

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
 Data File : VV028133.D  
 Acq On : 27 Sep 2022 12:40  
 Operator : SY/MD  
 Sample : N4820-02  
 Misc : 25.0mL/MSVOA\_V/WATER  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 MSVOA\_V  
 ClientSampleId :  
 CB8L0

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.1

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\SFAMVTR092222WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VV028133.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         | 1.298       | 73            | 80          | 93           | rVB      | 99794          | 108841        | 2.17%           | 0.394%        |
| 2         | 1.561       | 154           | 162         | 171          | rBV      | 85060          | 98518         | 1.97%           | 0.357%        |
| 3         | 2.098       | 318           | 329         | 343          | rBV      | 161964         | 232537        | 4.65%           | 0.842%        |
| 4         | 3.921       | 880           | 896         | 932          | rBV2     | 62903          | 336963        | 6.73%           | 1.221%        |
| 5         | 4.342       | 1010          | 1027        | 1053         | rBV2     | 121556         | 324518        | 6.48%           | 1.176%        |
| 6         | 5.040       | 1228          | 1244        | 1279         | rBV3     | 270565         | 690801        | 13.80%          | 2.502%        |
| 7         | 5.613       | 1410          | 1422        | 1449         | rBV      | 205315         | 421980        | 8.43%           | 1.529%        |
| 8         | 6.066       | 1549          | 1563        | 1577         | rBV      | 157071         | 332309        | 6.64%           | 1.204%        |
| 9         | 6.982       | 1838          | 1848        | 1873         | rBV      | 70334          | 141326        | 2.82%           | 0.512%        |
| 10        | 7.313       | 1938          | 1951        | 1966         | rBV      | 274313         | 485311        | 9.70%           | 1.758%        |
| 11        | 7.622       | 2030          | 2047        | 2067         | rBV      | 44997          | 90411         | 1.81%           | 0.328%        |
| 12        | 8.088       | 2179          | 2192        | 2224         | rBV      | 350148         | 713512        | 14.26%          | 2.585%        |
| 13        | 8.239       | 2226          | 2239        | 2252         | rBV6     | 29571          | 81001         | 1.62%           | 0.293%        |
| 14        | 8.876       | 2417          | 2437        | 2464         | rBV2     | 1003684        | 2176336       | 43.49%          | 7.884%        |
| 15        | 9.130       | 2503          | 2516        | 2531         | rVB9     | 29411          | 73002         | 1.46%           | 0.264%        |
| 16        | 9.799       | 2712          | 2724        | 2740         | rBV9     | 21945          | 57655         | 1.15%           | 0.209%        |
| 17        | 9.927       | 2752          | 2764        | 2777         | rBV      | 323536         | 508938        | 10.17%          | 1.844%        |
| 18        | 10.214      | 2843          | 2853        | 2869         | rVB      | 132852         | 230460        | 4.61%           | 0.835%        |
| 19        | 10.349      | 2881          | 2895        | 2904         | rBV      | 119104         | 204335        | 4.08%           | 0.740%        |
| 20        | 10.734      | 3006          | 3015        | 3036         | rVB4     | 38234          | 82728         | 1.65%           | 0.300%        |
| 21        | 10.860      | 3044          | 3054        | 3062         | rBV2     | 37304          | 59669         | 1.19%           | 0.216%        |
| 22        | 11.053      | 3104          | 3114        | 3118         | rBV      | 170758         | 253389        | 5.06%           | 0.918%        |
| 23        | 11.085      | 3118          | 3124        | 3139         | rVB      | 273685         | 418829        | 8.37%           | 1.517%        |
| 24        | 11.243      | 3162          | 3173        | 3195         | rVB3     | 515220         | 1036563       | 20.71%          | 3.755%        |
| 25        | 11.390      | 3210          | 3219        | 3228         | rBV7     | 29663          | 60319         | 1.21%           | 0.218%        |
| 26        | 11.448      | 3229          | 3237        | 3252         | rVB      | 96163          | 151759        | 3.03%           | 0.550%        |
| 27        | 11.542      | 3252          | 3266        | 3278         | rBV      | 849336         | 1324019       | 26.46%          | 4.796%        |
| 28        | 11.622      | 3280          | 3291        | 3296         | rBV2     | 386628         | 627393        | 12.54%          | 2.273%        |
| 29        | 11.741      | 3319          | 3328        | 3344         | rVB      | 83273          | 127856        | 2.55%           | 0.463%        |
| 30        | 12.011      | 3396          | 3412        | 3432         | rBV      | 3118479        | 5004450       | 100.00%         | 18.128%       |
| 31        | 12.111      | 3433          | 3443        | 3453         | rBV      | 736622         | 1147861       | 22.94%          | 4.158%        |
| 32        | 12.294      | 3492          | 3500        | 3511         | rVB      | 71393          | 104355        | 2.09%           | 0.378%        |
| 33        | 12.410      | 3519          | 3536        | 3548         | rBV      | 1907016        | 2793885       | 55.83%          | 10.121%       |
| 34        | 12.545      | 3571          | 3578        | 3588         | rVB      | 120583         | 176327        | 3.52%           | 0.639%        |
| 35        | 12.657      | 3600          | 3613        | 3624         | rBV5     | 92475          | 276544        | 5.53%           | 1.002%        |
| 36        | 12.718      | 3624          | 3632        | 3650         | rVB      | 517395         | 795112        | 15.89%          | 2.880%        |

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Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

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Stop Thrs : 0

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
 Title : TRACE VOA SFAM1.0

|    |        |      |      |      |      |         |         |        |         |
|----|--------|------|------|------|------|---------|---------|--------|---------|
| 37 | 12.876 | 3666 | 3681 | 3691 | rBV2 | 1925384 | 3005925 | 60.07% | 10.889% |
| 38 | 12.921 | 3692 | 3695 | 3710 | rVB3 | 80909   | 129046  | 2.58%  | 0.467%  |
| 39 | 12.995 | 3711 | 3718 | 3727 | rVV  | 135926  | 184878  | 3.69%  | 0.670%  |
| 40 | 13.213 | 3778 | 3786 | 3793 | rVB2 | 79473   | 115357  | 2.31%  | 0.418%  |
| 41 | 13.268 | 3793 | 3803 | 3813 | rBV  | 316039  | 485845  | 9.71%  | 1.760%  |
| 42 | 13.329 | 3817 | 3822 | 3827 | rVV  | 62624   | 96414   | 1.93%  | 0.349%  |
| 43 | 13.365 | 3828 | 3833 | 3837 | rVV  | 98801   | 132397  | 2.65%  | 0.480%  |
| 44 | 13.397 | 3838 | 3843 | 3858 | rVB  | 150921  | 219850  | 4.39%  | 0.796%  |
| 45 | 13.500 | 3861 | 3875 | 3881 | rBV3 | 39820   | 91825   | 1.83%  | 0.333%  |
| 46 | 13.541 | 3881 | 3888 | 3900 | rVB4 | 58173   | 101083  | 2.02%  | 0.366%  |
| 47 | 13.664 | 3918 | 3926 | 3932 | rBV2 | 31390   | 51932   | 1.04%  | 0.188%  |
| 48 | 13.763 | 3951 | 3957 | 3969 | rVB2 | 45760   | 73351   | 1.47%  | 0.266%  |
| 49 | 13.905 | 3991 | 4001 | 4016 | rBV2 | 75564   | 138189  | 2.76%  | 0.501%  |
| 50 | 14.088 | 4048 | 4058 | 4070 | rBV5 | 85811   | 192760  | 3.85%  | 0.698%  |
| 51 | 14.226 | 4094 | 4101 | 4112 | rVV  | 45567   | 74202   | 1.48%  | 0.269%  |
| 52 | 14.287 | 4114 | 4120 | 4134 | rVB3 | 40066   | 68710   | 1.37%  | 0.249%  |
| 53 | 14.426 | 4155 | 4163 | 4177 | rVB7 | 36337   | 70667   | 1.41%  | 0.256%  |
| 54 | 14.628 | 4217 | 4226 | 4235 | rBV3 | 33165   | 57611   | 1.15%  | 0.209%  |
| 55 | 14.815 | 4274 | 4284 | 4295 | rBV4 | 50143   | 94503   | 1.89%  | 0.342%  |
| 56 | 15.529 | 4488 | 4506 | 4517 | rBV2 | 28391   | 77225   | 1.54%  | 0.280%  |
| 57 | 15.660 | 4538 | 4547 | 4567 | rVB2 | 59492   | 114120  | 2.28%  | 0.413%  |
| 58 | 15.805 | 4577 | 4592 | 4601 | rBV  | 97002   | 160552  | 3.21%  | 0.582%  |
| 59 | 16.480 | 4787 | 4802 | 4812 | rBV2 | 70404   | 119720  | 2.39%  | 0.434%  |

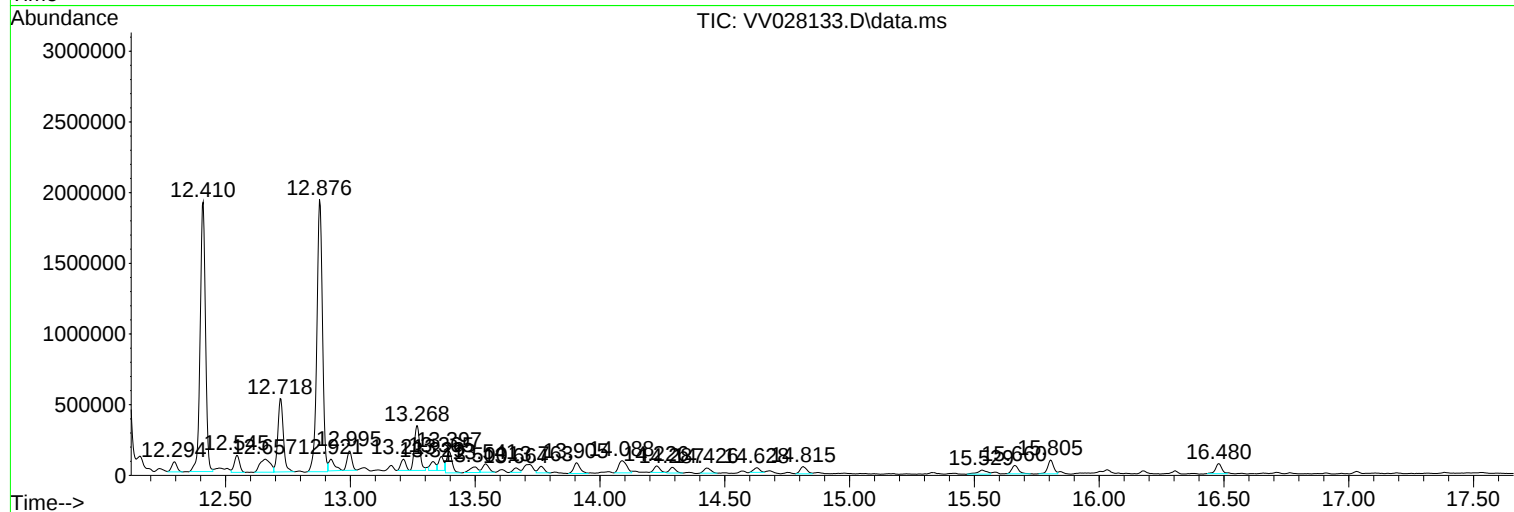
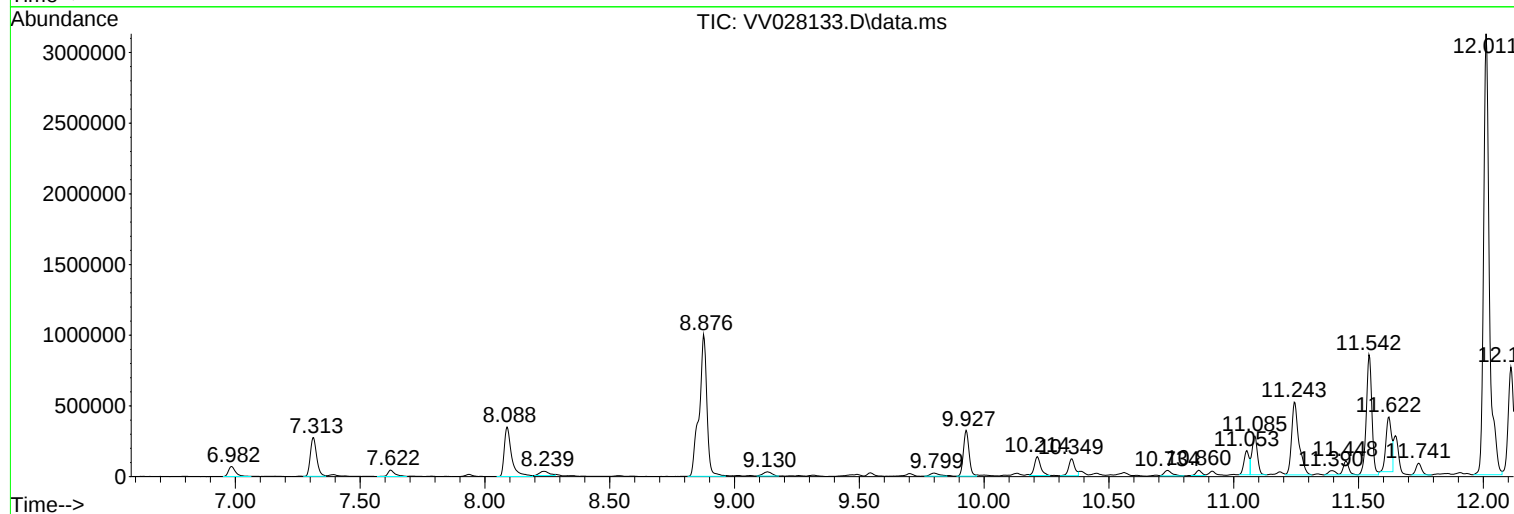
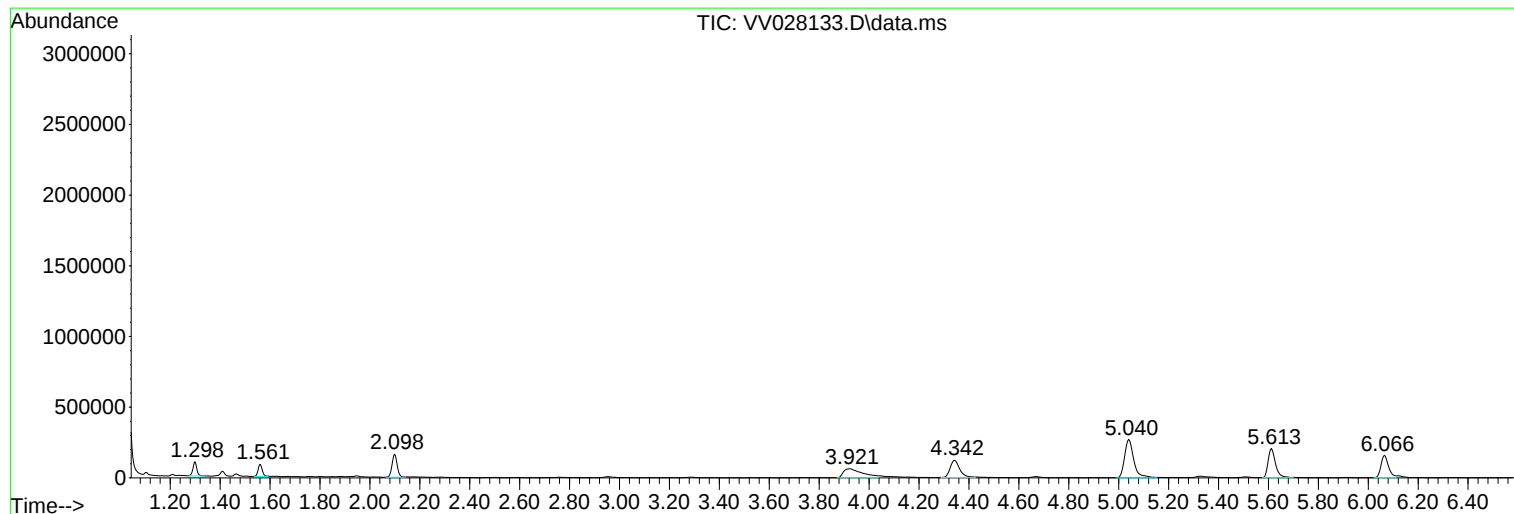
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ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

