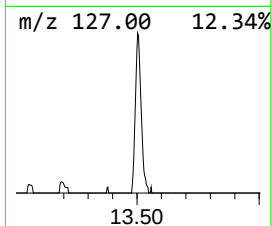
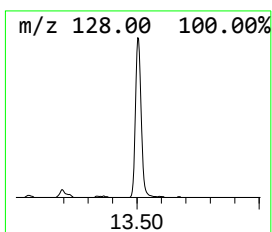
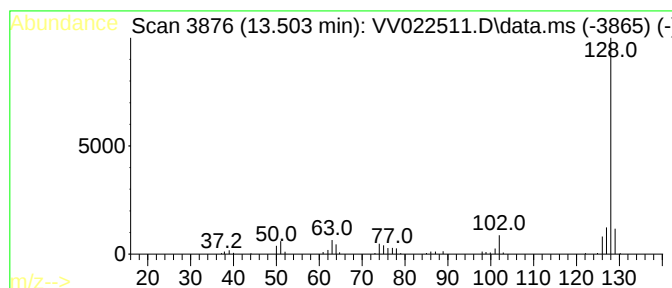
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Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 14 Azulene Concentration Rank 9

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 13.503 | 6.79 ug/L | 80171 | 1,4-Dichlorobenzene-d4 | 11.249 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene             | 128 | C10H8   | 000091-20-3 | 95   |
| 2    |    |   | Azulene                 | 128 | C10H8   | 000275-51-4 | 93   |
| 3    |    |   | Azulene                 | 128 | C10H8   | 000275-51-4 | 91   |
| 4    |    |   | Azulene                 | 128 | C10H8   | 000275-51-4 | 91   |
| 5    |    |   | 1H-Indene, 1-methylene- | 128 | C10H8   | 002471-84-3 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
Data File : VV022511.D  
Acq On : 02 Oct 2021 15:19  
Operator : SY/MD  
Sample : M3996-14  
Misc : 5.0mL/MSVOA\_V/WATER  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
F4A29

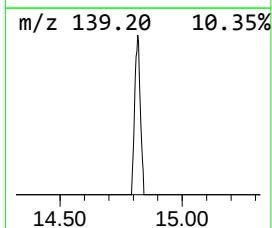
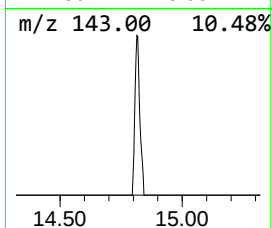
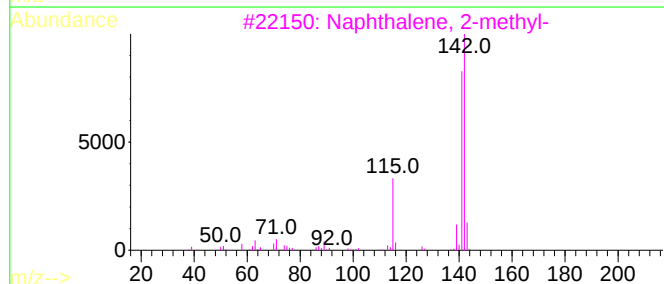
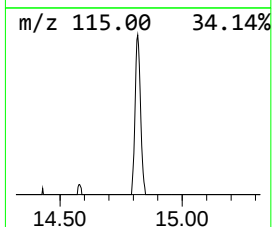
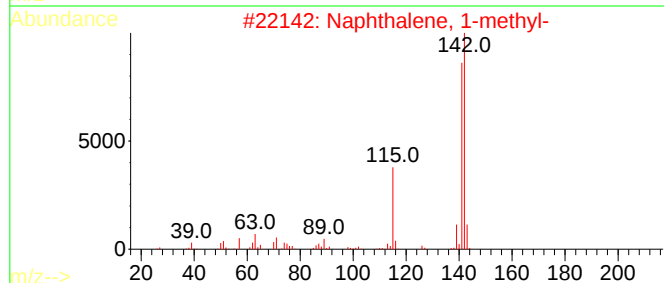
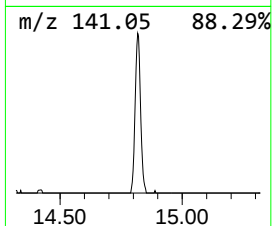
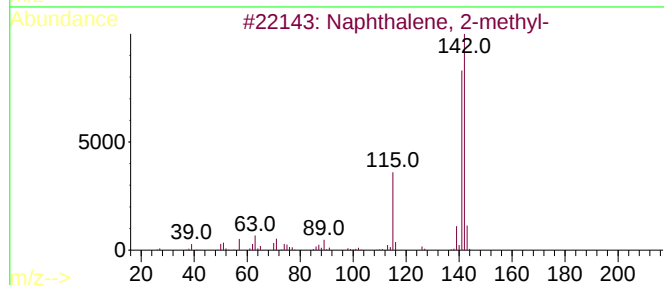
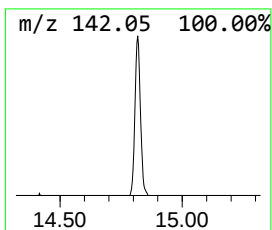
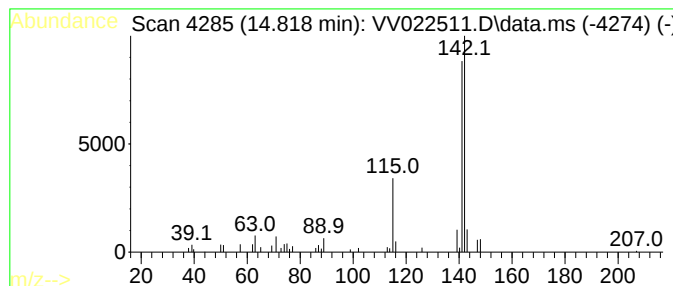
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Quant Title : VOC Analysis

