

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
 Data File : VU060027.D  
 Acq On : 18 Jul 2024 19:11  
 Operator : MD/SY  
 Sample : P3217-03DL 100X  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 YDZC5DL

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\_U\Method\SFAMUTR071824WMA.M

Title : TRACE VOA SFAM1.0

Signal : TIC: VU060027.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         | 1.595       | 86            | 95          | 112          | rVB2     | 83184          | 101740        | 3.23%           | 0.502%        |
| 2         | 1.911       | 184           | 193         | 207          | rBV      | 54477          | 76023         | 2.42%           | 0.375%        |
| 3         | 2.560       | 384           | 395         | 410          | rBV      | 97432          | 160972        | 5.12%           | 0.795%        |
| 4         | 2.679       | 413           | 432         | 449          | rVB3     | 19941          | 43052         | 1.37%           | 0.213%        |
| 5         | 4.656       | 1028          | 1047        | 1088         | rBV      | 1320587        | 3146216       | 100.00%         | 15.538%       |
| 6         | 5.055       | 1156          | 1171        | 1198         | rVB      | 97292          | 221498        | 7.04%           | 1.094%        |
| 7         | 5.717       | 1358          | 1377        | 1410         | rBV2     | 198430         | 538219        | 17.11%          | 2.658%        |
| 8         | 6.245       | 1527          | 1541        | 1568         | rBV2     | 159614         | 312453        | 9.93%           | 1.543%        |
| 9         | 6.537       | 1620          | 1632        | 1654         | rVB3     | 58925          | 113220        | 3.60%           | 0.559%        |
| 10        | 6.685       | 1664          | 1678        | 1697         | rBV2     | 123546         | 239932        | 7.63%           | 1.185%        |
| 11        | 7.560       | 1938          | 1950        | 1968         | rBV      | 54666          | 99861         | 3.17%           | 0.493%        |
| 12        | 7.894       | 2041          | 2054        | 2068         | rBV      | 237268         | 411282        | 13.07%          | 2.031%        |
| 13        | 7.965       | 2068          | 2076        | 2091         | rVB      | 47618          | 84433         | 2.68%           | 0.417%        |
| 14        | 8.177       | 2132          | 2142        | 2155         | rBV      | 33046          | 57724         | 1.83%           | 0.285%        |
| 15        | 8.547       | 2246          | 2257        | 2272         | rBV      | 133627         | 227636        | 7.24%           | 1.124%        |
| 16        | 8.627       | 2272          | 2282        | 2317         | rBV      | 286028         | 566111        | 17.99%          | 2.796%        |
| 17        | 9.412       | 2514          | 2526        | 2550         | rBV      | 237729         | 427473        | 13.59%          | 2.111%        |
| 18        | 9.566       | 2563          | 2574        | 2589         | rBV      | 99573          | 162457        | 5.16%           | 0.802%        |
| 19        | 9.688       | 2595          | 2612        | 2630         | rBV      | 278507         | 480664        | 15.28%          | 2.374%        |
| 20        | 10.093      | 2725          | 2738        | 2764         | rBV      | 374473         | 601589        | 19.12%          | 2.971%        |
| 21        | 10.749      | 2931          | 2942        | 2953         | rBV      | 116404         | 183632        | 5.84%           | 0.907%        |
| 22        | 10.897      | 2974          | 2988        | 3000         | rBV2     | 33786          | 66709         | 2.12%           | 0.329%        |
| 23        | 10.991      | 3006          | 3017        | 3023         | rBV      | 155010         | 260950        | 8.29%           | 1.289%        |
| 24        | 11.084      | 3035          | 3046        | 3063         | rVV      | 215125         | 351359        | 11.17%          | 1.735%        |
| 25        | 11.286      | 3098          | 3109        | 3123         | rBV      | 217680         | 345527        | 10.98%          | 1.706%        |
| 26        | 11.460      | 3153          | 3163        | 3186         | rVB      | 372115         | 590648        | 18.77%          | 2.917%        |
| 27        | 11.634      | 3200          | 3217        | 3229         | rBV6     | 41879          | 88734         | 2.82%           | 0.438%        |
| 28        | 11.740      | 3242          | 3250        | 3257         | rVV2     | 44932          | 62541         | 1.99%           | 0.309%        |
| 29        | 11.807      | 3257          | 3271        | 3286         | rVV3     | 292210         | 725980        | 23.07%          | 3.585%        |
| 30        | 11.888      | 3286          | 3296        | 3317         | rVB      | 559452         | 890725        | 28.31%          | 4.399%        |
| 31        | 12.090      | 3346          | 3359        | 3365         | rBV2     | 191766         | 323573        | 10.28%          | 1.598%        |
| 32        | 12.203      | 3377          | 3394        | 3413         | rVB4     | 642838         | 1584410       | 50.36%          | 7.825%        |
| 33        | 12.296      | 3414          | 3423        | 3437         | rBV5     | 28584          | 61117         | 1.94%           | 0.302%        |
| 34        | 12.370      | 3437          | 3446        | 3460         | rVB      | 83303          | 134868        | 4.29%           | 0.666%        |
| 35        | 12.460      | 3460          | 3474        | 3478         | rBV      | 93556          | 147540        | 4.69%           | 0.729%        |
| 36        | 12.489      | 3479          | 3483        | 3495         | rVV      | 97819          | 137511        | 4.37%           | 0.679%        |

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 Sample : P3217-03DL 100X  
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 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 YDZC5DL

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\_U\Method\SFAMUTR071824WMA.M  
 Title : TRACE VOA SFAM1.0

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 37 | 12.566 | 3496 | 3507 | 3514 | rVV  | 185639 | 282649  | 8.98%  | 1.396% |
| 38 | 12.605 | 3514 | 3519 | 3530 | rVB  | 51659  | 75465   | 2.40%  | 0.373% |
| 39 | 12.675 | 3530 | 3541 | 3561 | rBV2 | 137445 | 310950  | 9.88%  | 1.536% |
| 40 | 12.868 | 3591 | 3601 | 3612 | rVB2 | 93064  | 155931  | 4.96%  | 0.770% |
| 41 | 12.965 | 3614 | 3631 | 3640 | rBV2 | 125908 | 249894  | 7.94%  | 1.234% |
| 42 | 13.019 | 3640 | 3648 | 3664 | rVB  | 218051 | 335699  | 10.67% | 1.658% |
| 43 | 13.106 | 3664 | 3675 | 3679 | rBV5 | 20867  | 37876   | 1.20%  | 0.187% |
| 44 | 13.248 | 3710 | 3719 | 3723 | rBV3 | 25727  | 35867   | 1.14%  | 0.177% |
| 45 | 13.286 | 3723 | 3731 | 3741 | rVB  | 80463  | 125967  | 4.00%  | 0.622% |
| 46 | 13.399 | 3757 | 3766 | 3770 | rBV4 | 28968  | 44438   | 1.41%  | 0.219% |
| 47 | 13.447 | 3770 | 3781 | 3793 | rVV3 | 345108 | 618794  | 19.67% | 3.056% |
| 48 | 13.511 | 3793 | 3801 | 3809 | rVB4 | 55676  | 92280   | 2.93%  | 0.456% |
| 49 | 13.617 | 3824 | 3834 | 3852 | rVB  | 142692 | 239146  | 7.60%  | 1.181% |
| 50 | 13.730 | 3857 | 3869 | 3877 | rBV5 | 43865  | 78265   | 2.49%  | 0.387% |
| 51 | 13.778 | 3877 | 3884 | 3891 | rVV2 | 25764  | 34107   | 1.08%  | 0.168% |
| 52 | 13.833 | 3891 | 3901 | 3916 | rBV8 | 60768  | 124331  | 3.95%  | 0.614% |
| 53 | 13.936 | 3927 | 3933 | 3949 | rVB2 | 41780  | 91326   | 2.90%  | 0.451% |
| 54 | 14.080 | 3963 | 3978 | 3998 | rVB  | 225490 | 405532  | 12.89% | 2.003% |
| 55 | 14.183 | 3999 | 4010 | 4024 | rVB4 | 35322  | 62870   | 2.00%  | 0.310% |
| 56 | 14.283 | 4027 | 4041 | 4051 | rBV4 | 45986  | 91483   | 2.91%  | 0.452% |
| 57 | 14.341 | 4051 | 4059 | 4069 | rVB3 | 23964  | 38068   | 1.21%  | 0.188% |
| 58 | 14.479 | 4093 | 4102 | 4118 | rVB2 | 71510  | 122752  | 3.90%  | 0.606% |
| 59 | 14.666 | 4147 | 4160 | 4177 | rBV4 | 101079 | 256245  | 8.14%  | 1.266% |
| 60 | 14.804 | 4194 | 4203 | 4213 | rBV3 | 38759  | 63437   | 2.02%  | 0.313% |
| 61 | 14.868 | 4213 | 4223 | 4234 | rVB2 | 83627  | 134940  | 4.29%  | 0.666% |
| 62 | 15.010 | 4256 | 4267 | 4281 | rBV3 | 50495  | 91051   | 2.89%  | 0.450% |
| 63 | 15.215 | 4315 | 4331 | 4344 | rBV  | 706897 | 1115319 | 35.45% | 5.508% |
| 64 | 15.402 | 4377 | 4389 | 4402 | rBV  | 436789 | 703295  | 22.35% | 3.473% |
| 65 | 16.260 | 4642 | 4656 | 4676 | rBV3 | 39257  | 80342   | 2.55%  | 0.397% |
| 66 | 16.405 | 4690 | 4701 | 4708 | rBV2 | 48067  | 79887   | 2.54%  | 0.395% |
| 67 | 16.444 | 4708 | 4713 | 4728 | rVB4 | 26146  | 41512   | 1.32%  | 0.205% |