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



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Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

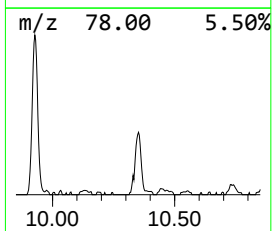
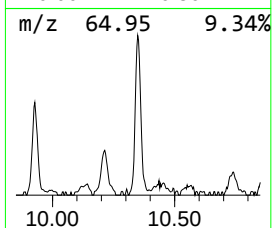
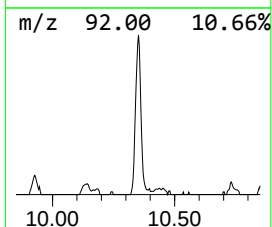
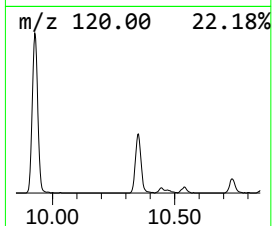
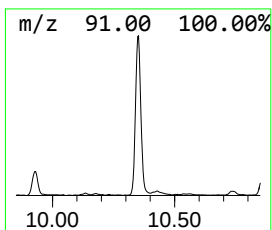
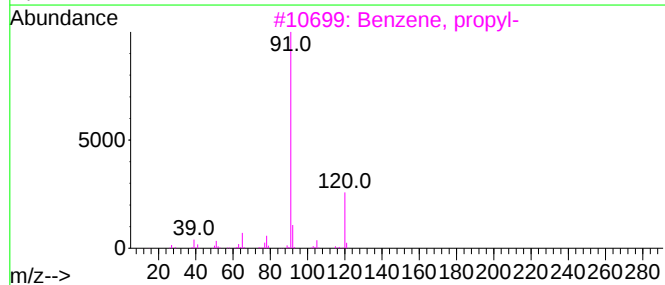
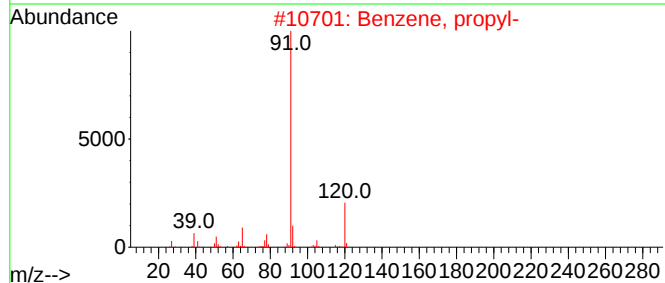
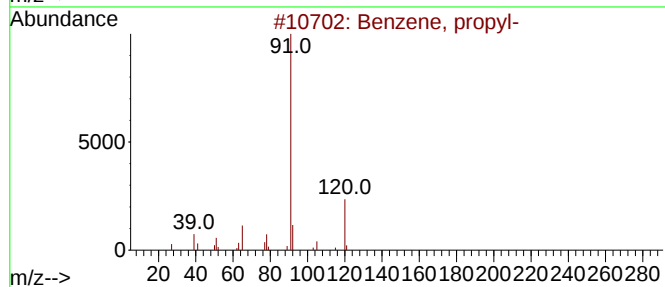
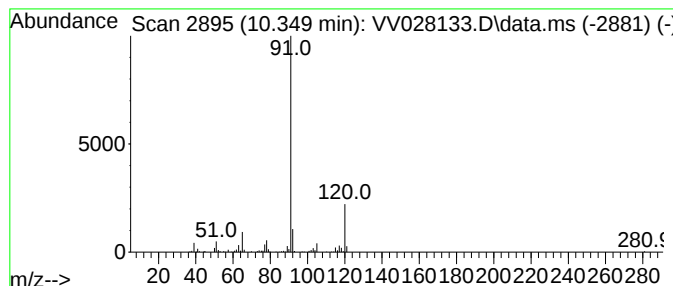
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 2 Benzene, propyl- Concentration Rank 13

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 10.349 | 0.99 ug/L | 204335 | 1,4-Dichlorobenzene-d4 | 11.246 |

| 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    |    |   | Benzene, propyl-                  | 120 | C9H12   | 000103-65-1 | 87   |
| 4    |    |   | N-Butylbenzylamine                | 163 | C11H17N | 002403-22-7 | 64   |
| 5    |    |   | Ethanol, 2-[(phenylmethyl)amino]- | 151 | C9H13NO | 000104-63-2 | 64   |



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Instrument :  
MSVOA\_V  
ClientSampleId :  
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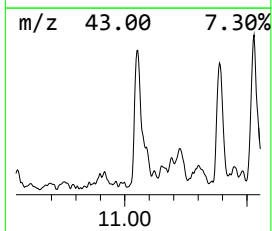
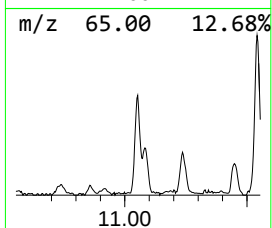
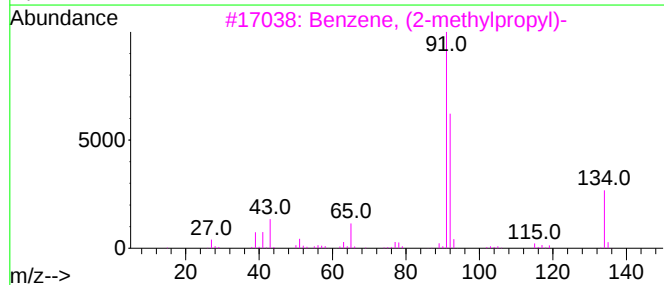
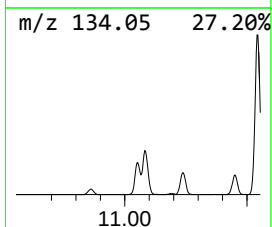
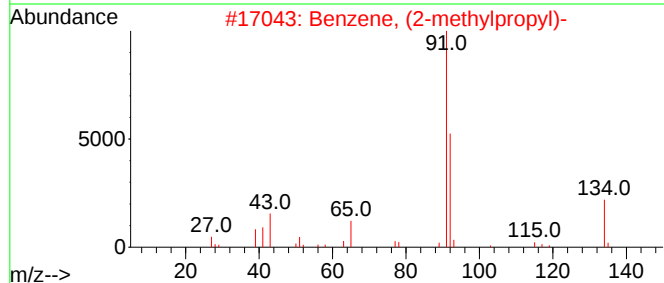
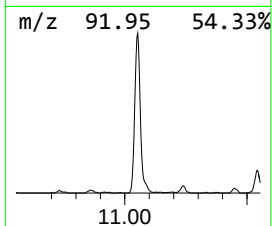
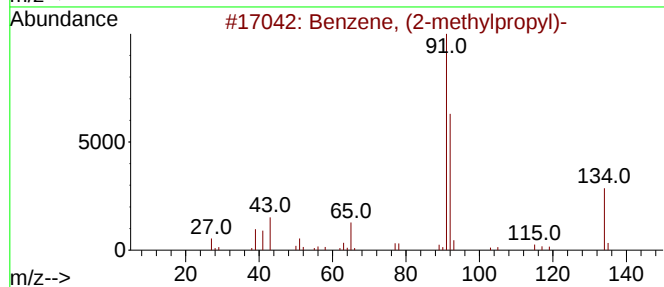
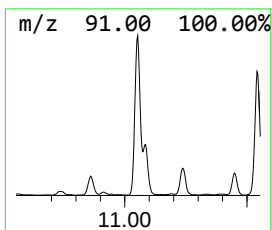
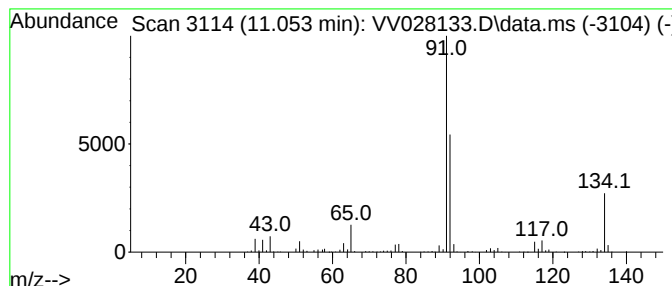
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 3 Benzene, (2-methylpropyl)- Concentration Rank 11

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 11.053 | 1.22 ug/L | 253389 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (2-methylpropyl)- | 134 | C10H14  | 000538-93-2 | 87   |
| 2    |    |   | Benzene, (2-methylpropyl)- | 134 | C10H14  | 000538-93-2 | 87   |
| 3    |    |   | Benzene, (2-methylpropyl)- | 134 | C10H14  | 000538-93-2 | 87   |
| 4    |    |   | Benzene, (2-methylpropyl)- | 134 | C10H14  | 000538-93-2 | 87   |
| 5    |    |   | Benzene, n-butyl-          | 134 | C10H14  | 000104-51-8 | 80   |



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ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :  
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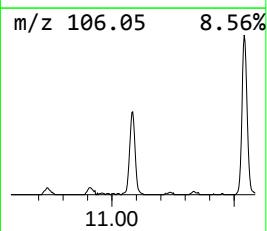
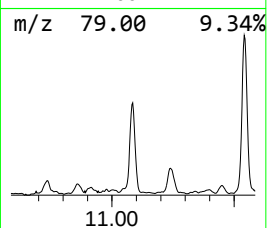
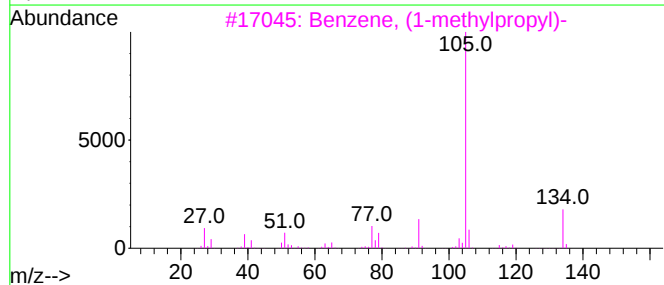
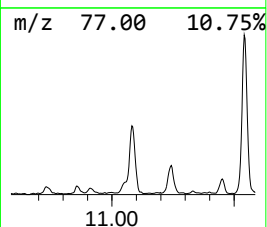
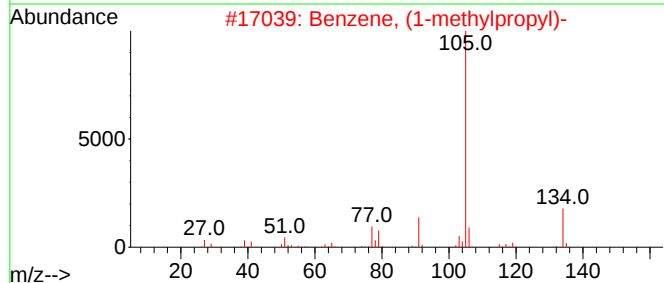
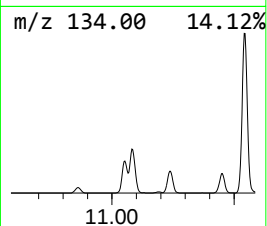
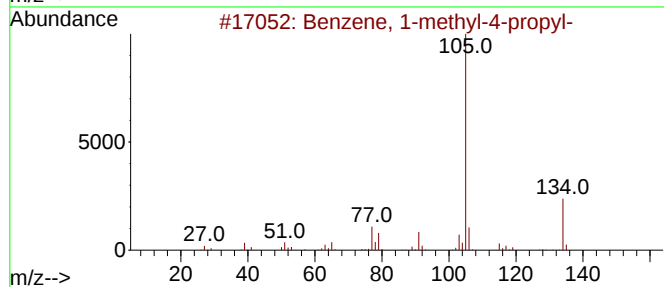
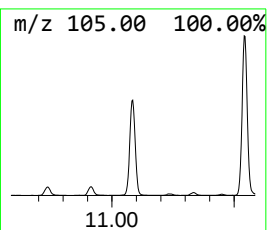
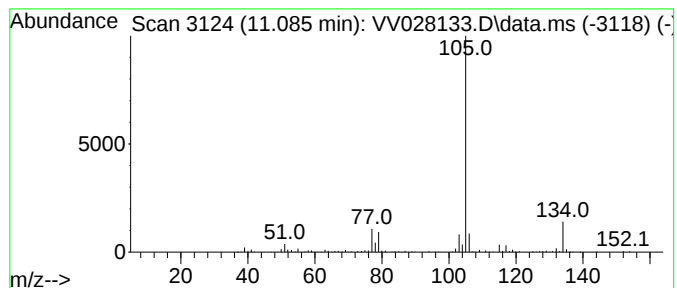
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 4 Benzene, 1-methyl-4-propyl- Concentration Rank 8

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 11.085 | 2.02 ug/L | 418829 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-4-propyl-         | 134 | C10H14  | 001074-55-1 | 91   |
| 2    |    |   | Benzene, (1-methylpropyl)-          | 134 | C10H14  | 000135-98-8 | 86   |
| 3    |    |   | Benzene, (1-methylpropyl)-          | 134 | C10H14  | 000135-98-8 | 78   |
| 4    |    |   | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O  | 000093-53-8 | 78   |
| 5    |    |   | Benzene, 1-methyl-2-propyl-         | 134 | C10H14  | 001074-17-5 | 78   |



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Instrument :  
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CB8L0

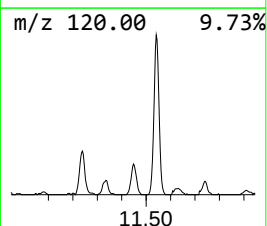
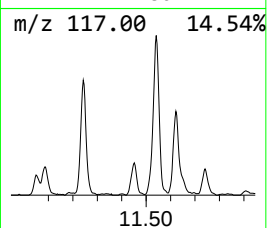
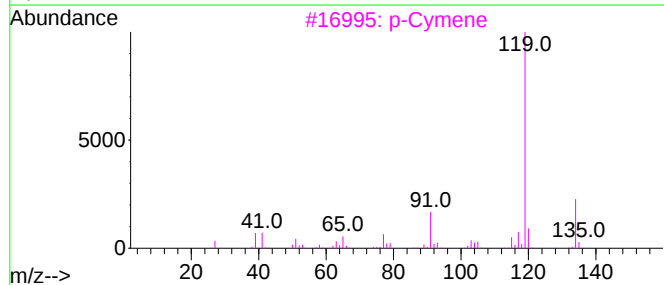
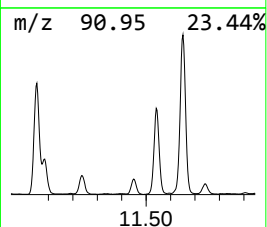
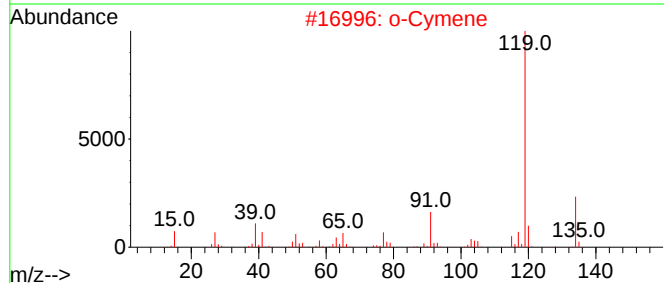
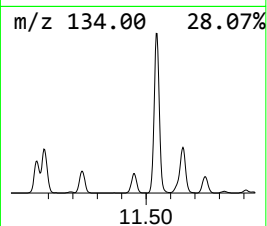
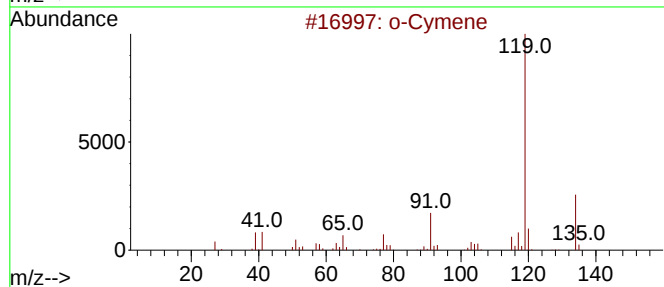
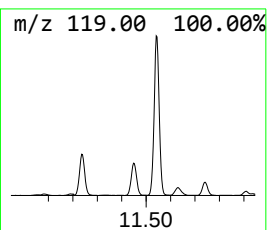
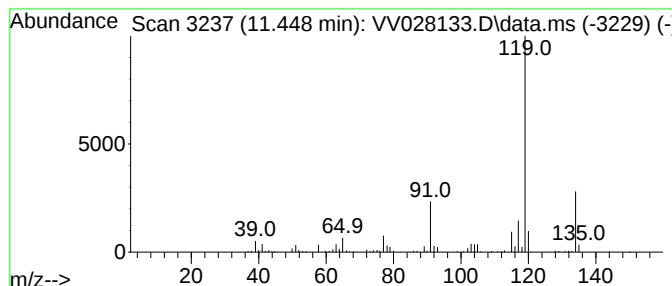
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 5 o-Cymene Concentration Rank 18