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

\*\*\*\*\*  
Peak Number 15 Naphthalene, 2-methyl- Concentration Rank 14

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 14.818 | 3.71 ug/L | 43824 | 1,4-Dichlorobenzene-d4 | 11.249 |

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV100221\  
 Data File : VV022511.D  
 Acq On : 02 Oct 2021 15:19  
 Operator : SY/MD  
 Sample : M3996-14  
 Misc : 5.0mL/MSVOA\_V/WATER  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 MSVOA\_V  
 ClientSampleId :  
 F4A29

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVLM100121WMA.M  
 Quant Title : VOC Analysis

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |        |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|--------|------|
|                    |        |         |       |          | #                     | RT     | Resp   | Conc |
| (DEL) Alkane: S... | 1.632  | 17.4    | ug/L  | 187163   | 1                     | 5.619  | 538537 | 50.0 |
| (DEL) Alkane: S... | 2.497  | 8.3     | ug/L  | 89273    | 1                     | 5.619  | 538537 | 50.0 |
| (DEL) Alkane: S... | 3.667  | 3.3     | ug/L  | 35138    | 1                     | 5.619  | 538537 | 50.0 |
| (DEL) Alkane: C... | 3.764  | 8.8     | ug/L  | 94752    | 1                     | 5.619  | 538537 | 50.0 |
| (DEL) Alkane: S... | 4.767  | 12.2    | ug/L  | 131013   | 1                     | 5.619  | 538537 | 50.0 |
| (DEL) Alkane: S... | 4.915  | 4.9     | ug/L  | 53255    | 1                     | 5.619  | 538537 | 50.0 |
| (DEL) Alkane: S... | 6.654  | 8.6     | ug/L  | 92063    | 1                     | 5.619  | 538537 | 50.0 |
| (DEL) Alkane: S... | 6.780  | 12.0    | ug/L  | 128927   | 1                     | 5.619  | 538537 | 50.0 |
| 3-Pentanone, 2,... | 8.169  | 8.2     | ug/L  | 96201    | 2                     | 8.854  | 585453 | 50.0 |
| Mesitylene         | 11.336 | 3.8     | ug/L  | 44660    | 3                     | 11.249 | 590583 | 50.0 |
| 1-Methyl-2-phen... | 12.117 | 4.2     | ug/L  | 49846    | 3                     | 11.249 | 590583 | 50.0 |
| Benzene, 1-ethe... | 12.886 | 4.1     | ug/L  | 48132    | 3                     | 11.249 | 590583 | 50.0 |
| Azulene            | 13.503 | 6.8     | ug/L  | 80171    | 3                     | 11.249 | 590583 | 50.0 |
| Naphthalene, 2-... | 14.818 | 3.7     | ug/L  | 43824    | 3                     | 11.249 | 590583 | 50.0 |