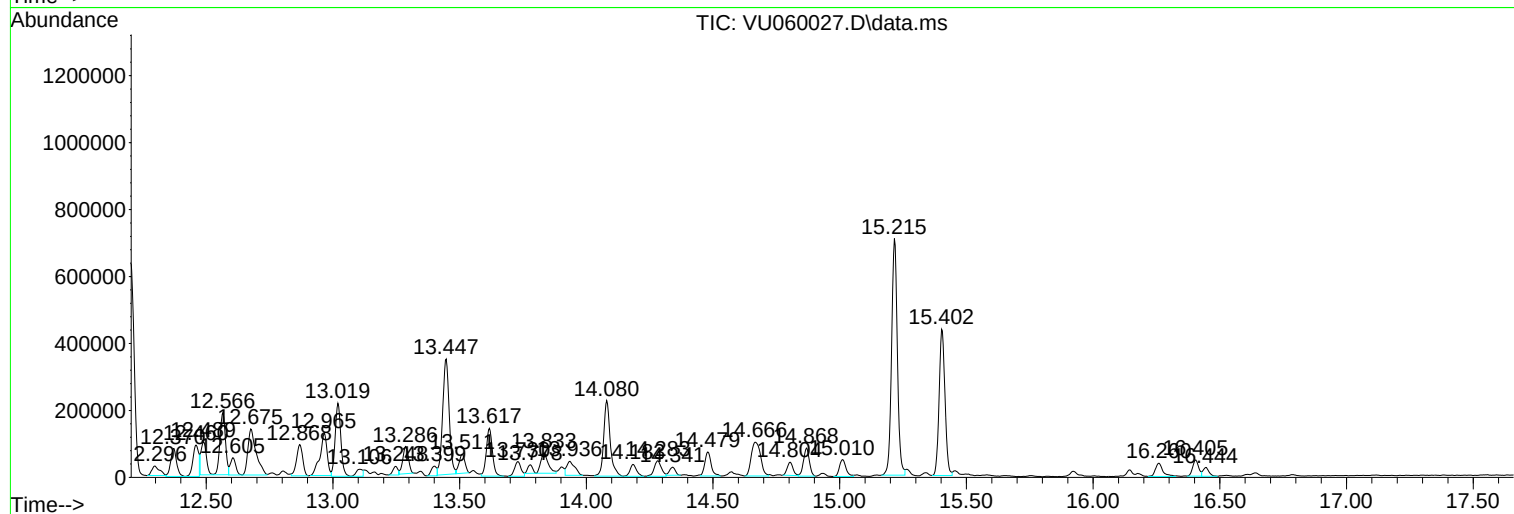
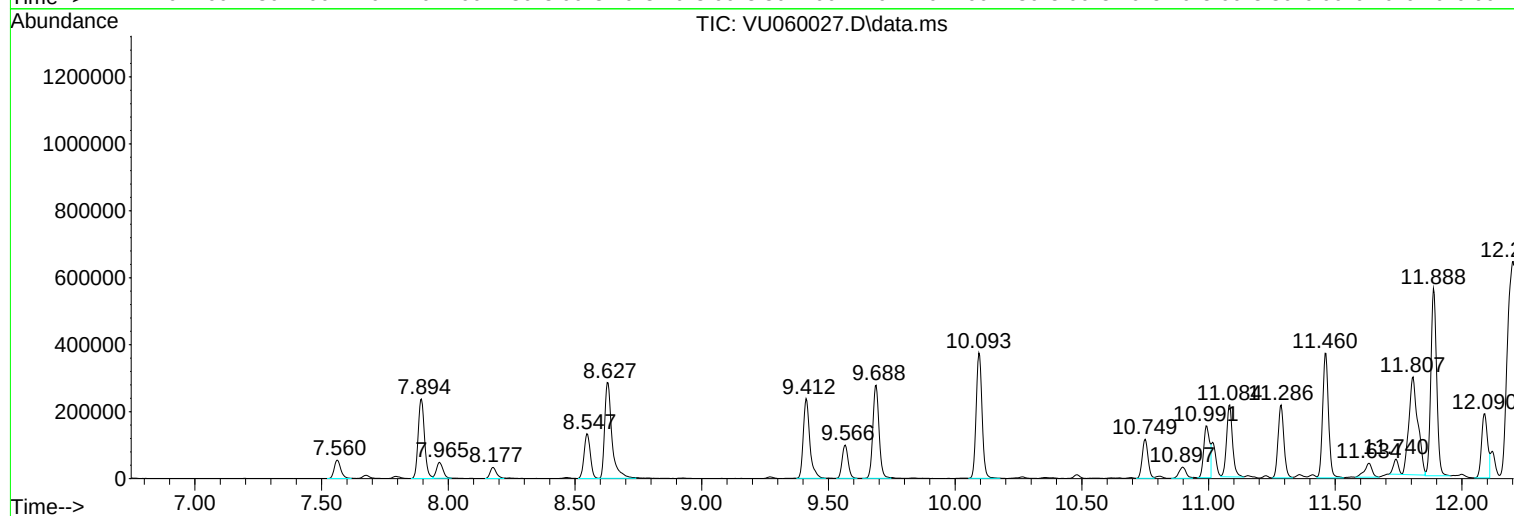
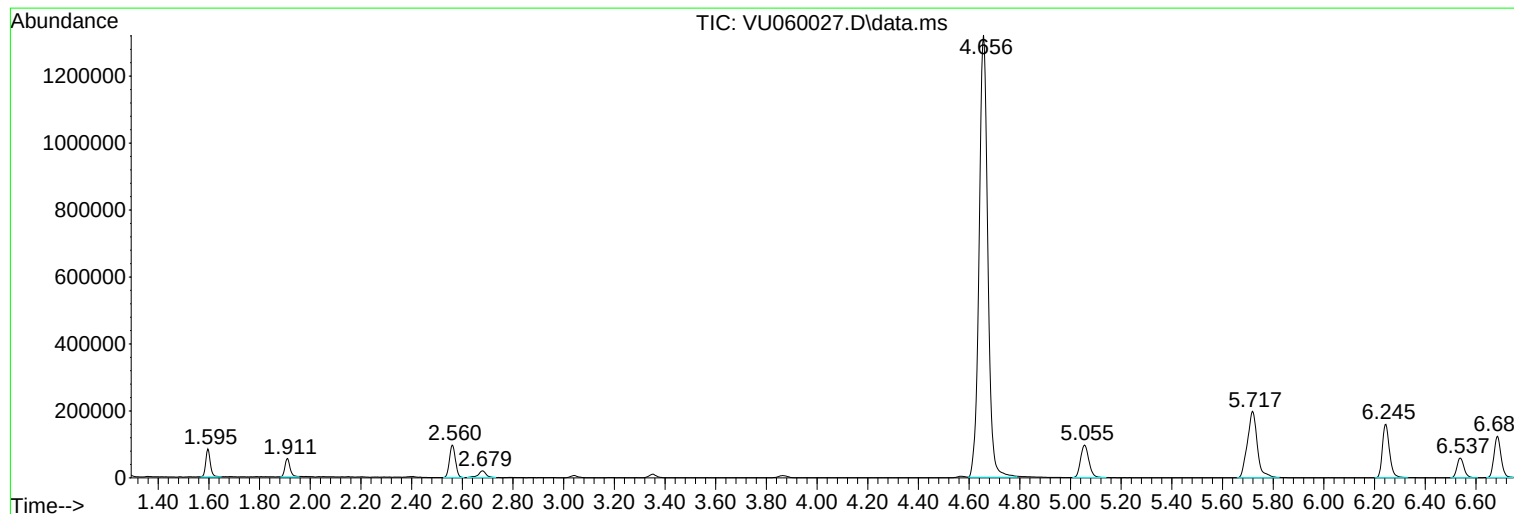
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Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

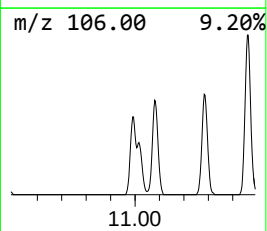
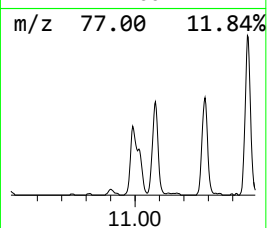
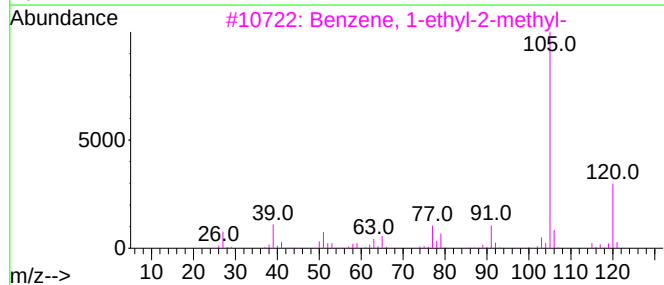
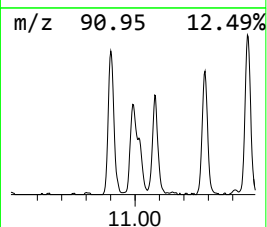
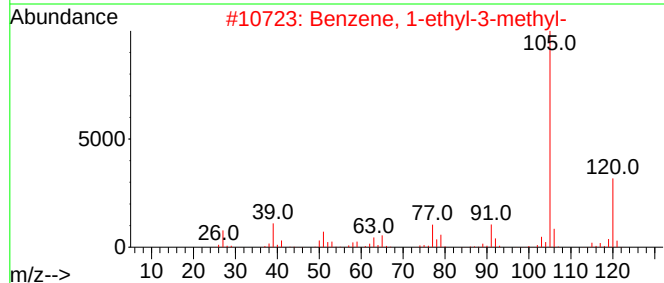
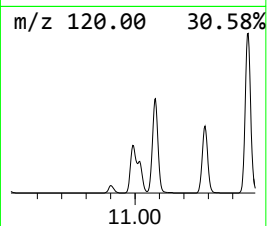
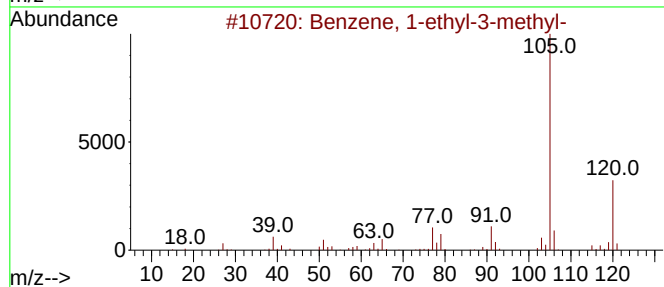
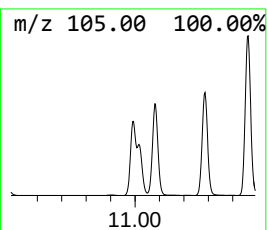
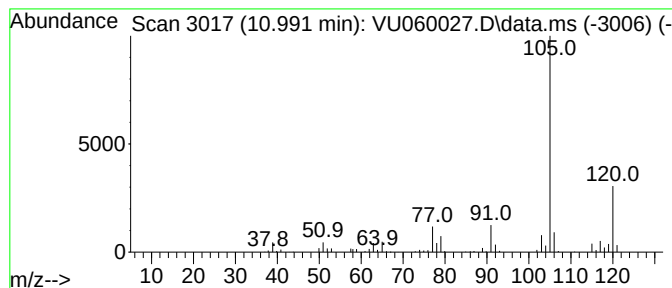
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 11

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 10.991 | 1.80 ug/L | 260950 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 95   |
| 2    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 94   |
| 3    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 4    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 5    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |



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ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