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 11.448 | 0.73 ug/L | 151759 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 97   |
| 2    |    |   | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 97   |
| 3    |    |   | p-Cymene                            | 134 | C10H14  | 000099-87-6 | 97   |
| 4    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 95   |
| 5    |    |   | p-Cymene                            | 134 | C10H14  | 000099-87-6 | 95   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

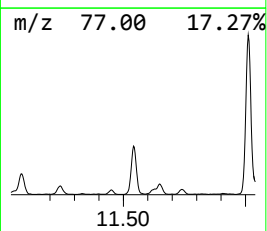
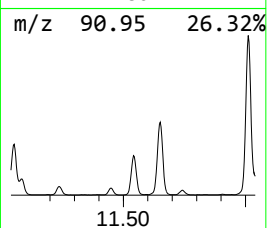
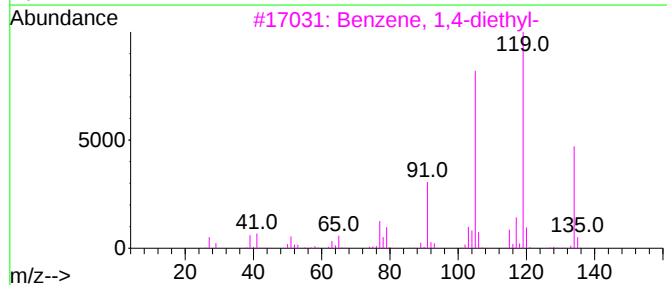
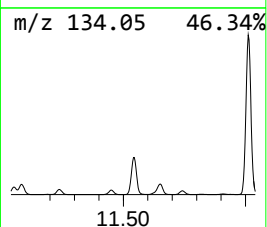
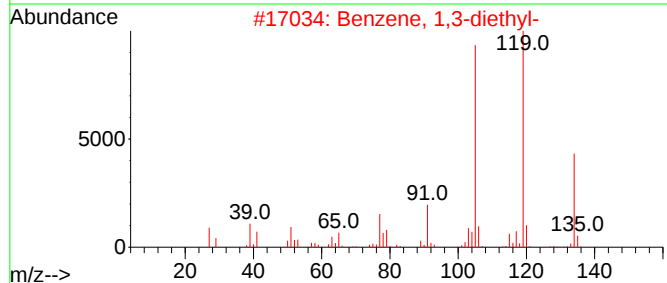
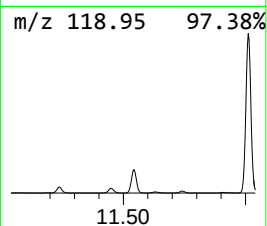
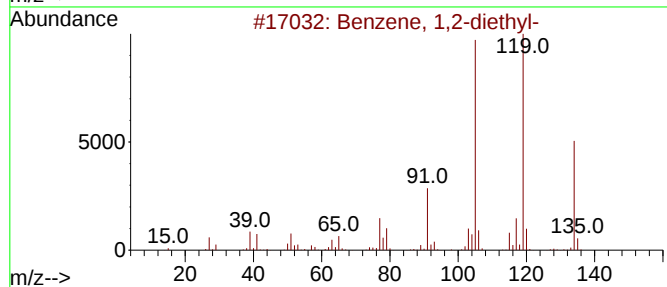
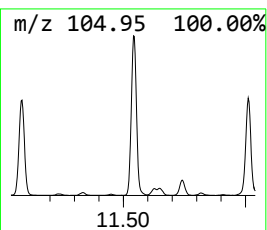
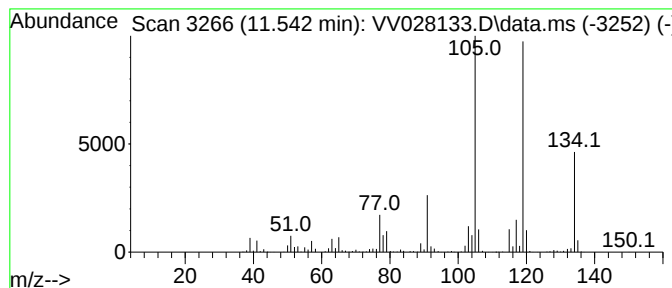
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 6 Benzene, 1,2-diethyl- Concentration Rank 4

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 11.541 | 6.39 ug/L | 1324020 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2-diethyl- | 134 | C10H14  | 000135-01-3 | 97   |
| 2    |    |   | Benzene, 1,3-diethyl- | 134 | C10H14  | 000141-93-5 | 97   |
| 3    |    |   | Benzene, 1,4-diethyl- | 134 | C10H14  | 000105-05-5 | 96   |
| 4    |    |   | Benzene, 1,3-diethyl- | 134 | C10H14  | 000141-93-5 | 95   |
| 5    |    |   | Benzene, 1,3-diethyl- | 134 | C10H14  | 000141-93-5 | 95   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

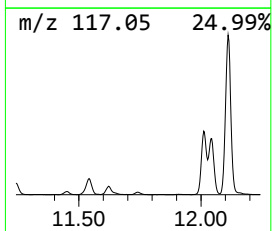
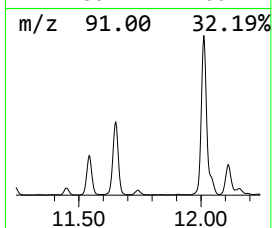
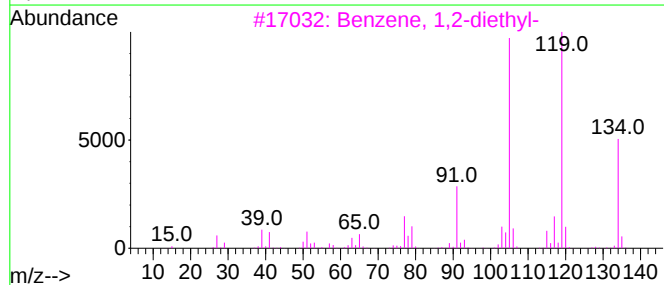
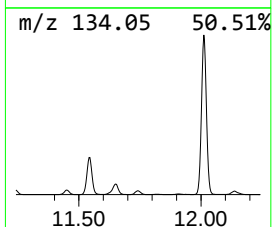
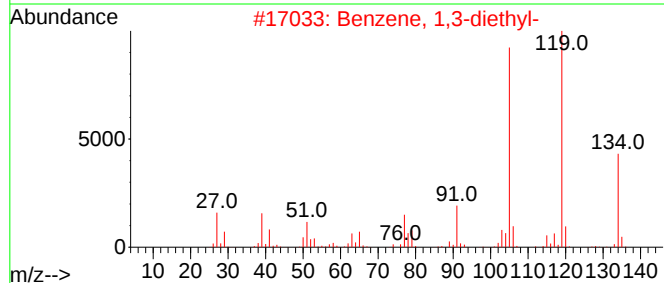
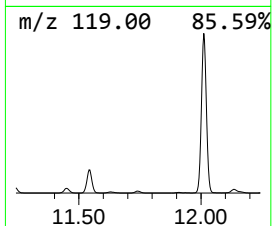
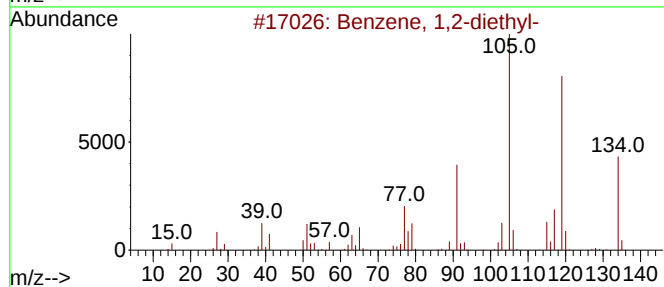
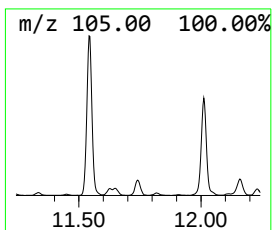
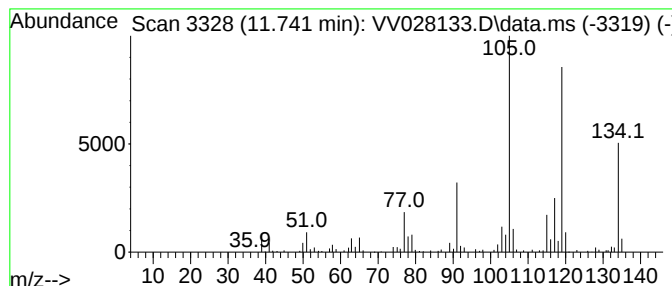
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 7 Benzene, 1,3-diethyl- Concentration Rank 22

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 11.741 | 0.62 ug/L | 127856 | 1,4-Dichlorobenzene-d4 | 11.246 |

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

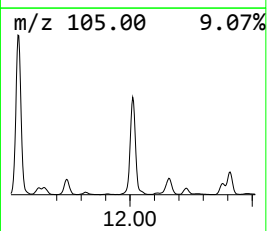
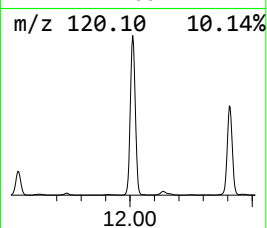
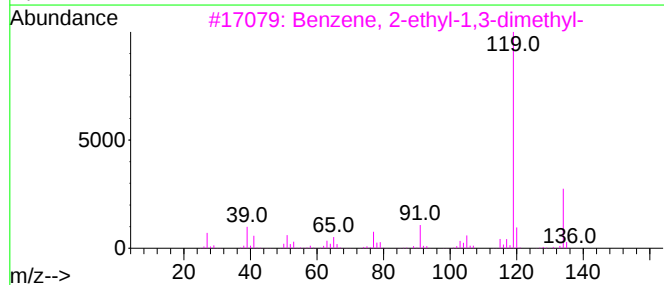
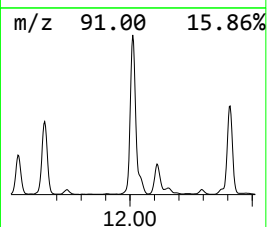
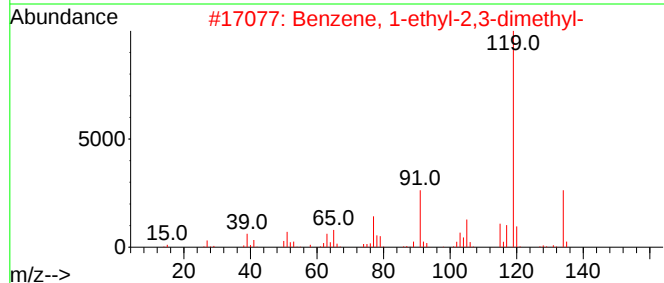
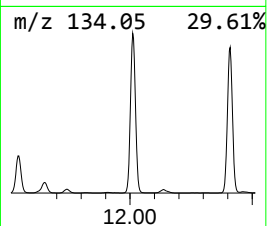
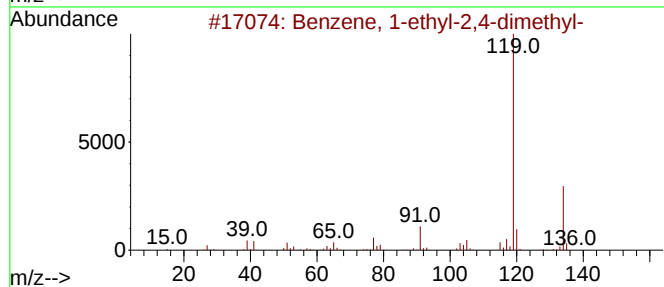
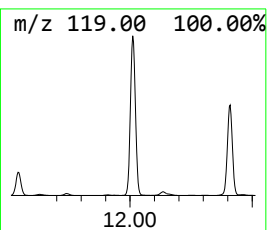
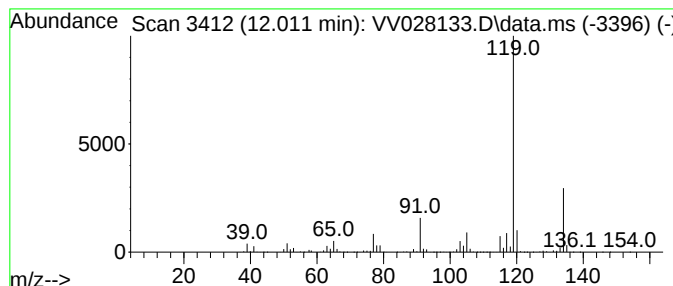
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 8 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 1

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 12.011 | 24.14 ug/L | 5004450 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 97   |
| 2    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 96   |
| 3    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 95   |
| 4    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 95   |
| 5    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 95   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

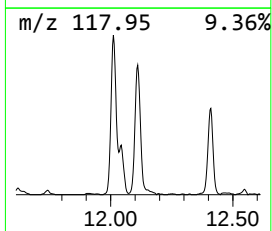
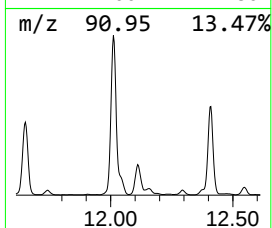
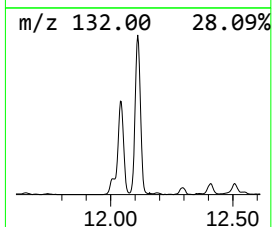
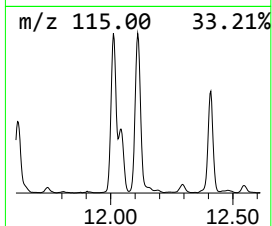
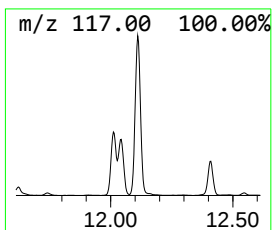
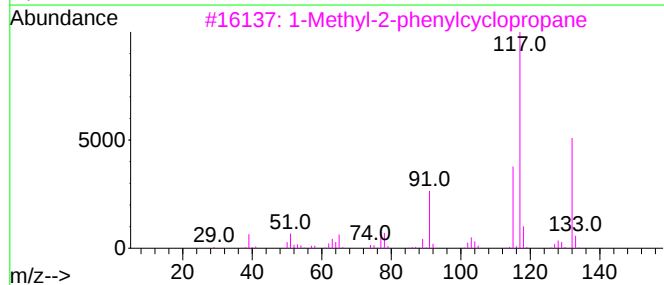
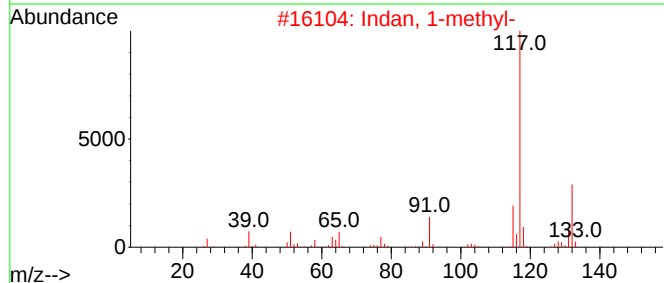
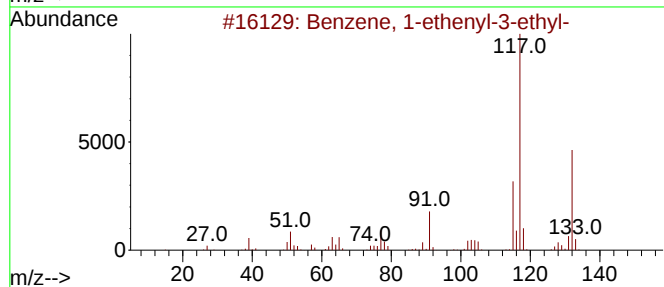
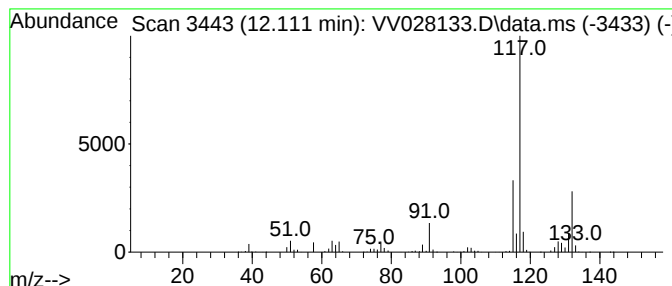
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 9 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 5

| R.T.   | EstConc   | Area    | Relative to ISTD       | R.T.   |
|--------|-----------|---------|------------------------|--------|
| 12.111 | 5.54 ug/L | 1147860 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethenyl-3-ethyl-   | 132 | C10H12  | 007525-62-4 | 91   |
| 2    |    |   | Indan, 1-methyl-              | 132 | C10H12  | 000767-58-8 | 90   |
| 3    |    |   | 1-Methyl-2-phenylcyclopropane | 132 | C10H12  | 003145-76-4 | 90   |
| 4    |    |   | 1-Phenyl-1-butene             | 132 | C10H12  | 000824-90-8 | 87   |
| 5    |    |   | Benzene, 1-ethenyl-4-ethyl-   | 132 | C10H12  | 003454-07-7 | 87   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

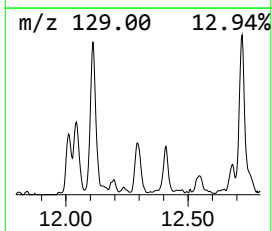
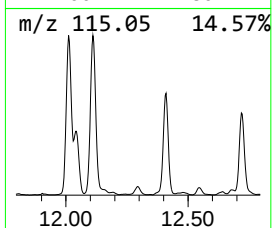
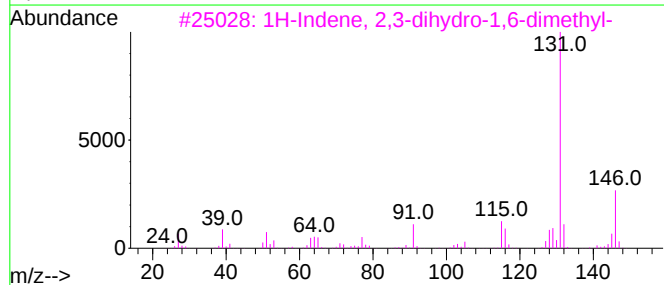
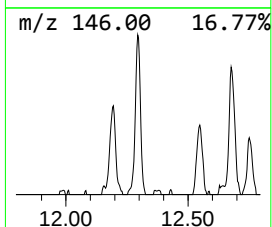
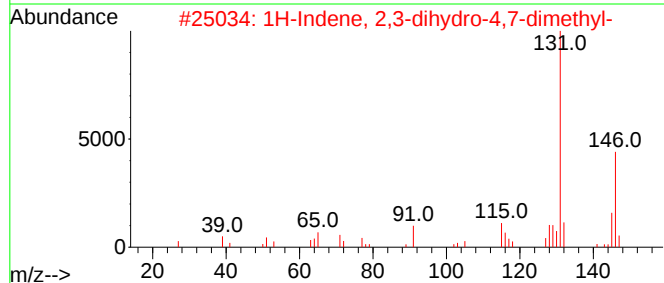
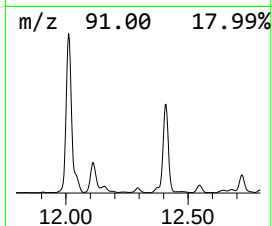
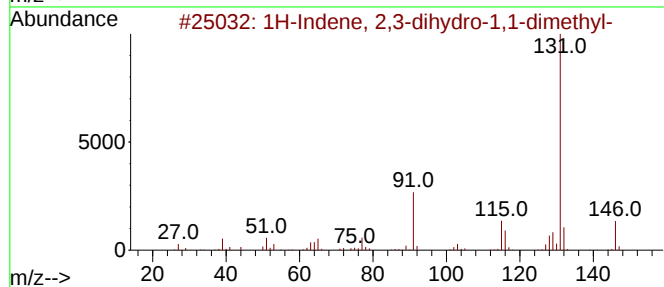
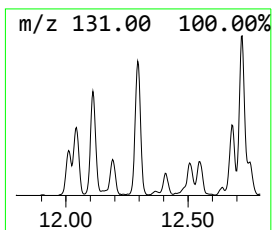
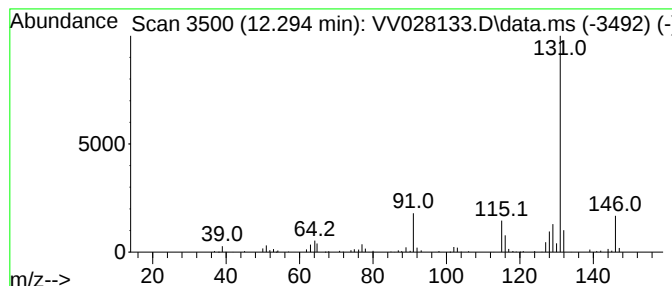
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 10 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 26

| R.T.      | EstConc                             | Area   | Relative to ISTD       |         |             | R.T.   |
|-----------|-------------------------------------|--------|------------------------|---------|-------------|--------|
| 12.294    | 0.50 ug/L                           | 104355 | 1,4-Dichlorobenzene-d4 |         |             | 11.246 |
| Hit# of 5 | Tentative ID                        |        | MW                     | MolForm | CAS#        | Qual   |
| 1         | 1H-Indene, 2,3-dihydro-1,1-dimet... |        | 146                    | C11H14  | 004912-92-9 | 93     |
| 2         | 1H-Indene, 2,3-dihydro-4,7-dimet... |        | 146                    | C11H14  | 006682-71-9 | 91     |
| 3         | 1H-Indene, 2,3-dihydro-1,6-dimet... |        | 146                    | C11H14  | 017059-48-2 | 91     |
| 4         | 1H-Indene, 2,3-dihydro-1,3-dimet... |        | 146                    | C11H14  | 004175-53-5 | 90     |
| 5         | 1H-Indene, 2,3-dihydro-1,1-dimet... |        | 146                    | C11H14  | 004912-92-9 | 90     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

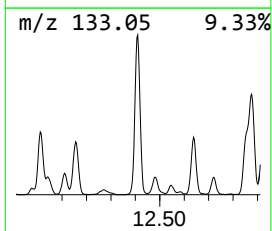
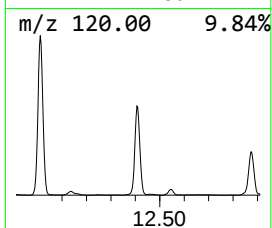
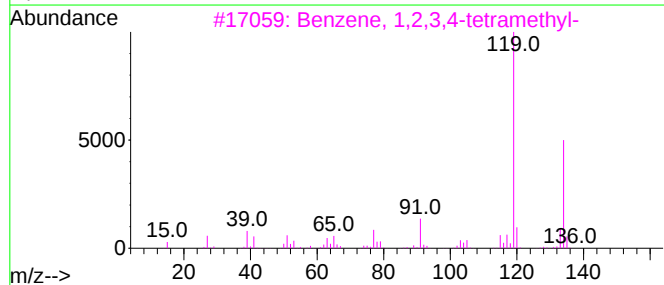
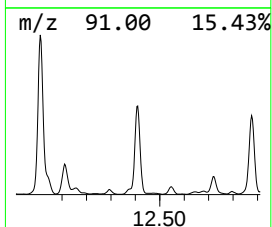
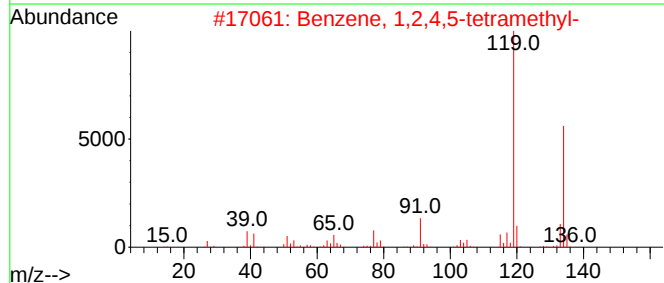
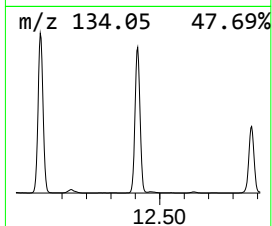
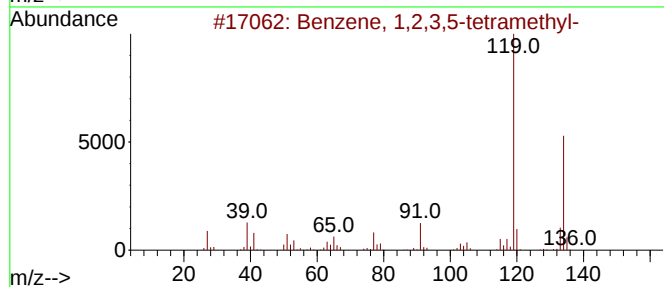
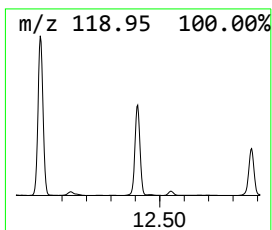
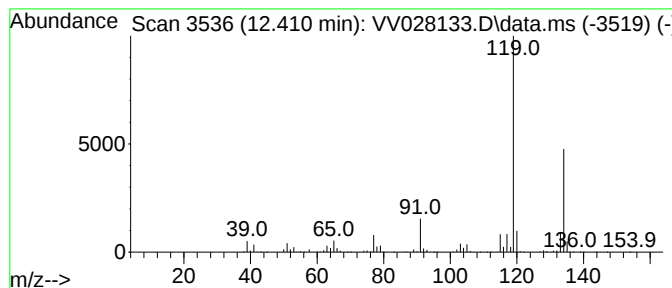
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 3

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 12.410 | 13.48 ug/L | 2793890 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 97   |
| 2    |    |   | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 97   |
| 3    |    |   | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 97   |
| 4    |    |   | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 97   |
| 5    |    |   | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 96   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

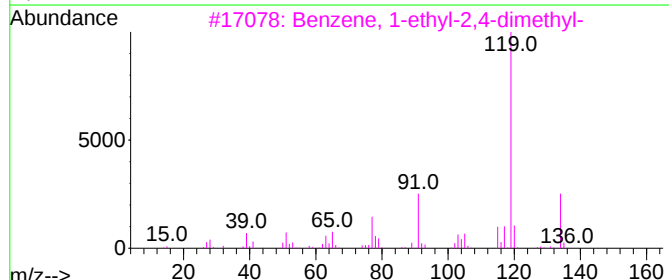
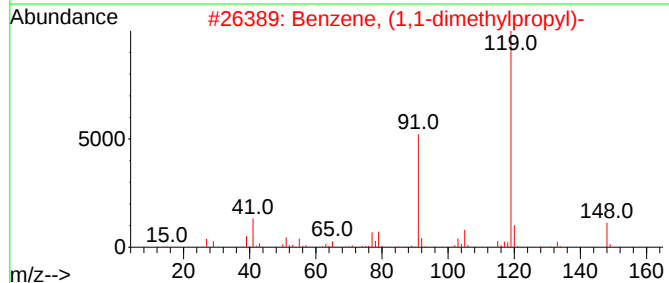
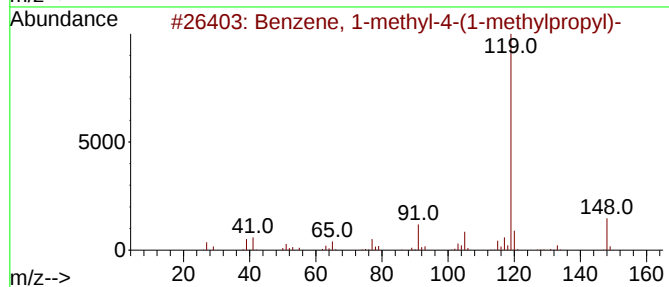
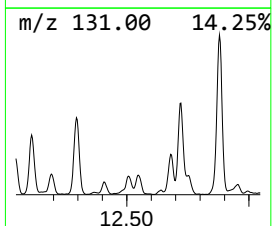
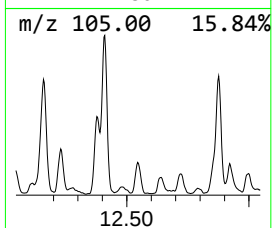
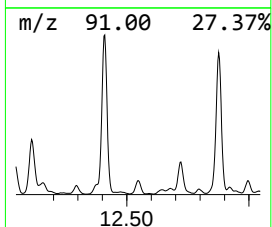
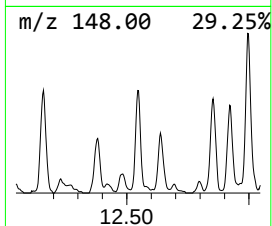
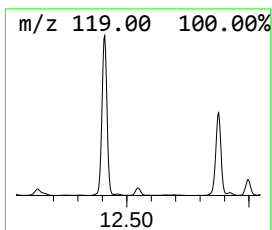
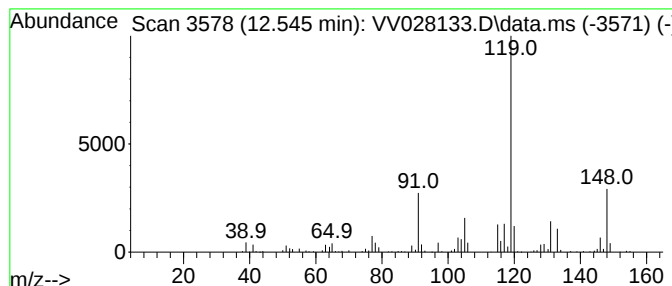
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Benzene, 1-methyl-4-(1-meth... Concentration Rank 16

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.545 | 0.85 ug/L | 176327 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16  | 001595-16-0 | 76   |
| 2    |    |   | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8 | 64   |
| 3    |    |   | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9 | 60   |
| 4    |    |   | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 60   |
| 5    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 60   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