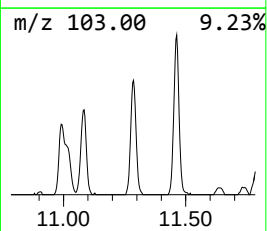
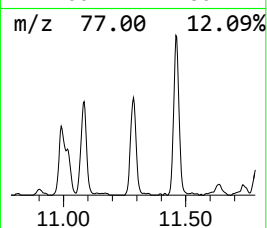
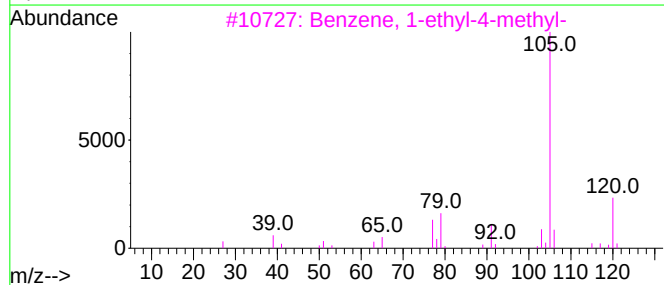
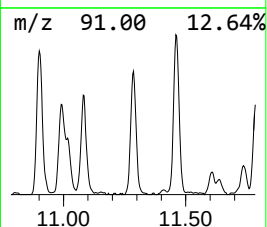
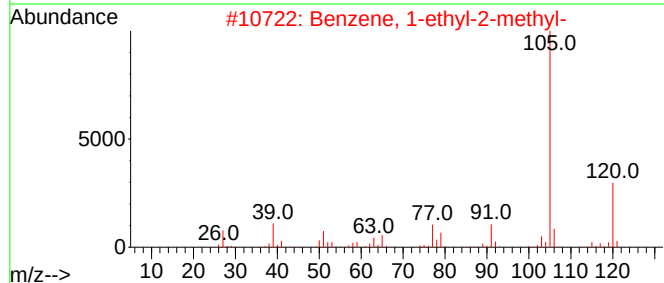
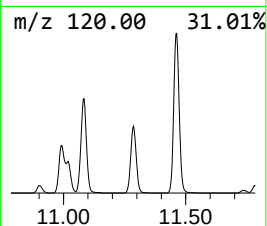
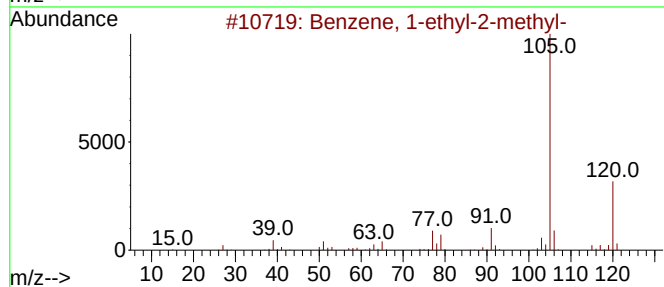
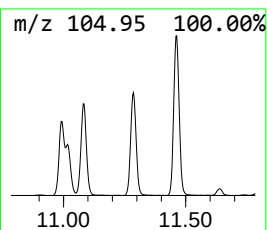
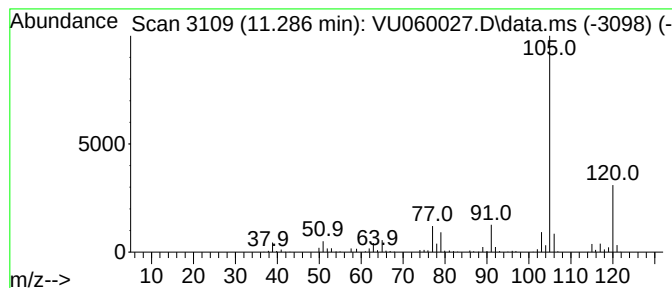
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 11.286 | 2.38 ug/L | 345527 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 2    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 3    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 93   |
| 4    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |
| 5    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |





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ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

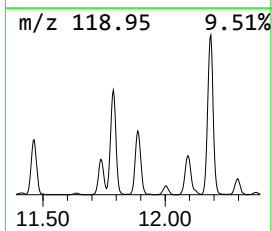
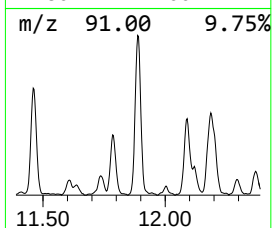
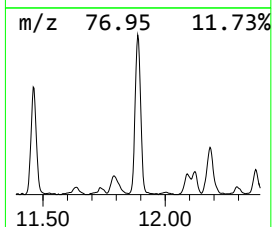
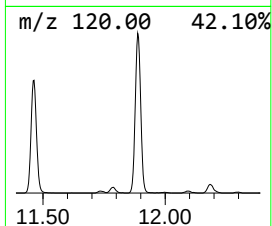
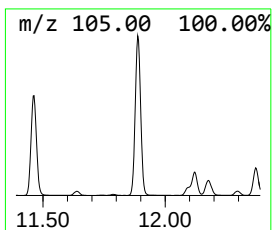
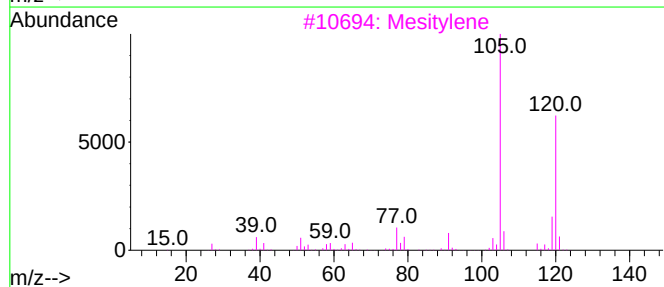
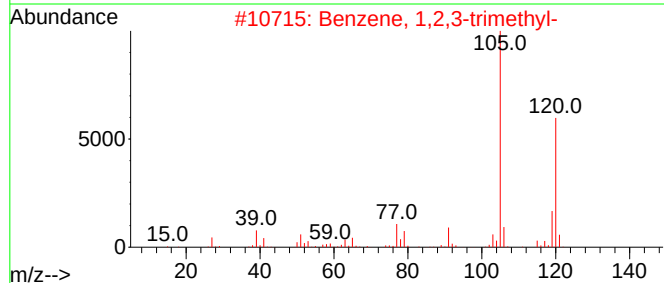
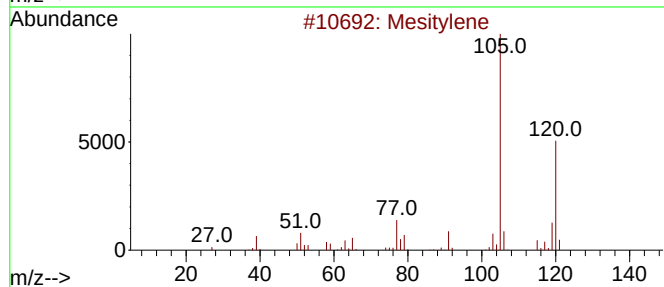
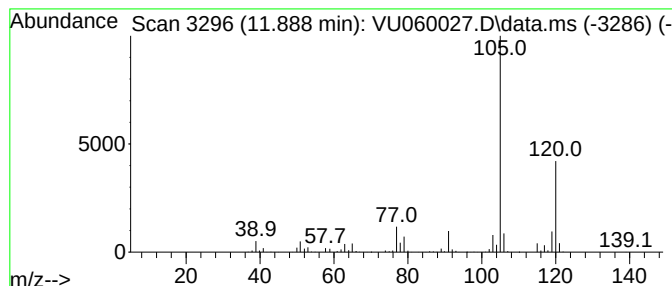
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 2

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 11.888 | 6.13 ug/L | 890725 | 1,4-Dichlorobenzene-d4 | 11.807 |

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

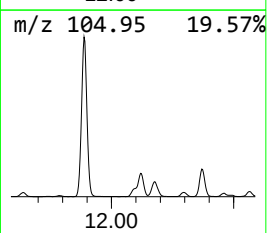
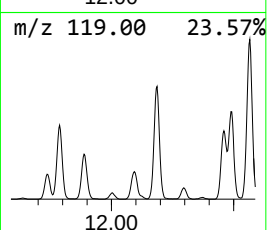
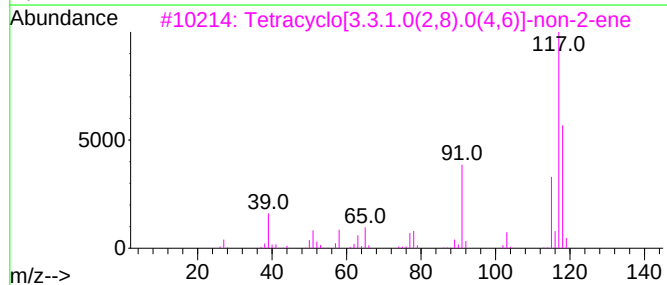
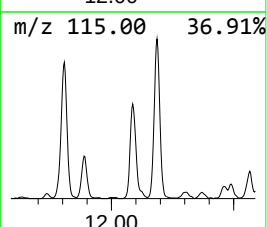
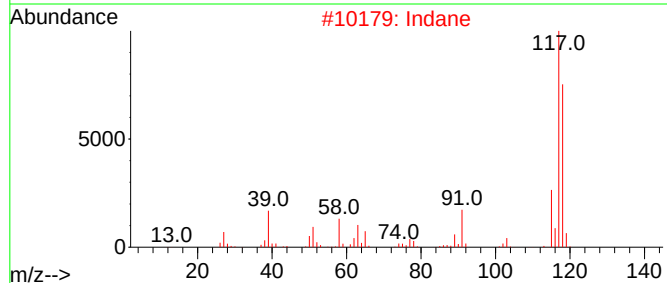
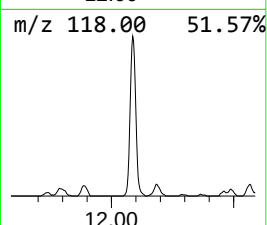
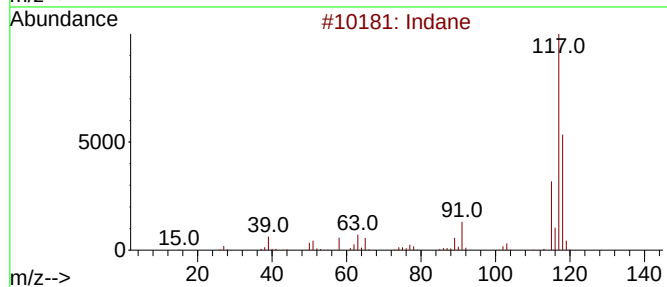
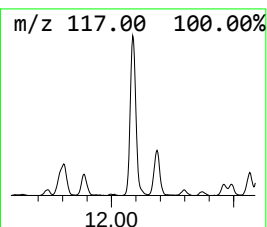
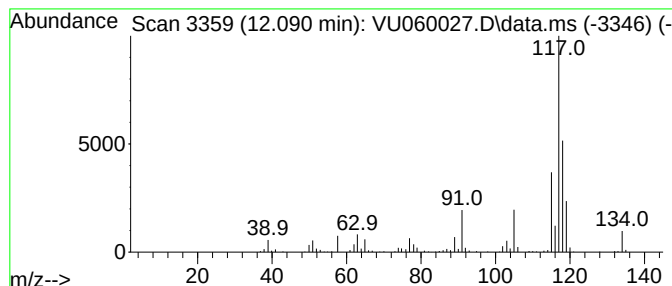
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 6 Indane Concentration Rank 8

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.090 | 2.23 ug/L | 323573 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Indane                              | 118 | C9H10   | 000496-11-7  | 70   |
| 2    |    |   | Indane                              | 118 | C9H10   | 000496-11-7  | 70   |
| 3    |    |   | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118 | C9H10   | 1000191-13-7 | 64   |
| 4    |    |   | Benzene, cyclopropyl-               | 118 | C9H10   | 000873-49-4  | 60   |
| 5    |    |   | Benzene, cyclopropyl-               | 118 | C9H10   | 000873-49-4  | 58   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

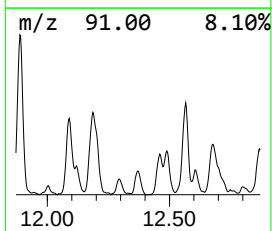
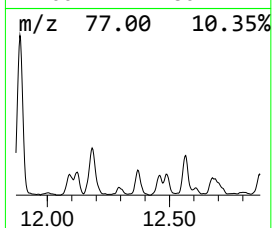
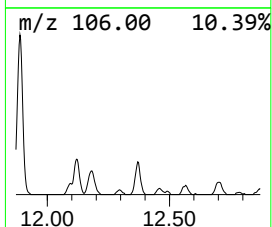
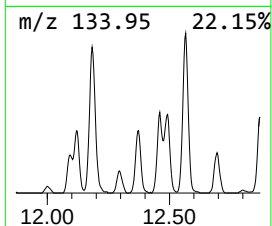
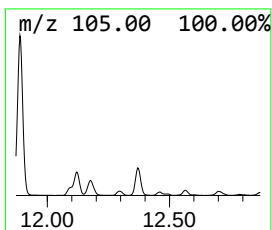
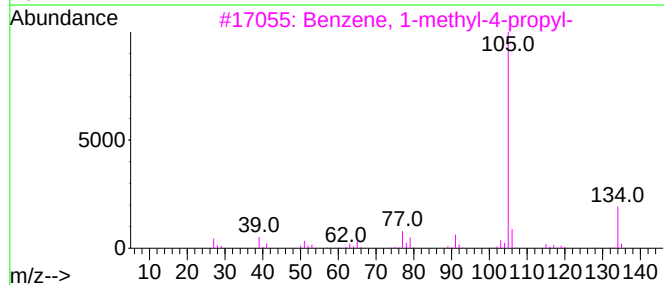
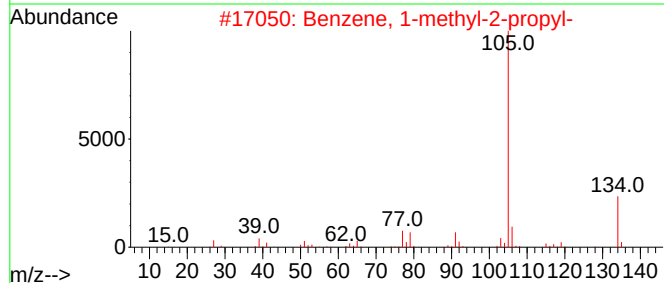
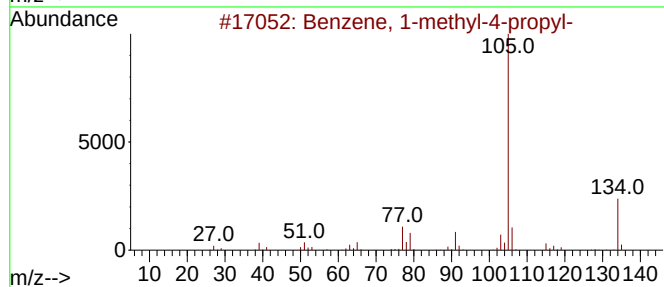
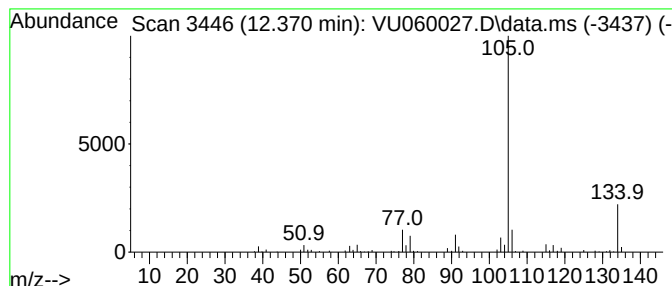
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 7 Benzene, 1-methyl-4-propyl- Concentration Rank 20

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.370 | 0.93 ug/L | 134868 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 94   |
| 2    |    |   | Benzene, 1-methyl-2-propyl- | 134 | C10H14  | 001074-17-5 | 91   |
| 3    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 91   |
| 4    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 91   |
| 5    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 90   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

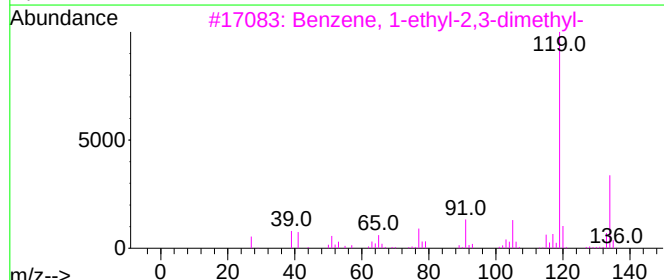
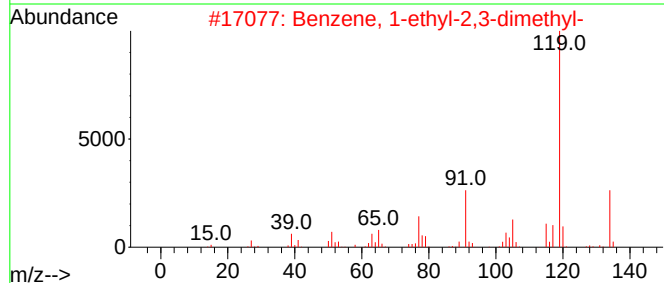
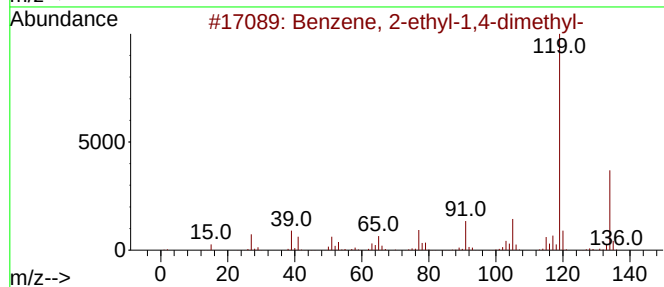
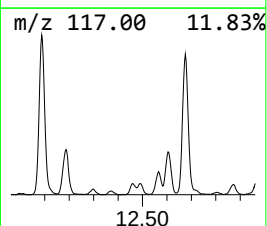
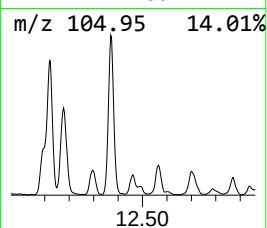
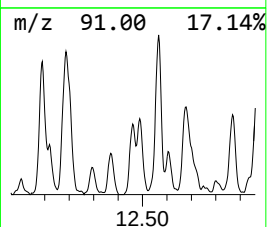
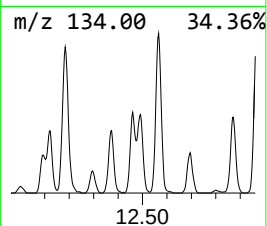
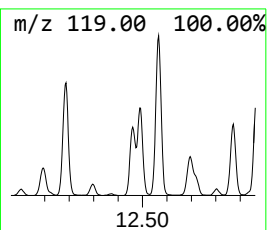
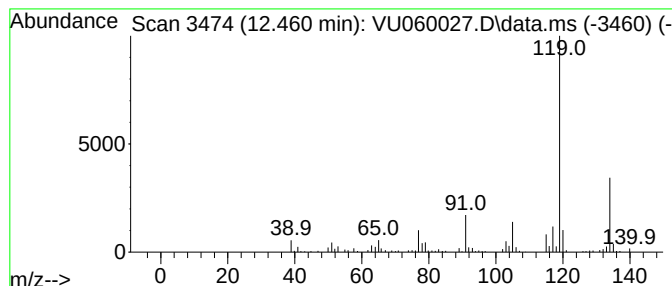
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 8 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 17

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.460 | 1.02 ug/L | 147540 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 97   |
| 2    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 96   |
| 3    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 95   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 95   |
| 5    |    |   | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 95   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

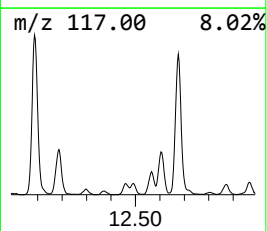
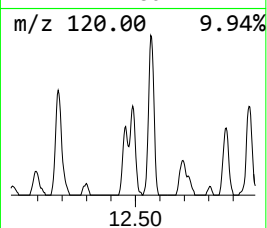
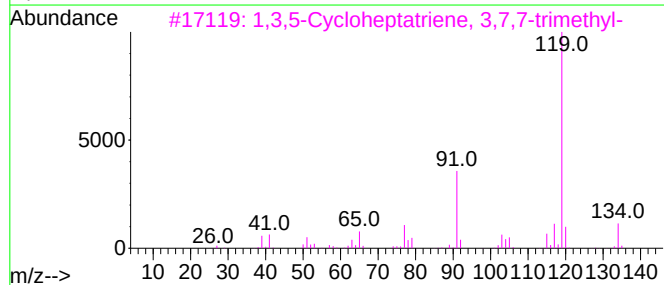
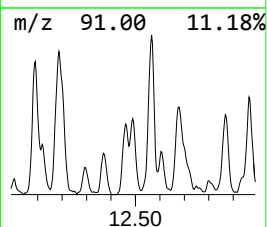
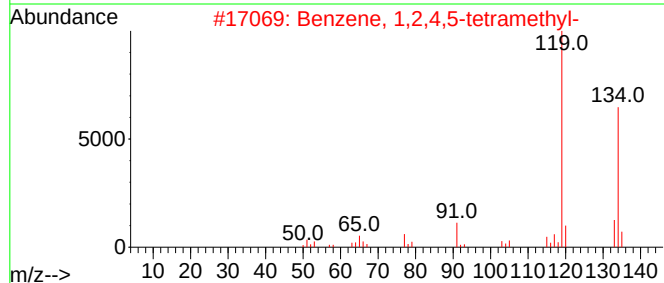
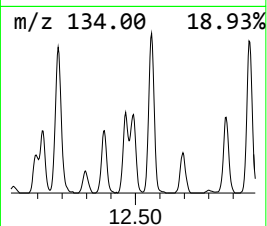
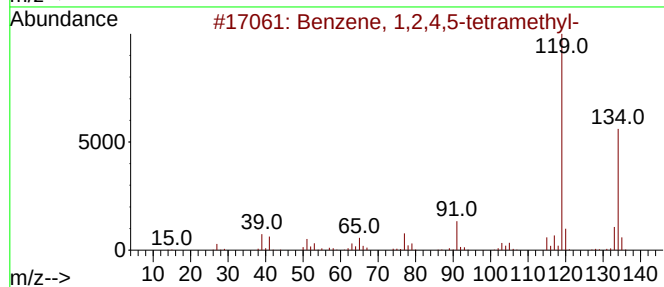
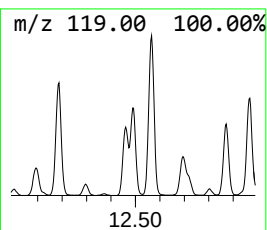
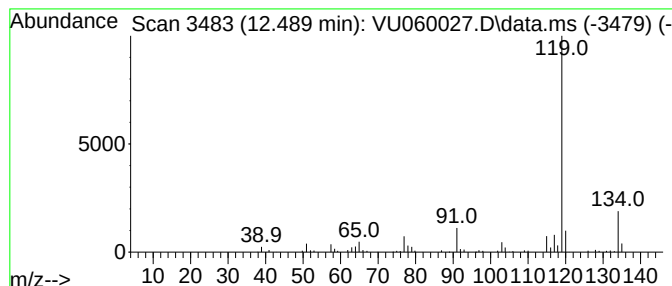
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 18

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.489 | 0.95 ug/L | 137511 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 91   |
| 2    |    |   | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 91   |
| 3    |    |   | 1,3,5-Cycloheptatriene, 3,7,7-tr... | 134 | C10H14  | 003479-89-8 | 91   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 91   |
| 5    |    |   | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