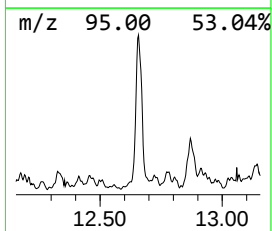
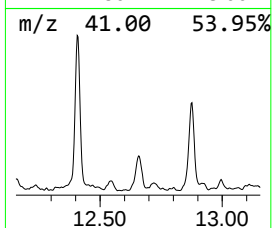
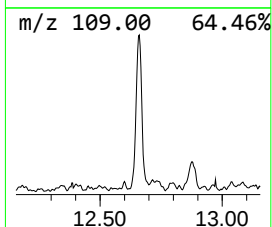
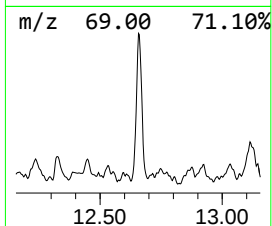
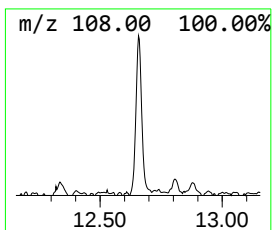
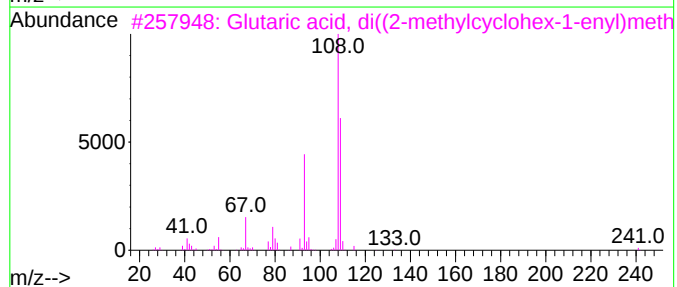
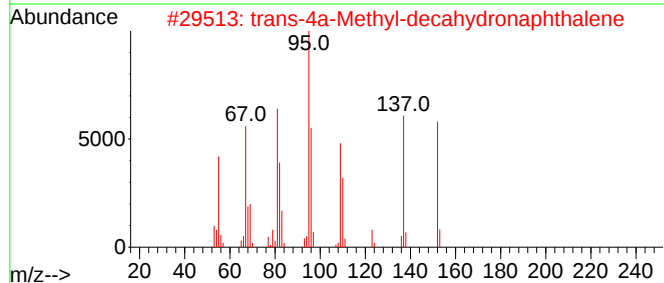
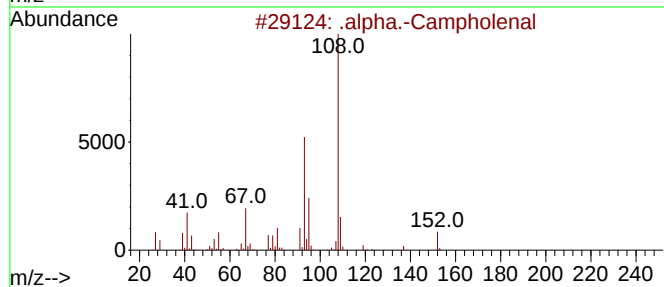
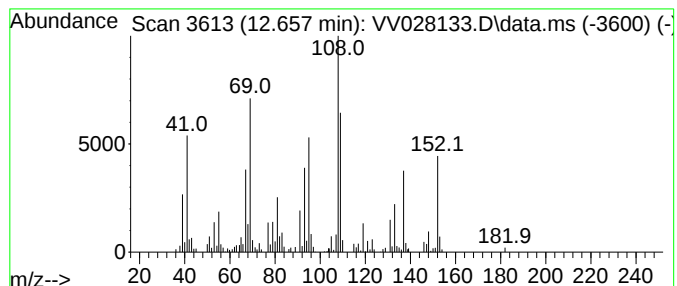
Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 13 .alpha.-Campholenal Concentration Rank 10

| R.T.      | EstConc                             | Area   | Relative to ISTD       |          |              | R.T.   |
|-----------|-------------------------------------|--------|------------------------|----------|--------------|--------|
| 12.657    | 1.33 ug/L                           | 276544 | 1,4-Dichlorobenzene-d4 |          |              | 11.246 |
| Hit# of 5 | Tentative ID                        |        | MW                     | MolForm  | CAS#         | Qual   |
| 1         | .alpha.-Campholenal                 |        | 152                    | C10H16O  | 004501-58-0  | 50     |
| 2         | trans-4a-Methyl-decahydronaphtha... |        | 152                    | C11H20   | 002547-27-5  | 41     |
| 3         | Glutaric acid, di((2-methylcyclo... |        | 348                    | C21H32O4 | 1000405-51-2 | 38     |
| 4         | Ethanone, 1-(1,4-dimethyl-3-cycl... |        | 152                    | C10H16O  | 043219-68-7  | 38     |
| 5         | Ethanol, 2-(4-methylphenoxy)-       |        | 152                    | C9H12O2  | 015149-10-7  | 30     |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

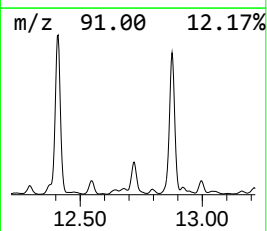
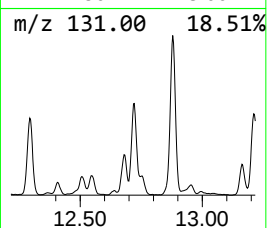
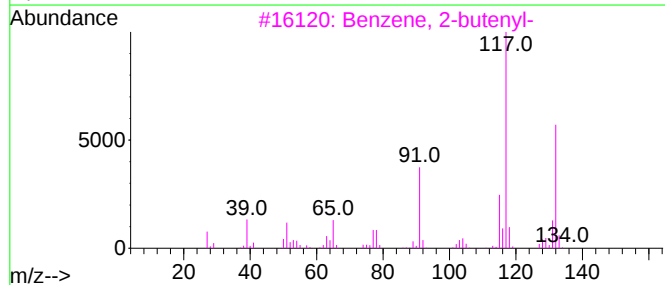
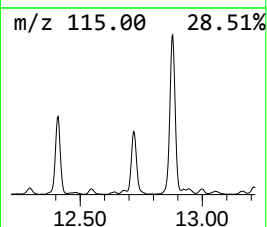
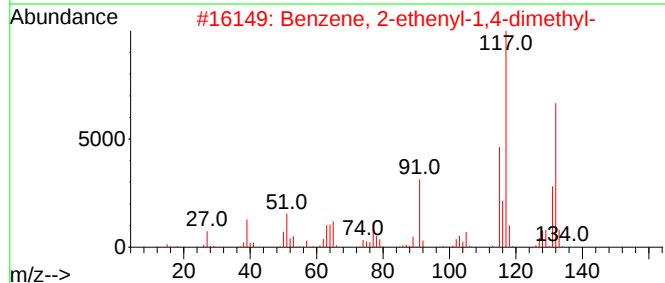
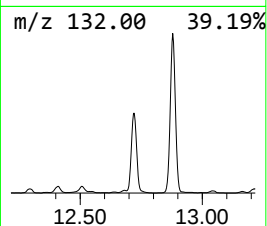
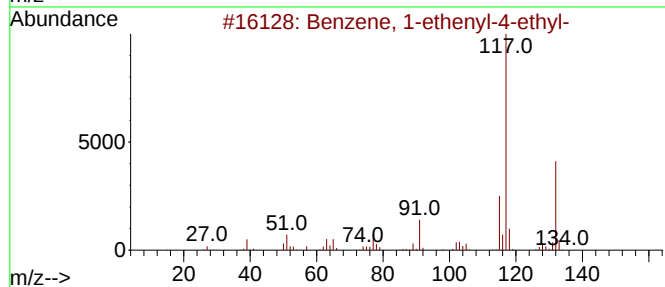
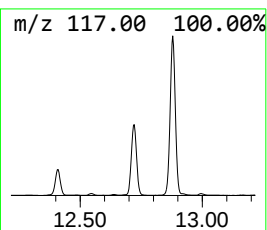
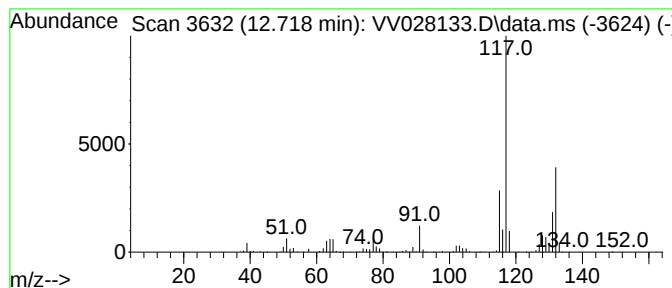
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 14 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 6

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.718 | 3.84 ug/L | 795112 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethenyl-4-ethyl-      | 132 | C10H12  | 003454-07-7 | 94   |
| 2    |    |   | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12  | 002039-89-6 | 92   |
| 3    |    |   | Benzene, 2-butenyl-              | 132 | C10H12  | 001560-06-1 | 91   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-4-methyl- | 132 | C10H12  | 000824-22-6 | 90   |
| 5    |    |   | Benzene, 1-ethenyl-3-ethyl-      | 132 | C10H12  | 007525-62-4 | 90   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

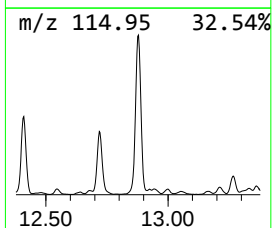
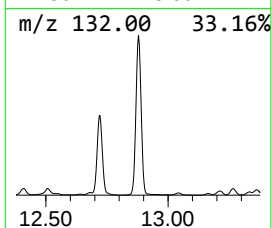
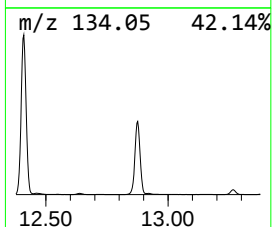
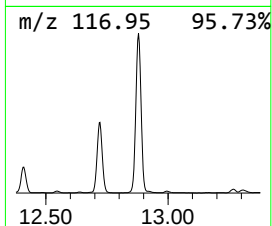
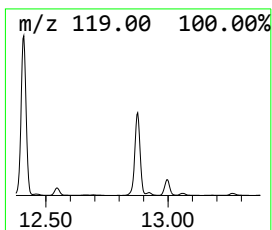
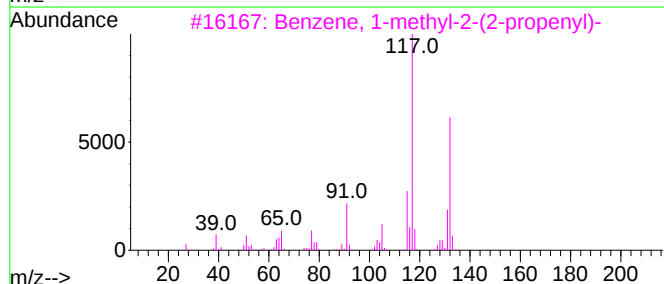
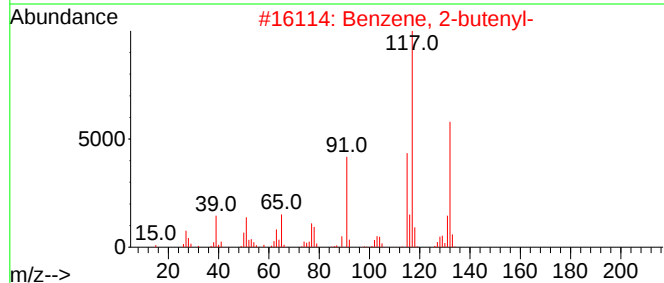
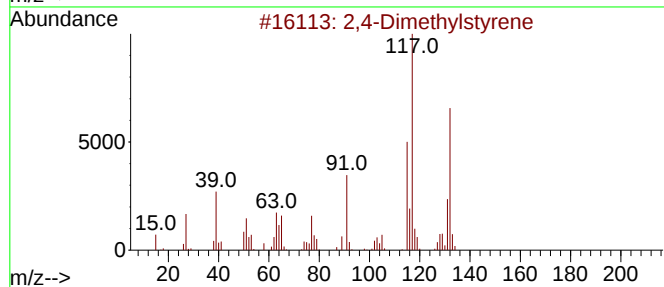
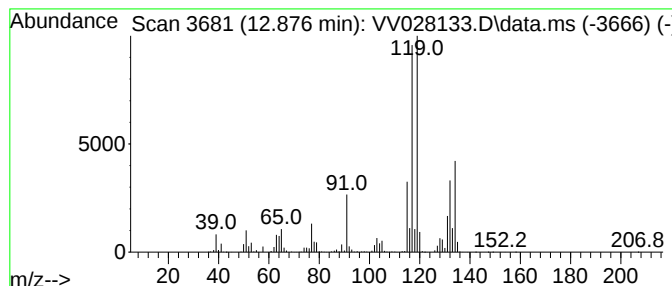
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 15 2,4-Dimethylstyrene Concentration Rank 2

| R.T.   | EstConc    | Area    | Relative to ISTD       | R.T.   |
|--------|------------|---------|------------------------|--------|
| 12.876 | 14.50 ug/L | 3005930 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2,4-Dimethylstyrene               | 132 | C10H12  | 002234-20-0 | 86   |
| 2    |    |   | Benzene, 2-butenyl-               | 132 | C10H12  | 001560-06-1 | 64   |
| 3    |    |   | Benzene, 1-methyl-2-(2-propenyl)- | 132 | C10H12  | 001587-04-8 | 50   |
| 4    |    |   | Benzene, 1,2,3,5-tetramethyl-     | 134 | C10H14  | 000527-53-7 | 50   |
| 5    |    |   | 1-Methyl-2-phenylcyclopropane     | 132 | C10H12  | 003145-76-4 | 50   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

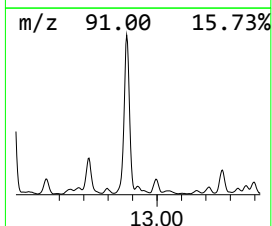
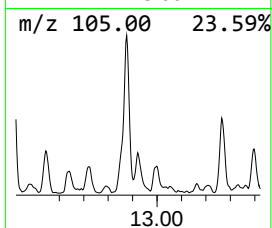
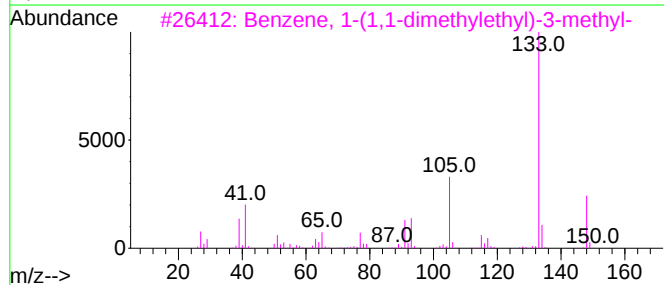
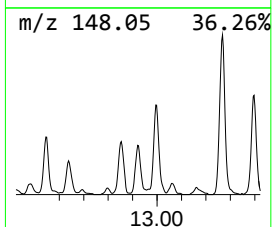
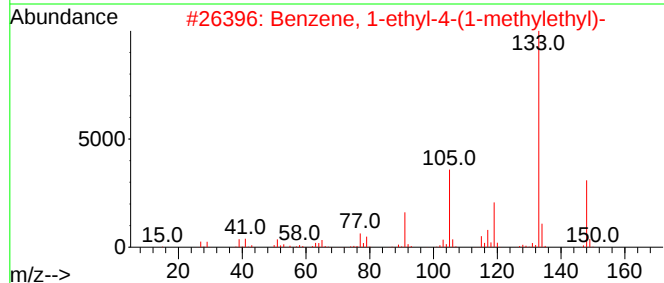
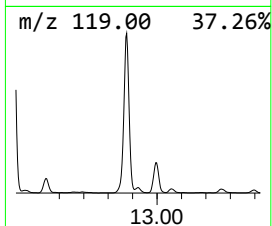
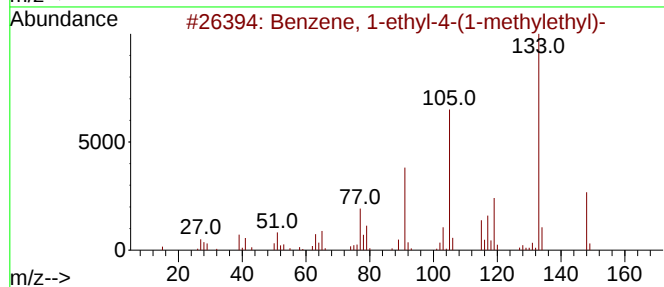
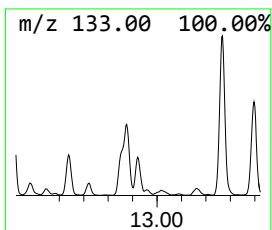
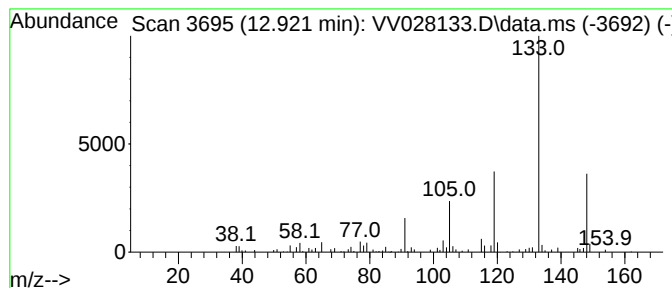
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 21