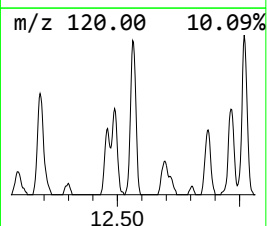
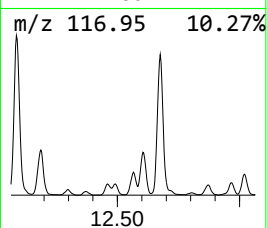
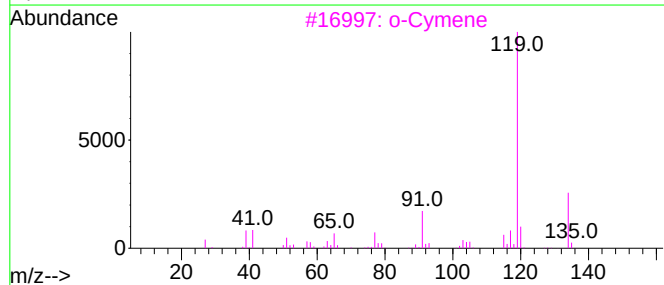
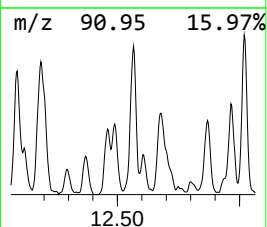
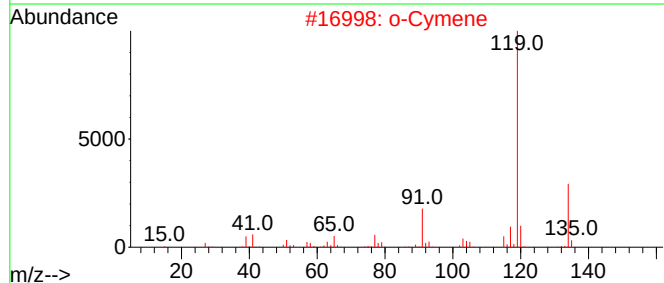
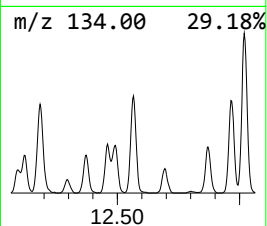
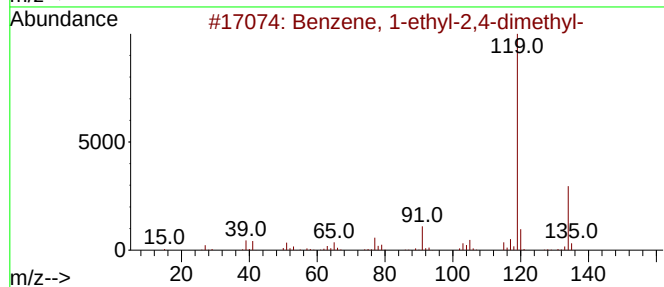
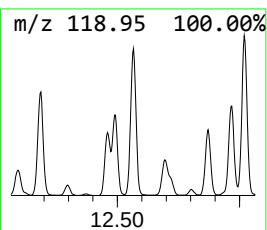
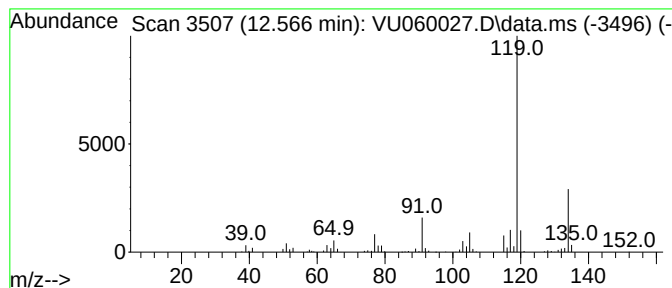
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Quant Title : TRACE VOA SFAM1.0

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

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

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.566 | 1.95 ug/L | 282649 | 1,4-Dichlorobenzene-d4 | 11.807 |

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

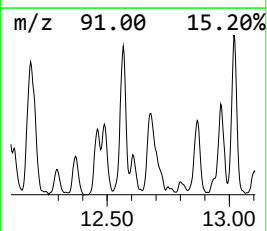
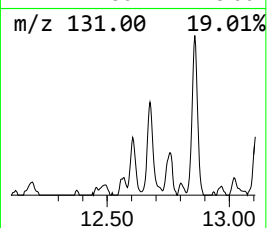
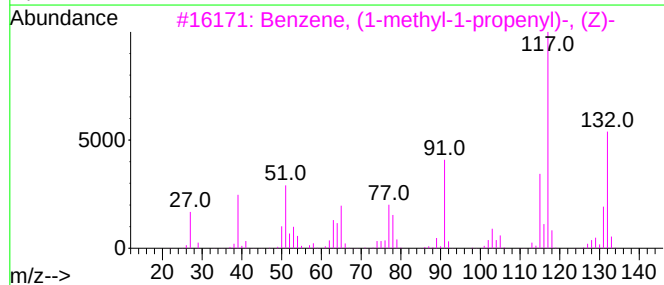
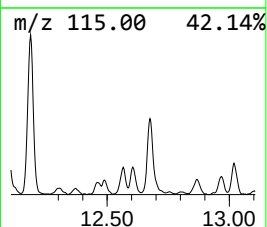
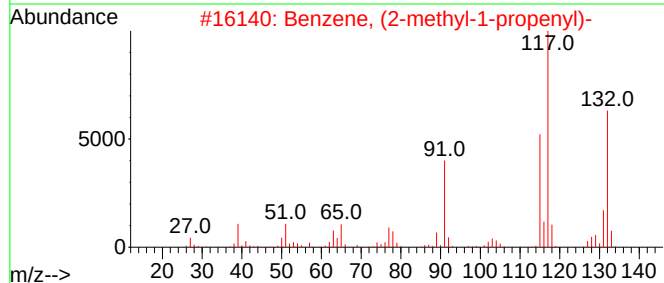
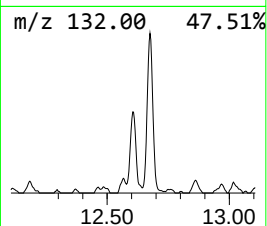
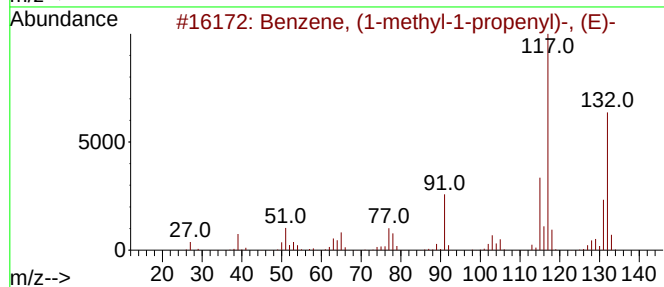
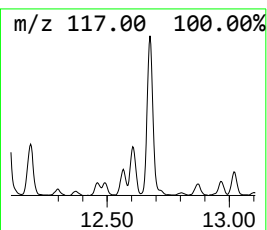
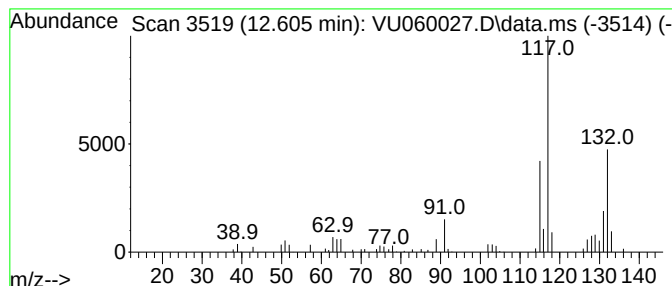
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 11 Benzene, (1-methyl-1-propen... Concentration Rank 32

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 12.605 | 0.52 ug/L | 75465 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 91   |
| 2    |    |   | Benzene, (2-methyl-1-propenyl)-     | 132 | C10H12  | 000768-49-0 | 91   |
| 3    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000767-99-7 | 91   |
| 4    |    |   | Benzene, (2-methyl-2-propenyl)-     | 132 | C10H12  | 003290-53-7 | 91   |
| 5    |    |   | Benzene, 1-methyl-4-(2-propenyl)-   | 132 | C10H12  | 003333-13-9 | 90   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

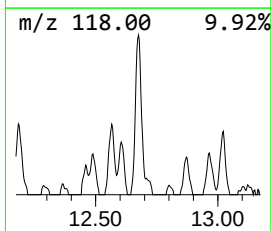
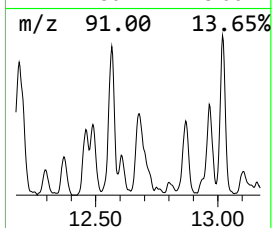
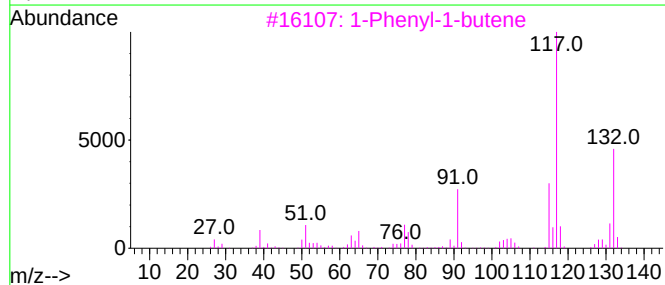
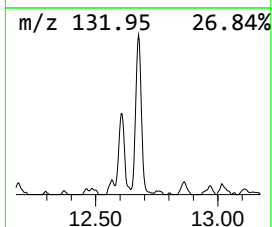
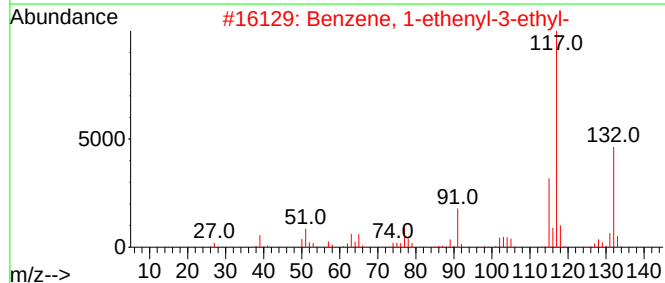
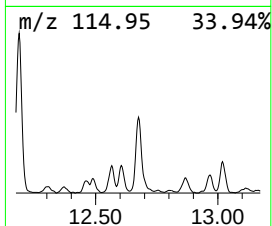
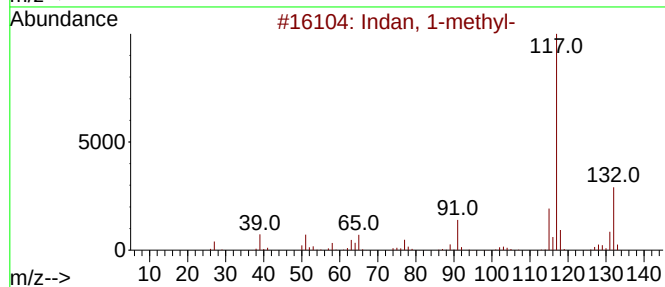
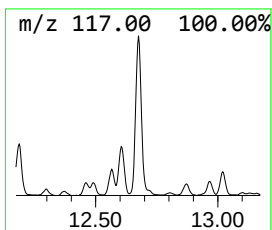
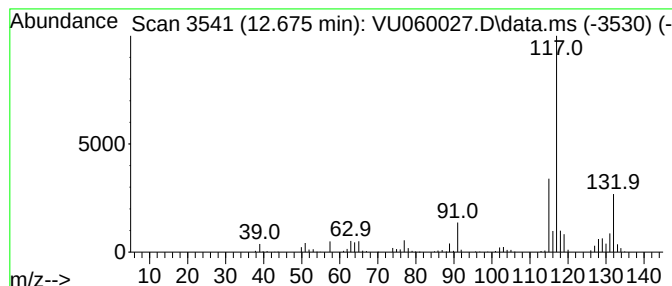
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 12 Indan, 1-methyl- Concentration Rank 9