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.921 | 0.62 ug/L | 129046 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 | C11H16  | 004218-48-8  | 83   |
| 2    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 | C11H16  | 004218-48-8  | 72   |
| 3    |    |   | Benzene, 1-(1,1-dimethylethyl)-3... | 148 | C11H16  | 001075-38-3  | 64   |
| 4    |    |   | Benzene, 1,3-dimethyl-5-(1-methy... | 148 | C11H16  | 004706-90-5  | 64   |
| 5    |    |   | 3,4-Dimethylcumene                  | 148 | C11H16  | 1000370-34-1 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

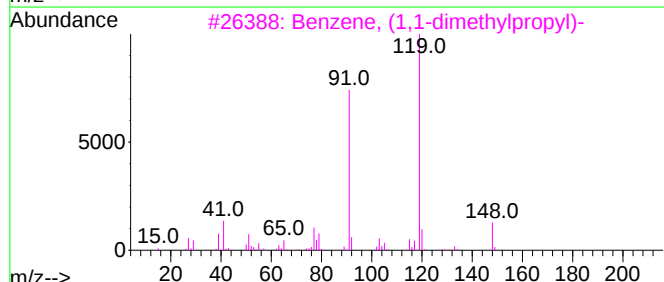
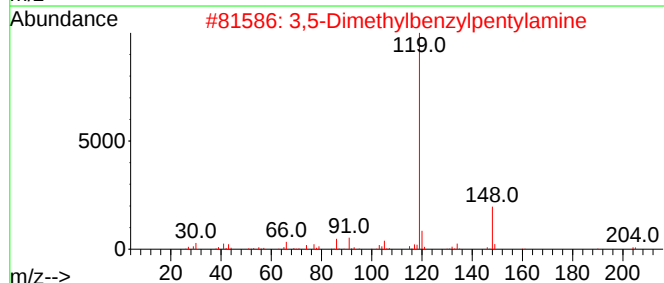
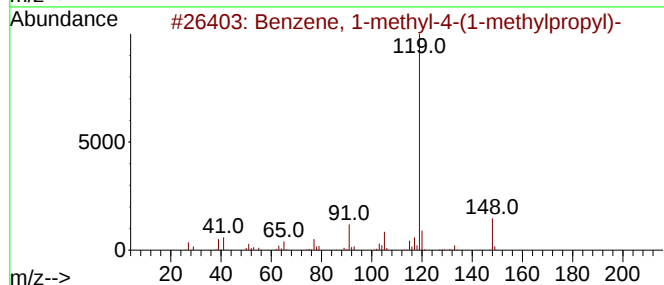
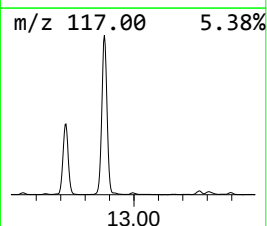
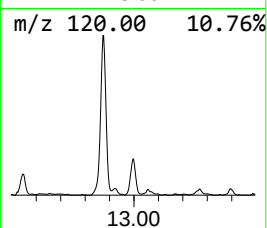
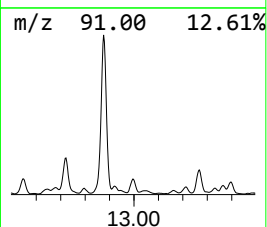
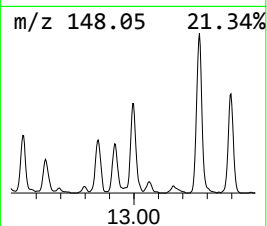
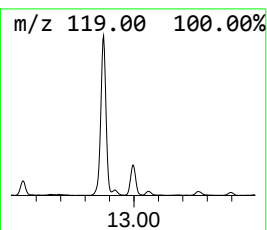
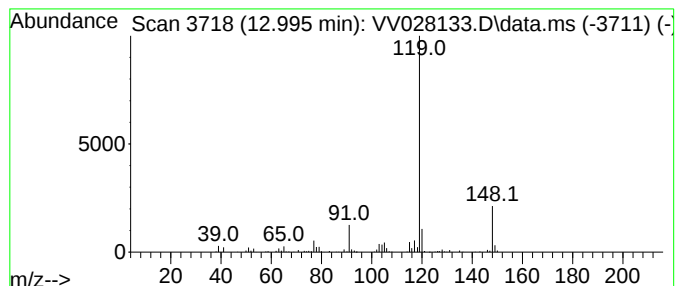
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 3,5-Dimethylbenzylpentylamine Concentration Rank 15

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.995 | 0.89 ug/L | 184878 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16  | 001595-16-0  | 91   |
| 2    |    |   | 3,5-Dimethylbenzylpentylamine       | 205 | C14H23N | 1000491-60-8 | 78   |
| 3    |    |   | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8  | 78   |
| 4    |    |   | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9  | 72   |
| 5    |    |   | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5  | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

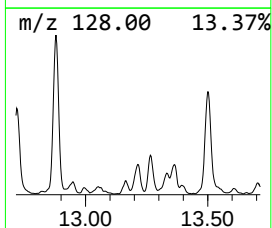
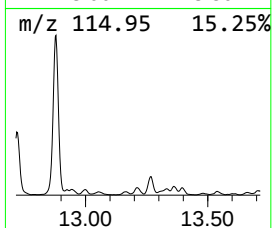
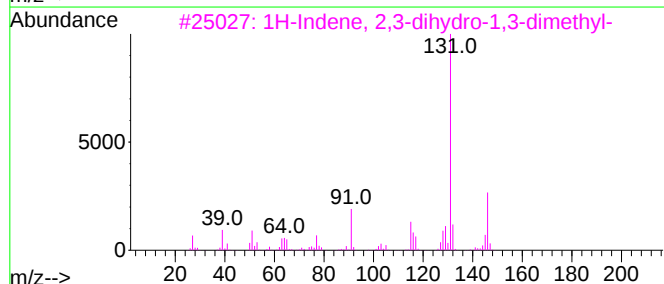
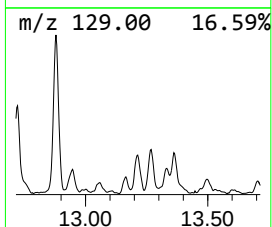
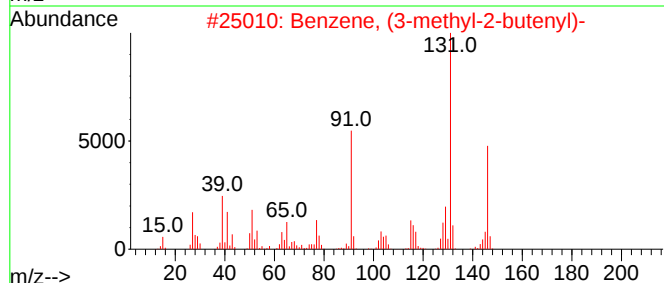
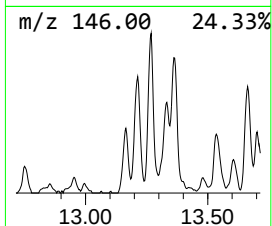
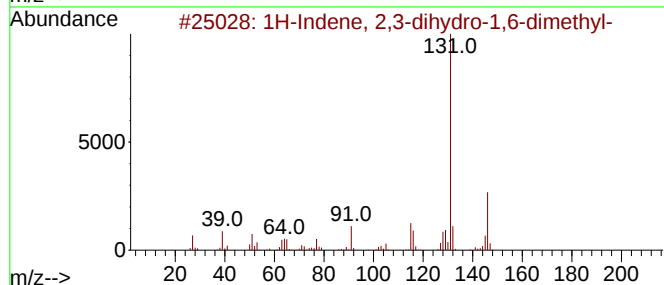
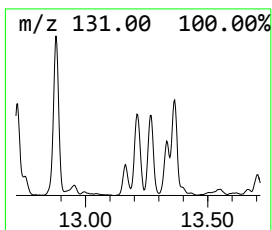
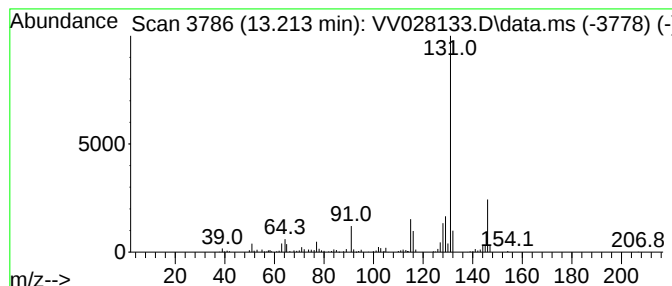
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 24

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.214 | 0.56 ug/L | 115357 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14  | 017059-48-2 | 94   |
| 2    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 94   |
| 3    |    |   | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146 | C11H14  | 004175-53-5 | 91   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 91   |
| 5    |    |   | 2,2-Dimethylindene, 2,3-dihydro-    | 146 | C11H14  | 020836-11-7 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

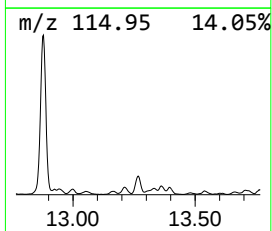
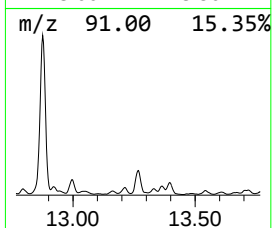
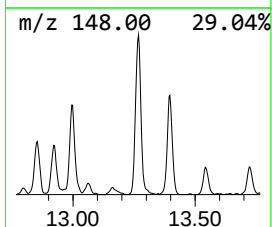
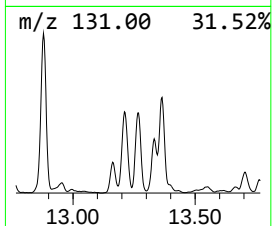
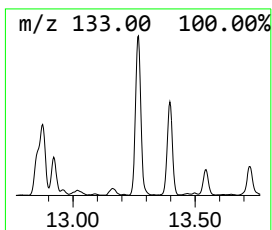
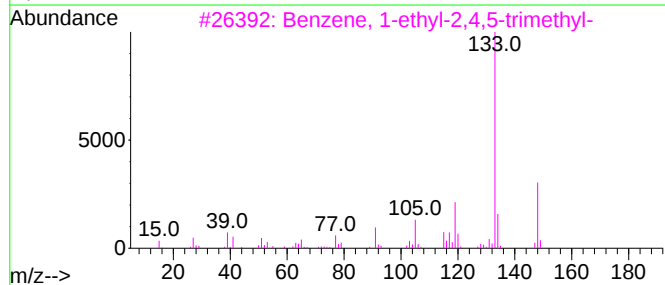
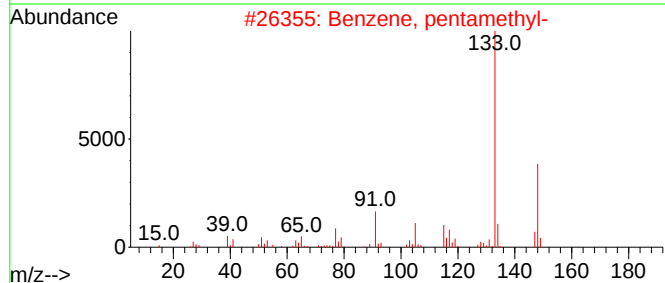
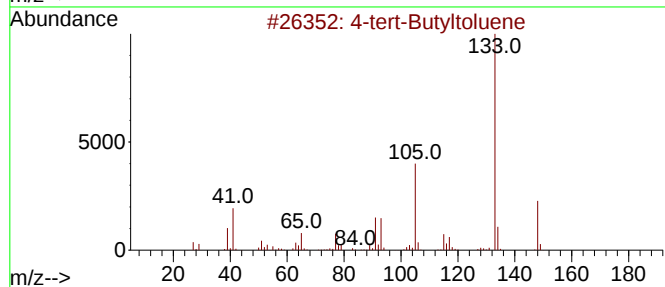
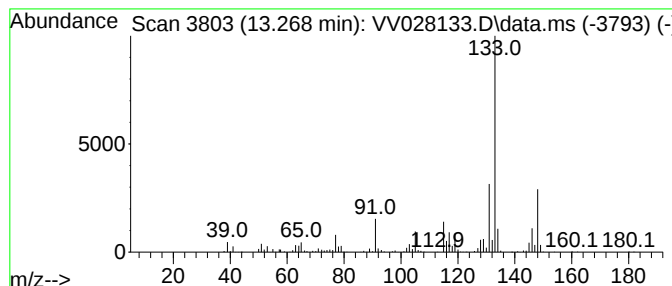
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 19 4-tert-Butyltoluene Concentration Rank 7

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.268 | 2.34 ug/L | 485845 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | 4-tert-Butyltoluene                 | 148 | C11H16  | 000098-51-1  | 70   |
| 2    |    |   | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9  | 70   |
| 3    |    |   | Benzene, 1-ethyl-2,4,5-trimethyl-   | 148 | C11H16  | 017851-27-3  | 68   |
| 4    |    |   | 3,4-Dimethylcumene                  | 148 | C11H16  | 1000370-34-1 | 68   |
| 5    |    |   | Benzene, 1-(1,1-dimethylethyl)-3... | 148 | C11H16  | 001075-38-3  | 62   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