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.675 | 2.14 ug/L | 310950 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Indan, 1-methyl-                  | 132 | C10H12  | 000767-58-8 | 87   |
| 2    |    |   | Benzene, 1-ethenyl-3-ethyl-       | 132 | C10H12  | 007525-62-4 | 87   |
| 3    |    |   | 1-Phenyl-1-butene                 | 132 | C10H12  | 000824-90-8 | 86   |
| 4    |    |   | Benzene, 1-methyl-2-(2-propenyl)- | 132 | C10H12  | 001587-04-8 | 86   |
| 5    |    |   | Indan, 1-methyl-                  | 132 | C10H12  | 000767-58-8 | 81   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

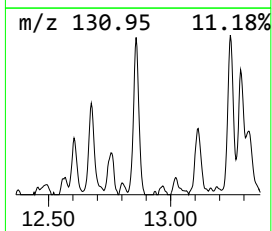
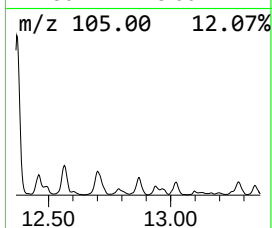
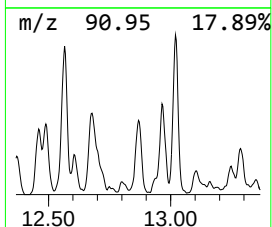
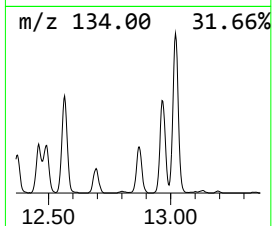
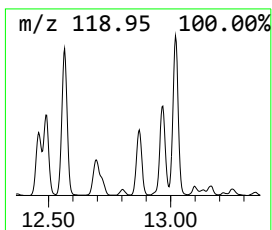
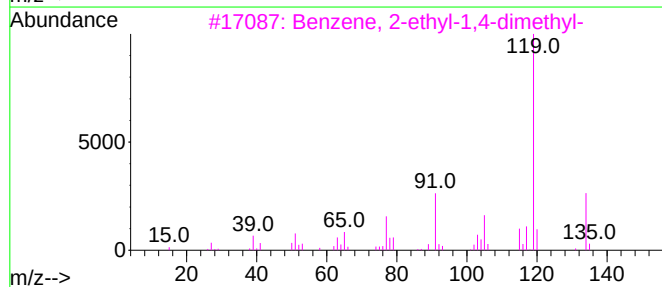
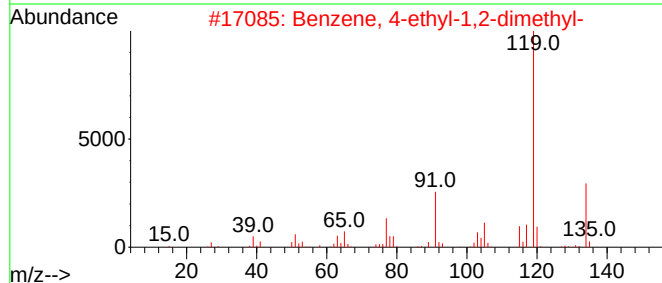
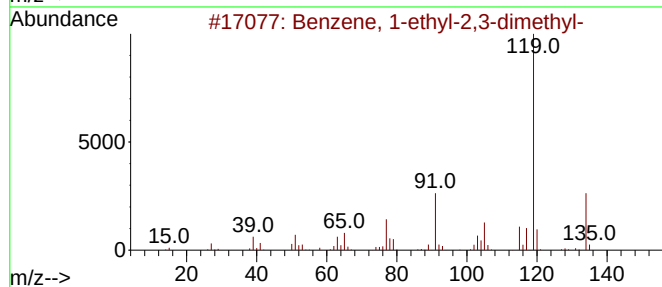
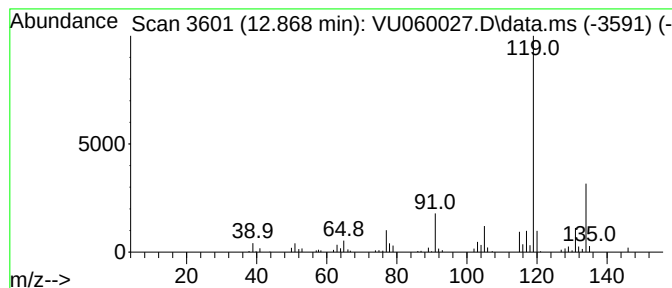
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 13 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 16

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.868 | 1.07 ug/L | 155931 | 1,4-Dichlorobenzene-d4 | 11.807 |

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

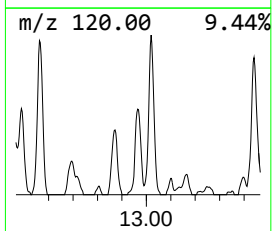
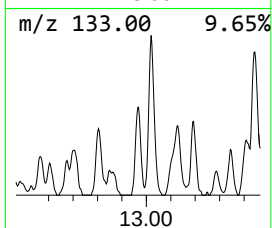
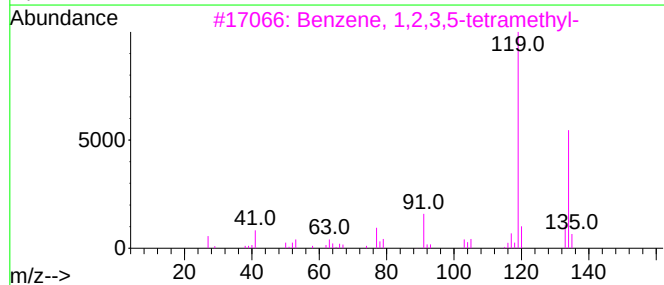
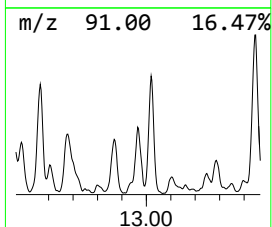
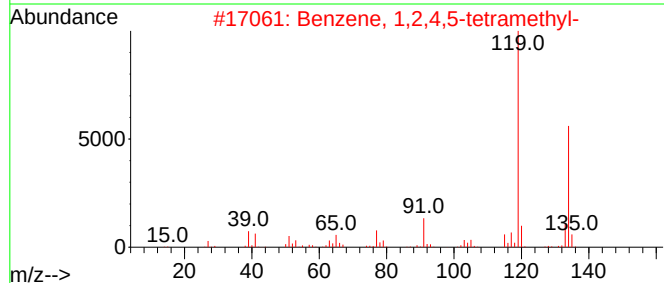
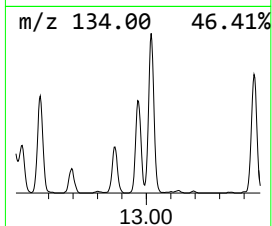
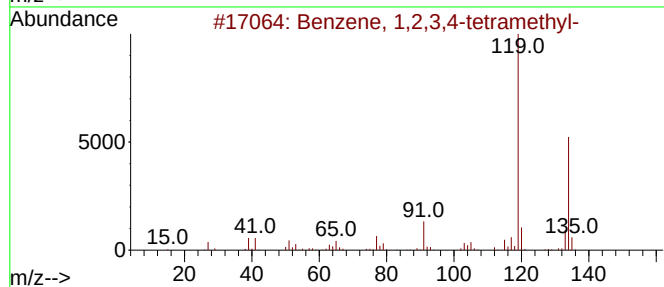
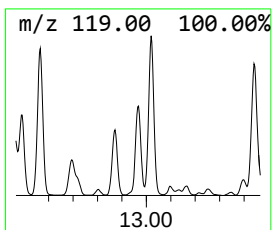
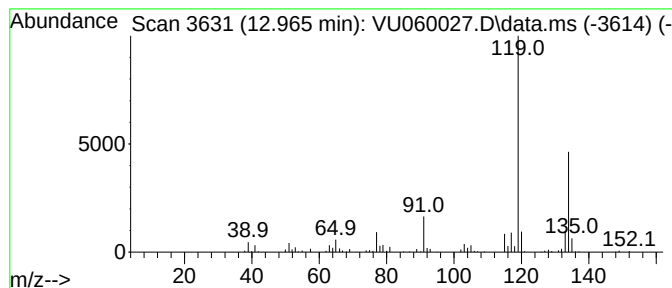
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 14 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 13

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 12.965 | 1.72 ug/L | 249894 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 97   |
| 2    |    |   | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 97   |
| 3    |    |   | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 96   |
| 4    |    |   | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 95   |
| 5    |    |   | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14  | 000095-93-2 | 95   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

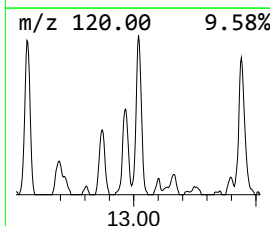
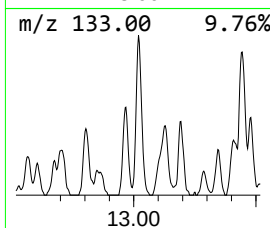
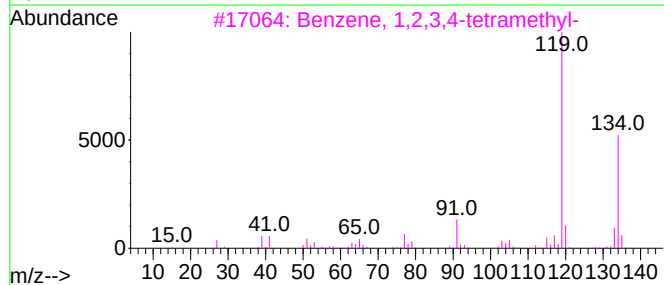
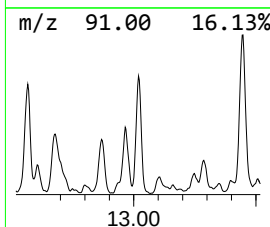
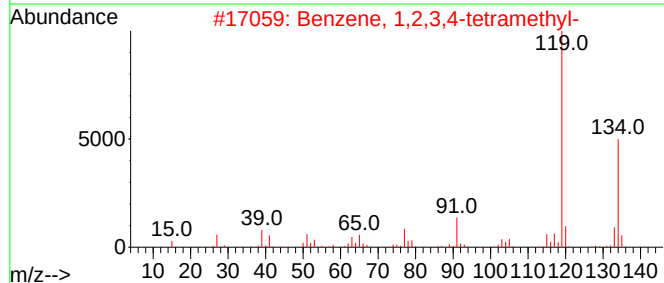
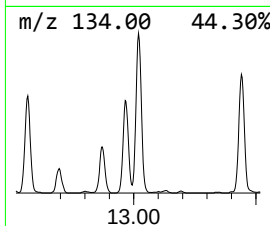
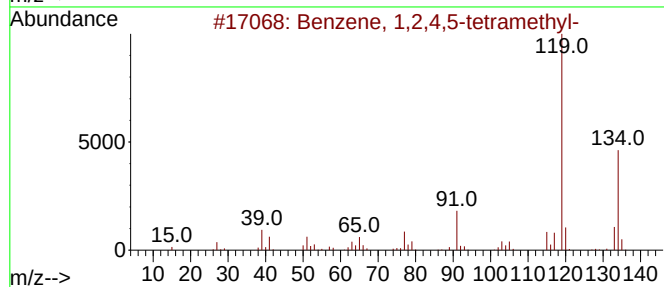
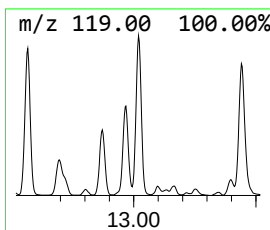
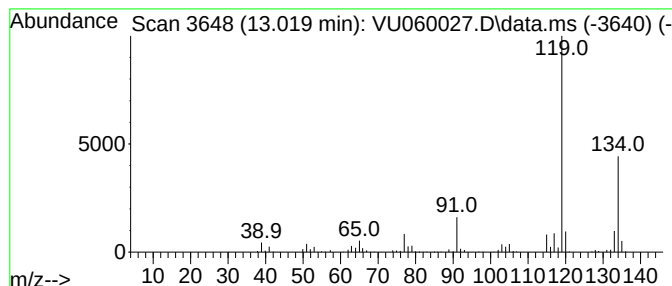
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 15 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 7

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.019 | 2.31 ug/L | 335699 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 97   |
| 2    |    |   | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 97   |
| 3    |    |   | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 97   |
| 4    |    |   | 1,3-Cyclopentadiene, 1,2,3,4-tet... | 134 | C10H14  | 076089-59-3 | 97   |
| 5    |    |   | Benzene, 1,2,3,5-tetramethyl-       | 134 | C10H14  | 000527-53-7 | 96   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

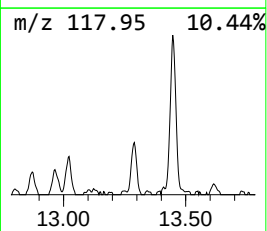
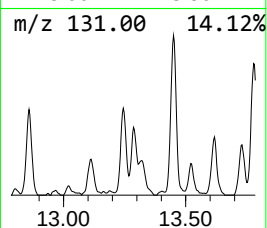
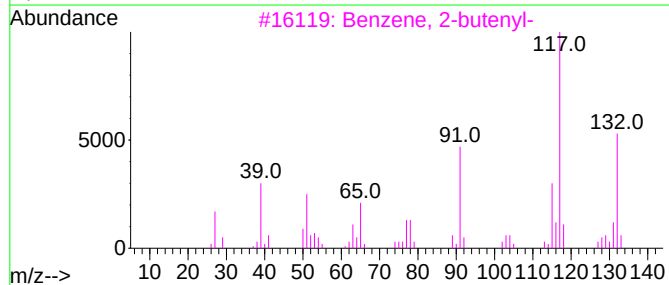
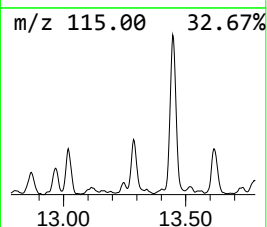
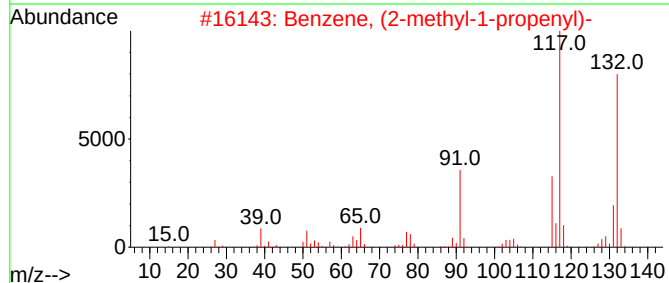
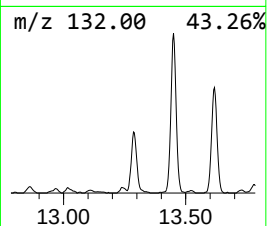
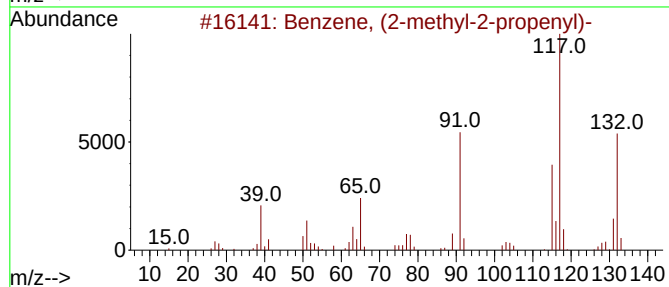
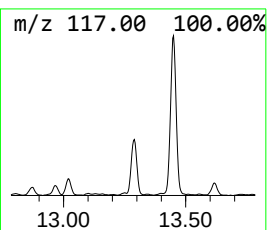
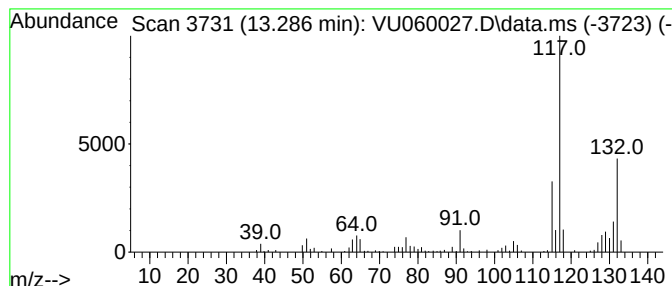
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzene, (2-methyl-2-propen... Concentration Rank 21

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.286 | 0.87 ug/L | 125967 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (2-methyl-2-propenyl)-  | 132 | C10H12  | 003290-53-7 | 90   |
| 2    |    |   | Benzene, (2-methyl-1-propenyl)-  | 132 | C10H12  | 000768-49-0 | 90   |
| 3    |    |   | Benzene, 2-butenyl-              | 132 | C10H12  | 001560-06-1 | 90   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-2-methyl- | 132 | C10H12  | 000824-63-5 | 86   |
| 5    |    |   | Benzene, 2-butenyl-              | 132 | C10H12  | 001560-06-1 | 86   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

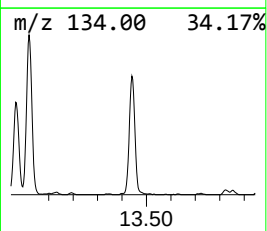
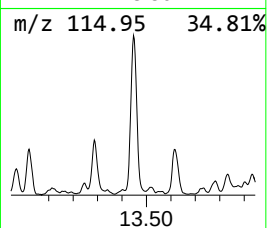
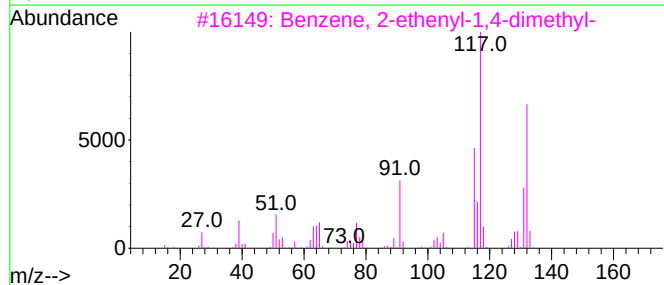
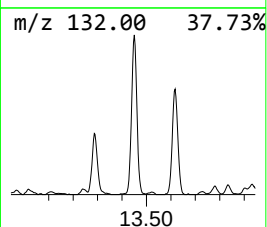
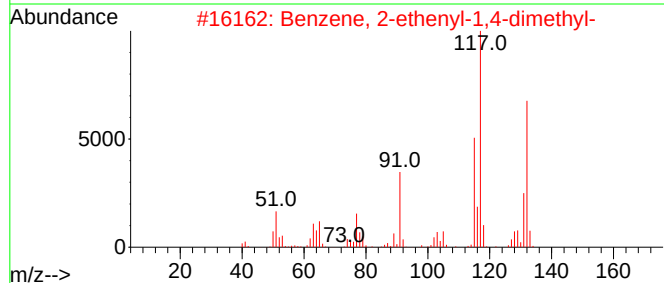
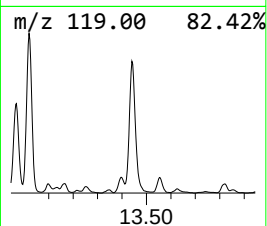
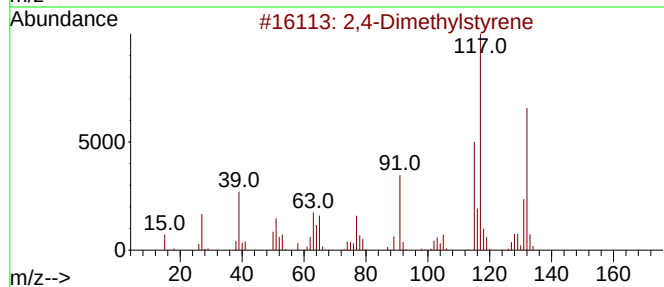
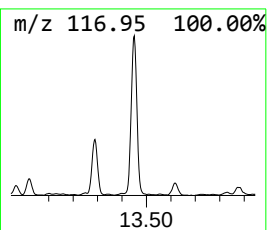
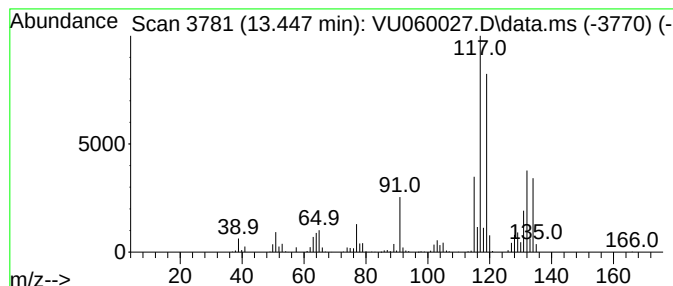
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 17 2,4-Dimethylstyrene Concentration Rank 4

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.447 | 4.26 ug/L | 618794 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2,4-Dimethylstyrene              | 132 | C10H12  | 002234-20-0 | 94   |
| 2    |    |   | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12  | 002039-89-6 | 94   |
| 3    |    |   | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12  | 002039-89-6 | 94   |
| 4    |    |   | 3-Phenylbut-1-ene                | 132 | C10H12  | 000934-10-1 | 60   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-4-methyl- | 132 | C10H12  | 000824-22-6 | 55   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