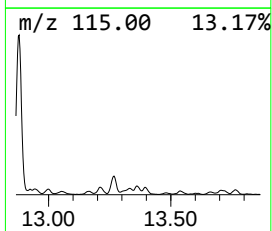
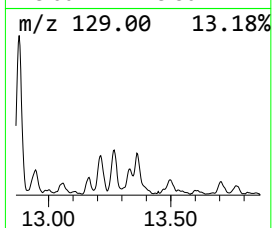
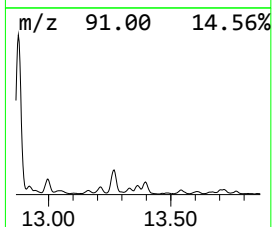
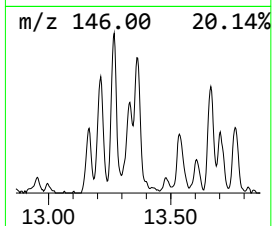
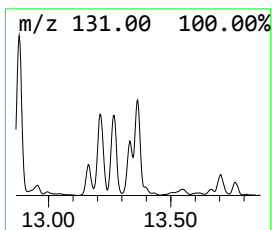
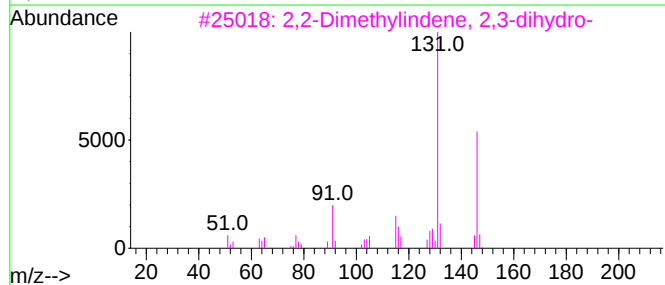
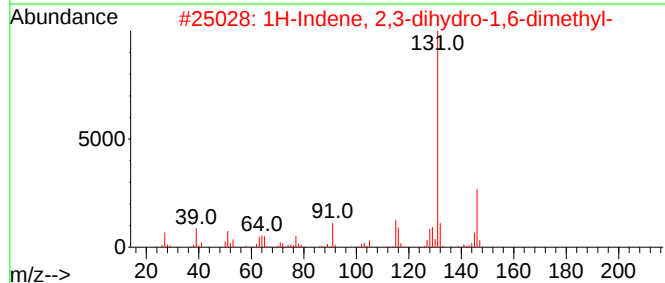
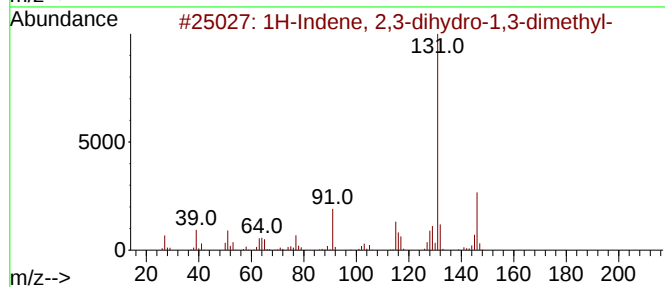
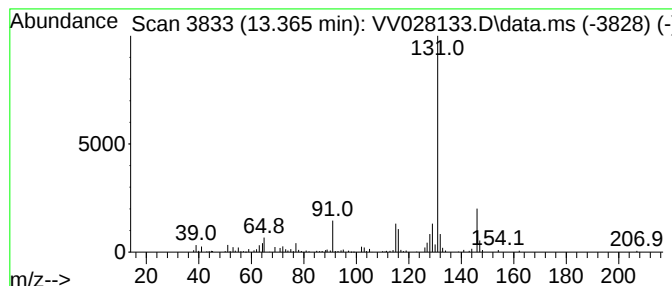
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 20

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.365 | 0.64 ug/L | 132397 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of 5                                | Tentative ID | MW     | MolForm | CAS#        | Qual |
|------|-------------------------------------|--------------|--------|---------|-------------|------|
| 1    | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146          | C11H14 |         | 004175-53-5 | 91   |
| 2    | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146          | C11H14 |         | 017059-48-2 | 91   |
| 3    | 2,2-Dimethylindene, 2,3-dihydro-    | 146          | C11H14 |         | 020836-11-7 | 91   |
| 4    | Benzene, (1,2-dimethyl-1-propenyl)- | 146          | C11H14 |         | 000769-57-3 | 91   |
| 5    | 1H-Indene, 2,3-dihydro-1,1-dimet... | 146          | C11H14 |         | 004912-92-9 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

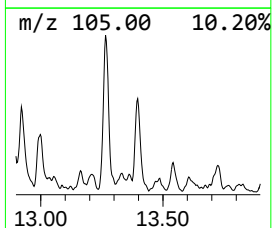
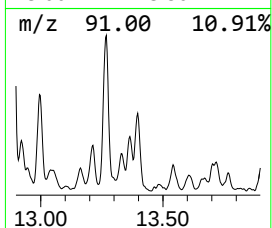
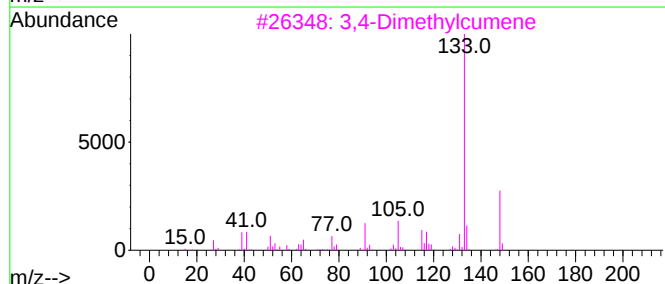
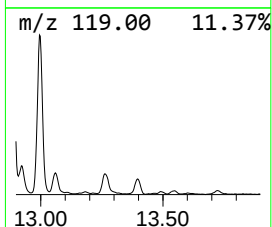
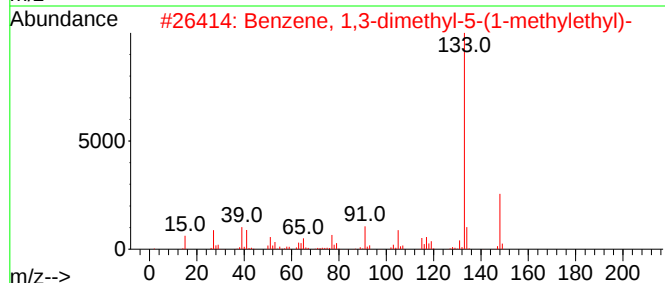
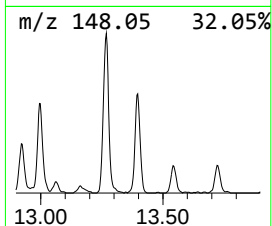
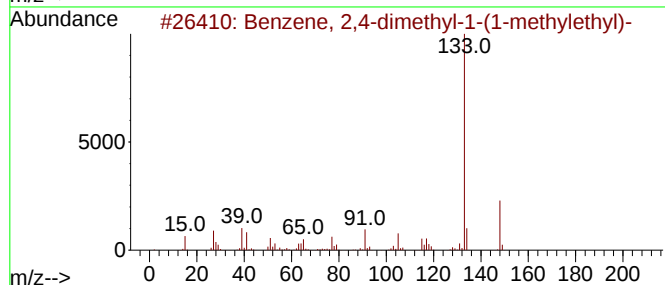
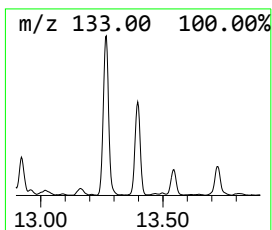
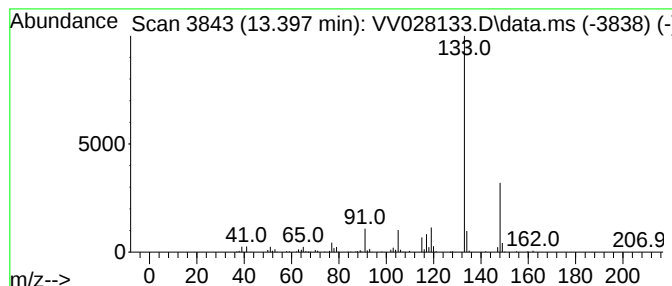
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 Benzene, 2,4-dimethyl-1-(1-... Concentration Rank 12

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.397 | 1.06 ug/L | 219850 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Benzene, 2,4-dimethyl-1-(1-methy... | 148 | C11H16  | 004706-89-2  | 87   |
| 2    |    |   | Benzene, 1,3-dimethyl-5-(1-methy... | 148 | C11H16  | 004706-90-5  | 87   |
| 3    |    |   | 3,4-Dimethylcumene                  | 148 | C11H16  | 1000370-34-1 | 86   |
| 4    |    |   | Benzene, pentamethyl-               | 148 | C11H16  | 000700-12-9  | 80   |
| 5    |    |   | Benzene, 1-ethyl-3-(1-methylethyl)- | 148 | C11H16  | 004920-99-4  | 78   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

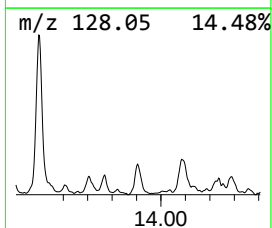
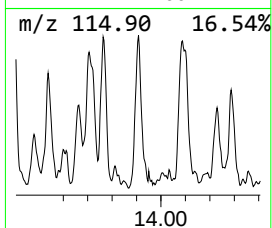
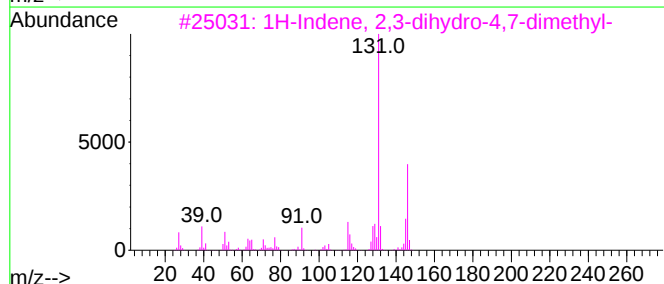
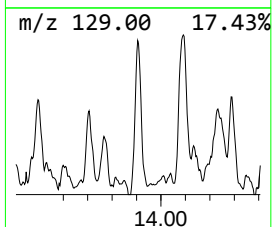
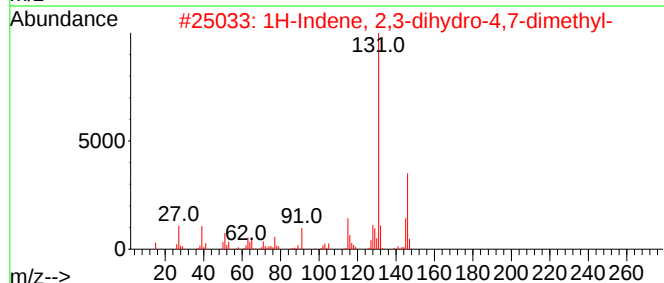
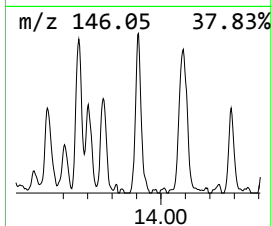
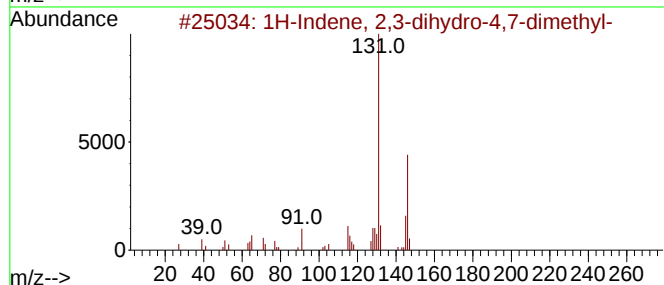
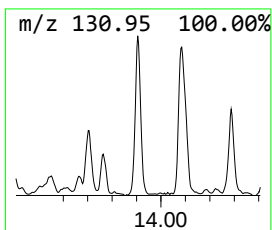
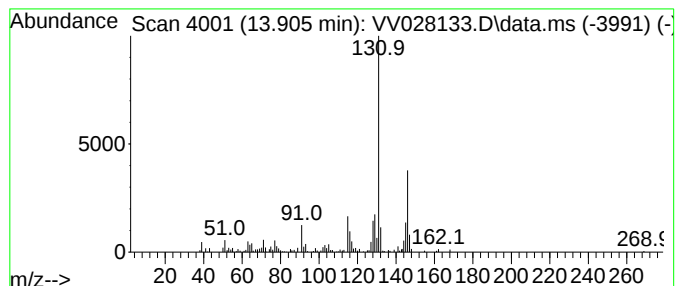
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 22 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 19

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.905 | 0.67 ug/L | 138189 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of 5                                | Tentative ID | MW     | MolForm | CAS#        | Qual |
|------|-------------------------------------|--------------|--------|---------|-------------|------|
| 1    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146          | C11H14 |         | 006682-71-9 | 97   |
| 2    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146          | C11H14 |         | 006682-71-9 | 97   |
| 3    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146          | C11H14 |         | 006682-71-9 | 94   |
| 4    | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146          | C11H14 |         | 017057-82-8 | 94   |
| 5    | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146          | C11H14 |         | 004175-53-5 | 94   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

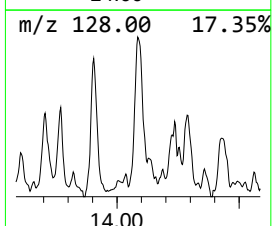
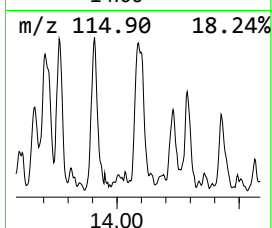
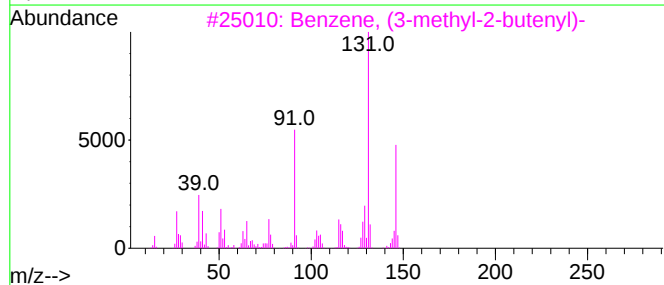
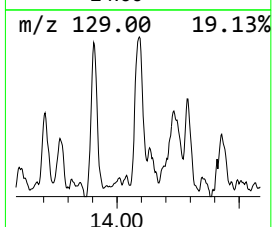
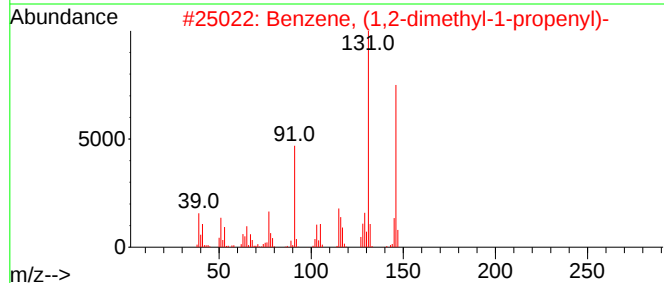
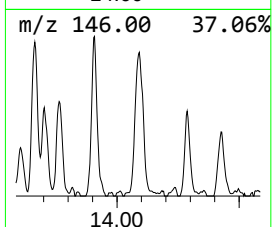
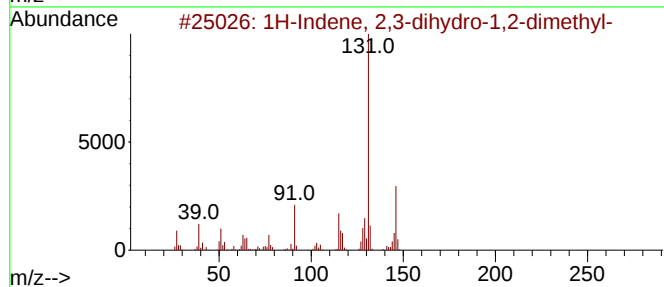
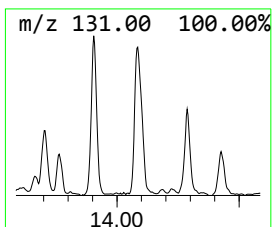
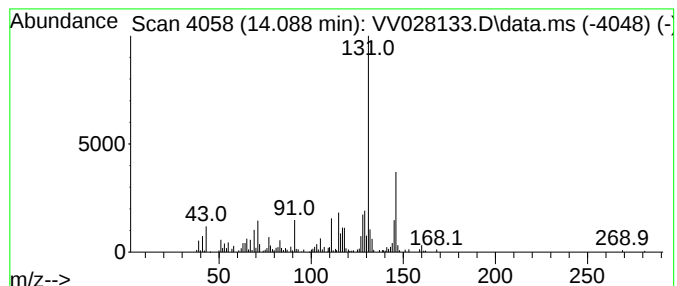
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 23 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 14

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 14.088 | 0.93 ug/L | 192760 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 93   |
| 2    |    |   | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 90   |
| 3    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 83   |
| 4    |    |   | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 83   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146 | C11H14  | 004175-53-5 | 81   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

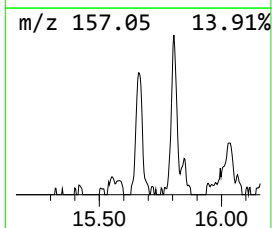
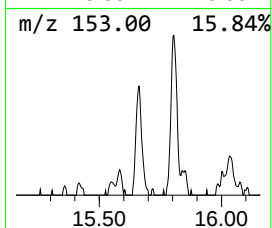
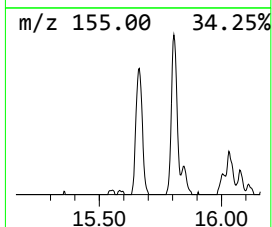
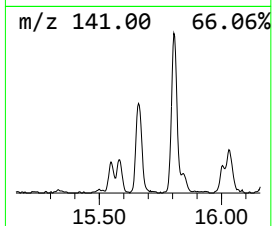
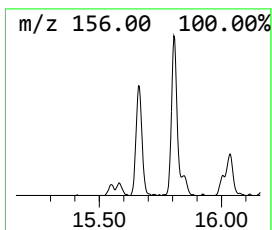
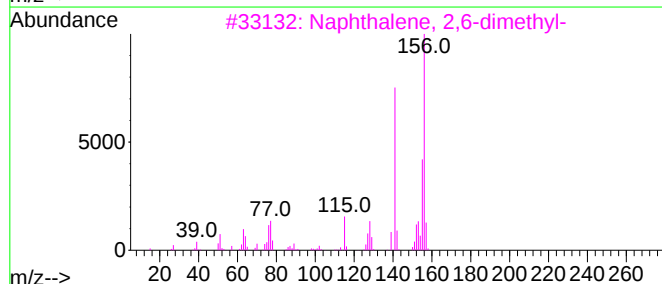
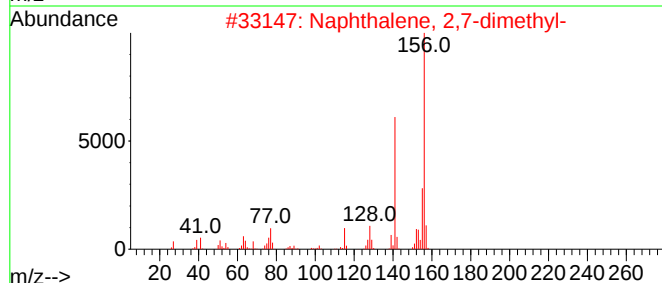
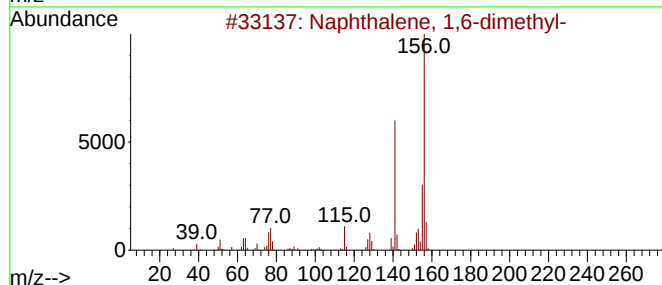
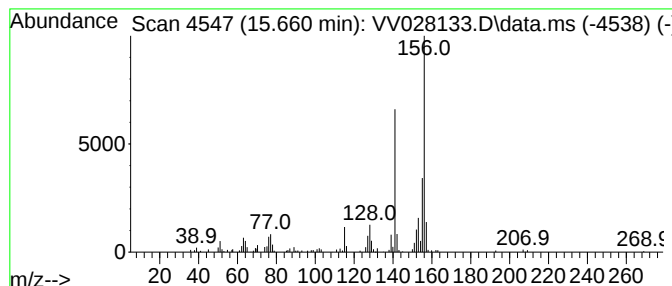
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 24 Naphthalene, 1,6-dimethyl- Concentration Rank 25

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 15.660 | 0.55 ug/L | 114120 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |
| 2    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 3    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 97   |
| 4    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 97   |
| 5    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

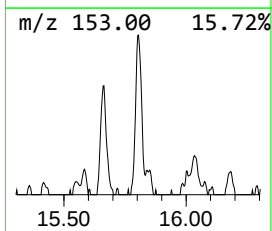
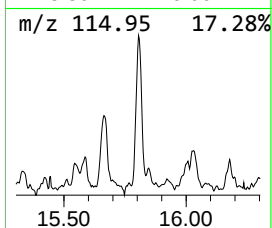
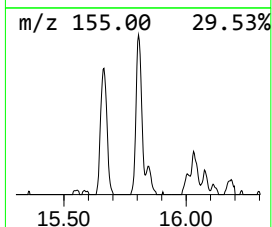
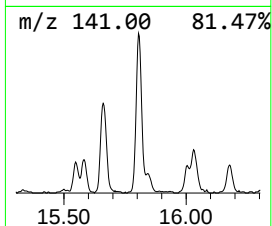
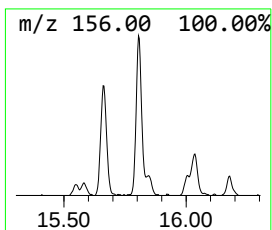
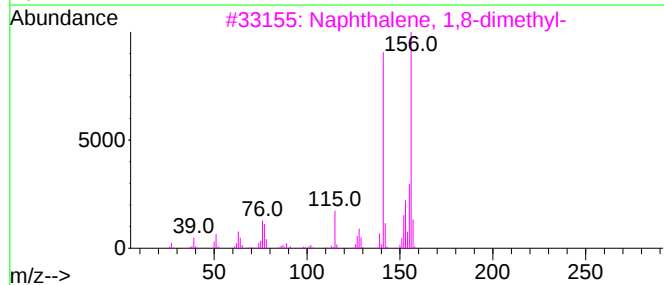
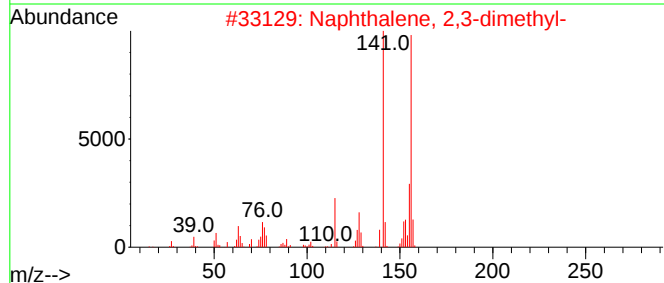
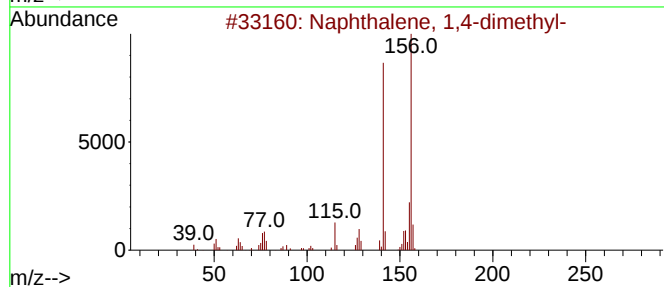
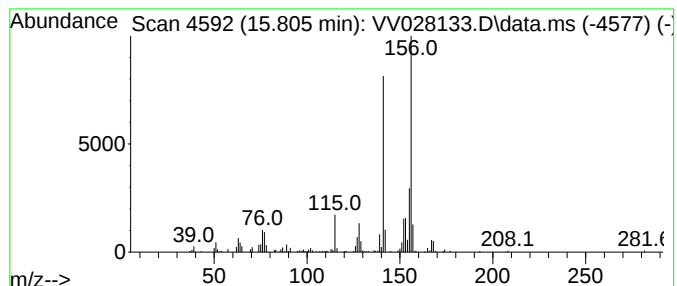
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 25 Naphthalene, 1,4-dimethyl- Concentration Rank 17

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 15.805 | 0.77 ug/L | 160552 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 97   |
| 2    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 3    |    |   | Naphthalene, 1,8-dimethyl- | 156 | C12H12  | 000569-41-5 | 97   |
| 4    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |
| 5    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV092722\  
Data File : VV028133.D  
Acq On : 27 Sep 2022 12:40  
Operator : SY/MD  
Sample : N4820-02  
Misc : 25.0mL/MSVOA\_V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
MSVOA\_V  
ClientSampleId :  
CB8L0

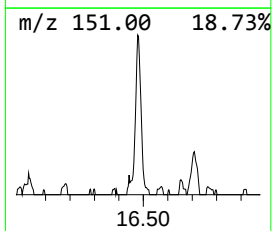
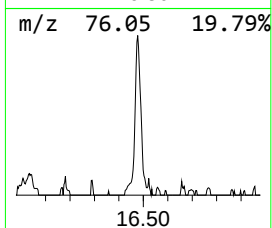
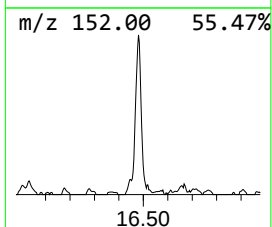
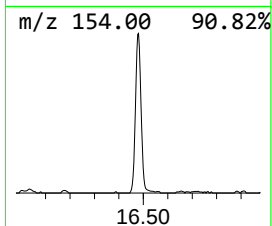
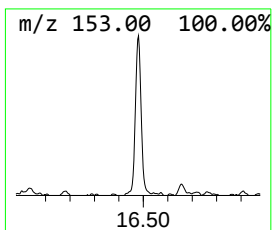
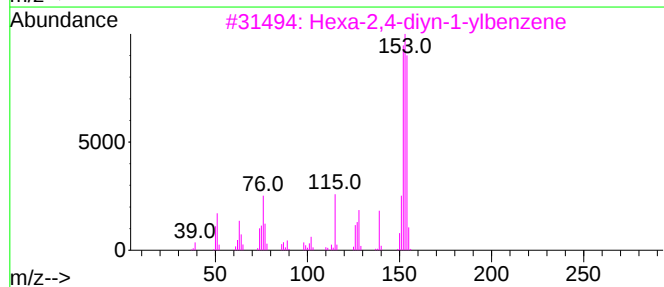
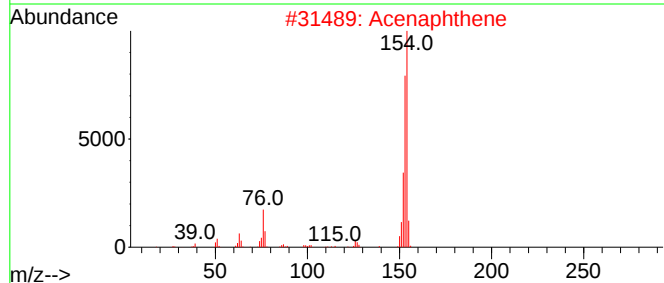
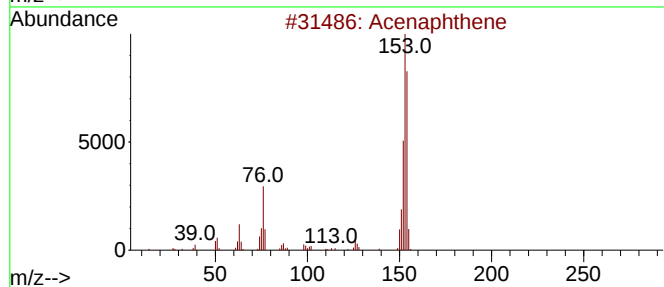
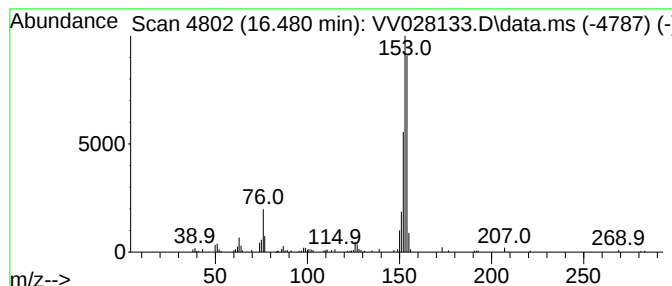
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 26 Acenaphthene Concentration Rank 23

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 16.480 | 0.58 ug/L | 119720 | 1,4-Dichlorobenzene-d4 | 11.246 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Acenaphthene                        | 154 | C12H10  | 000083-32-9 | 89   |
| 2    |    |   | Acenaphthene                        | 154 | C12H10  | 000083-32-9 | 81   |
| 3    |    |   | Hexa-2,4-diyn-1-ylbenzene           | 154 | C12H10  | 000520-74-1 | 59   |
| 4    |    |   | 1,4-Ethenonaphthalene, 1,4-dihydro- | 154 | C12H10  | 007322-47-6 | 53   |
| 5    |    |   | 2,4,6-Trihydroxybenzaldehyde        | 154 | C7H6O4  | 000487-70-7 | 52   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VW092722\  
 Data File : VW028133.D  
 Acq On : 27 Sep 2022 12:40  
 Operator : SY/MD  
 Sample : N4820-02  
 Misc : 25.0mL/MSVOA\_V/WATER  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 MSVOA\_V  
 ClientSampleId :  
 CB8L0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR092222WMA.M  
 Quant Title : TRACE VOA SFAM1.0

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| Benzene, propyl-   | 10.349 | 1.0     | ug/L  | 204335   | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, (2-met... | 11.053 | 1.2     | ug/L  | 253389   | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, 1-meth... | 11.085 | 2.0     | ug/L  | 418829   | 3                     | 11.246 | 1036560 | 5.0  |
| o-Cymene           | 11.448 | 0.7     | ug/L  | 151759   | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, 1,2-di... | 11.541 | 6.4     | ug/L  | 1324020  | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, 1,3-di... | 11.741 | 0.6     | ug/L  | 127856   | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, 1-ethy... | 12.011 | 24.1    | ug/L  | 5004450  | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, 1-ethe... | 12.111 | 5.5     | ug/L  | 1147860  | 3                     | 11.246 | 1036560 | 5.0  |
| 1H-Indene, 2,3-... | 12.294 | 0.5     | ug/L  | 104355   | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, 1,2,3,... | 12.410 | 13.5    | ug/L  | 2793890  | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, 1-meth... | 12.545 | 0.8     | ug/L  | 176327   | 3                     | 11.246 | 1036560 | 5.0  |
| .alpha.-Camphol... | 12.657 | 1.3     | ug/L  | 276544   | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, 1-ethe... | 12.718 | 3.8     | ug/L  | 795112   | 3                     | 11.246 | 1036560 | 5.0  |
| 2,4-Dimethylsty... | 12.876 | 14.5    | ug/L  | 3005930  | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, 1-ethy... | 12.921 | 0.6     | ug/L  | 129046   | 3                     | 11.246 | 1036560 | 5.0  |
| 3,5-Dimethylben... | 12.995 | 0.9     | ug/L  | 184878   | 3                     | 11.246 | 1036560 | 5.0  |
| 1H-Indene, 2,3-... | 13.214 | 0.6     | ug/L  | 115357   | 3                     | 11.246 | 1036560 | 5.0  |
| 4-tert-Butyltol... | 13.268 | 2.3     | ug/L  | 485845   | 3                     | 11.246 | 1036560 | 5.0  |
| 1H-Indene, 2,3-... | 13.365 | 0.6     | ug/L  | 132397   | 3                     | 11.246 | 1036560 | 5.0  |
| Benzene, 2,4-di... | 13.397 | 1.1     | ug/L  | 219850   | 3                     | 11.246 | 1036560 | 5.0  |
| 1H-Indene, 2,3-... | 13.905 | 0.7     | ug/L  | 138189   | 3                     | 11.246 | 1036560 | 5.0  |
| 1H-Indene, 2,3-... | 14.088 | 0.9     | ug/L  | 192760   | 3                     | 11.246 | 1036560 | 5.0  |
| Naphthalene, 1,... | 15.660 | 0.6     | ug/L  | 114120   | 3                     | 11.246 | 1036560 | 5.0  |
| Naphthalene, 1,... | 15.805 | 0.8     | ug/L  | 160552   | 3                     | 11.246 | 1036560 | 5.0  |
| Acenaphthene       | 16.480 | 0.6     | ug/L  | 119720   | 3                     | 11.246 | 1036560 | 5.0  |