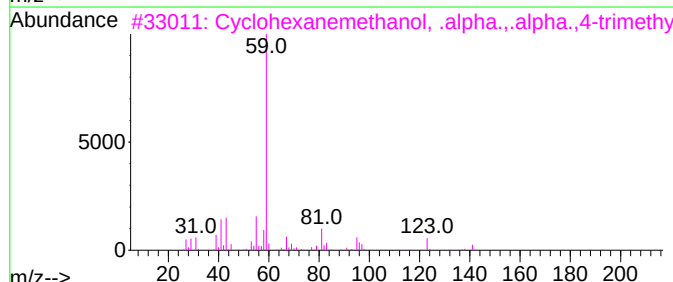
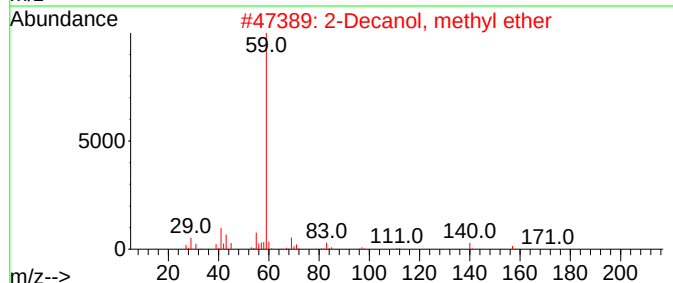
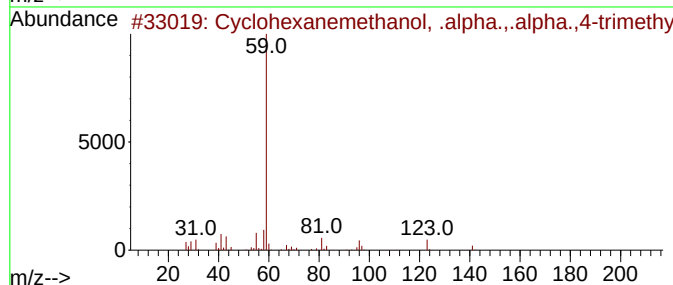
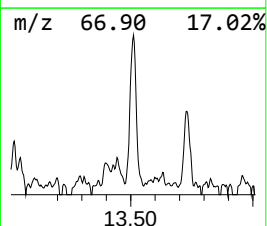
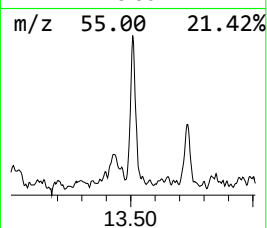
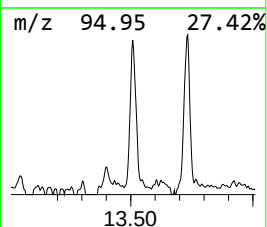
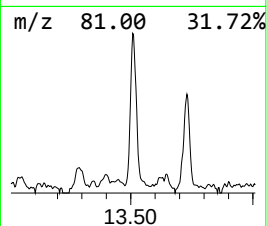
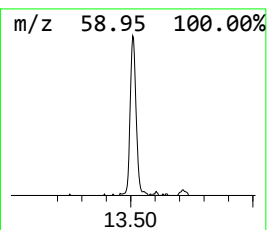
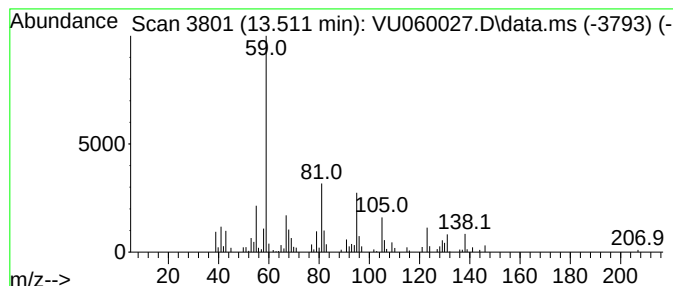
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 18 unknown-01 Concentration Rank 24

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 13.511 | 0.64 ug/L | 92280 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Cyclohexanemethanol, .alpha.,.al... | 156 | C10H20O | 007322-63-6  | 43   |
| 2    |    |   | 2-Decanol, methyl ether             | 172 | C11H24O | 1000333-83-9 | 38   |
| 3    |    |   | Cyclohexanemethanol, .alpha.,.al... | 156 | C10H20O | 000498-81-7  | 38   |
| 4    |    |   | Heptane, 1-ethoxy-                  | 144 | C9H20O  | 001969-43-3  | 38   |
| 5    |    |   | Pentane, 2-methoxy-                 | 102 | C6H14O  | 006795-88-6  | 38   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

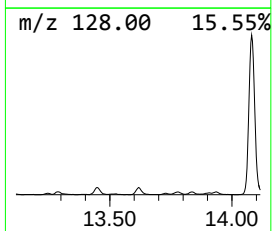
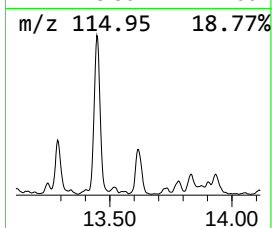
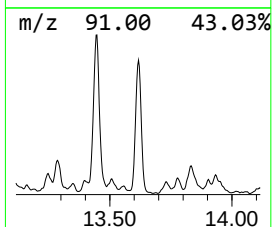
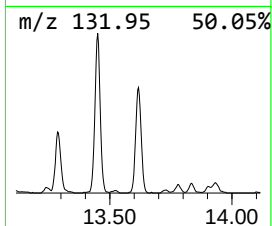
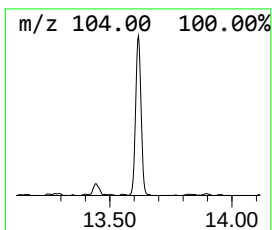
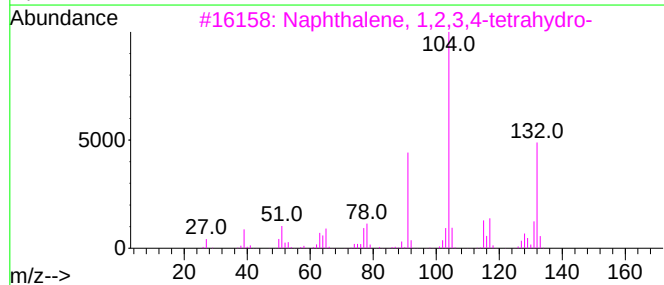
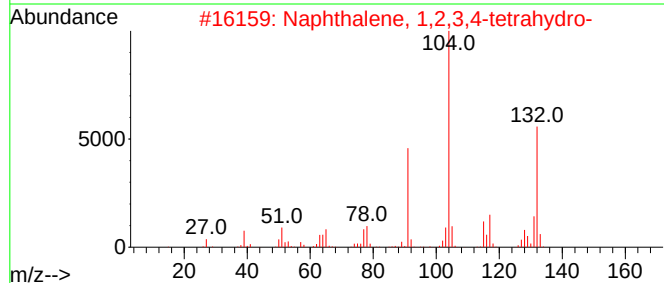
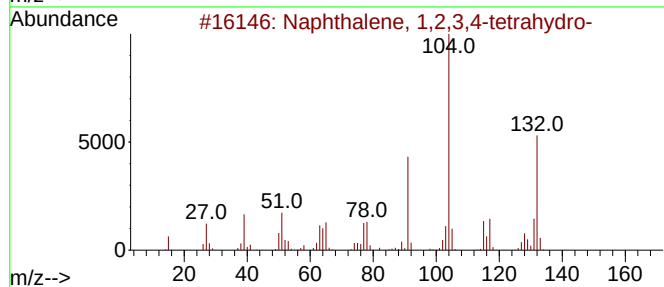
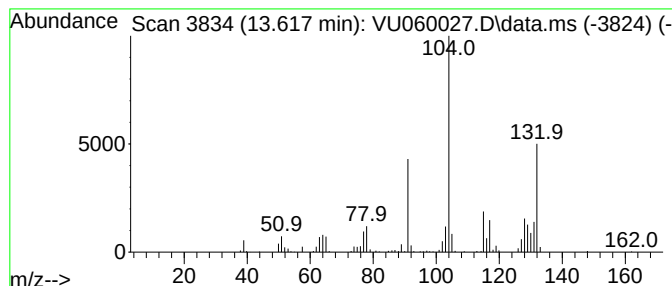
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 14

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.617 | 1.65 ug/L | 239146 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------------------|-----|---------|-------------|------|
| 1    |      | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 91   |
| 2    |      | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 91   |
| 3    |      | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 91   |
| 4    |      | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 90   |
| 5    |      | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

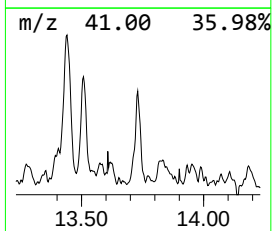
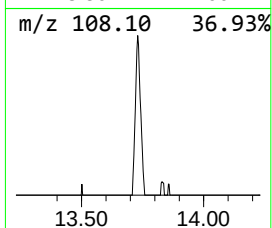
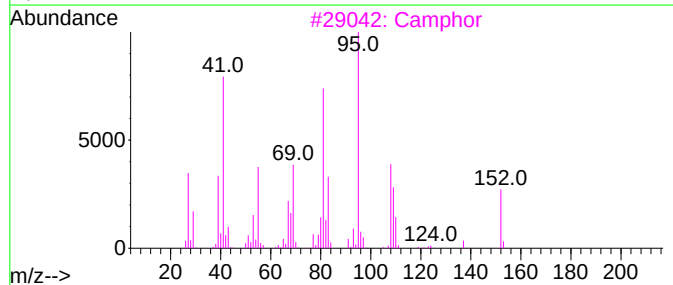
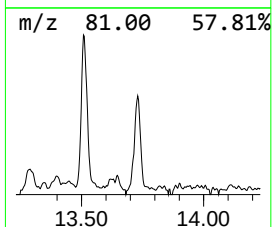
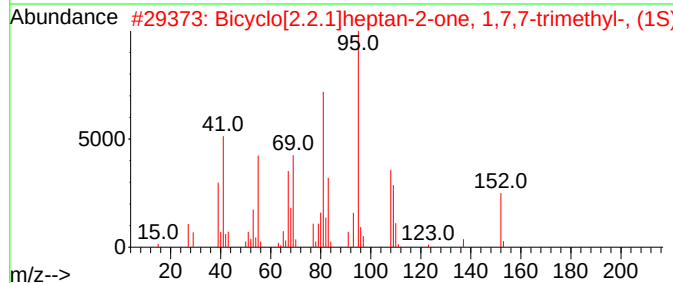
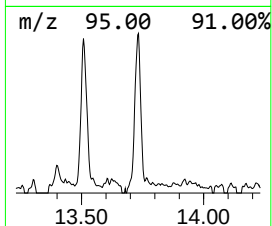
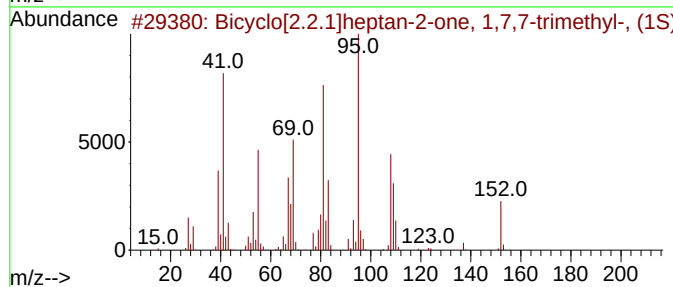
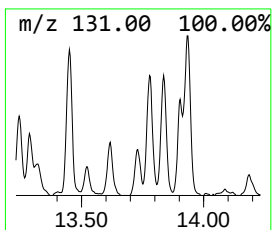
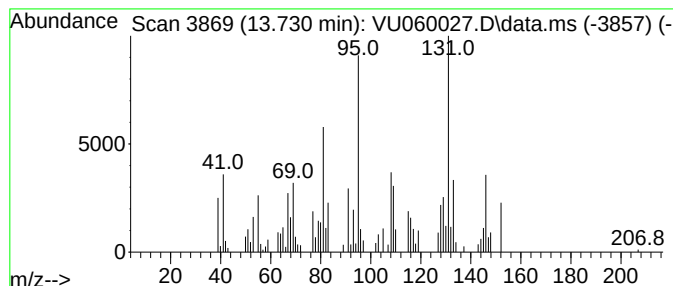
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 31

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 13.730 | 0.54 ug/L | 78265 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 000464-48-2 | 95   |
| 2    |    |   | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 000464-48-2 | 95   |
| 3    |    |   | Camphor                             | 152 | C10H16O | 000076-22-2 | 86   |
| 4    |    |   | (+)-2-Bornanone                     | 152 | C10H16O | 000464-49-3 | 84   |
| 5    |    |   | (+)-2-Bornanone                     | 152 | C10H16O | 000464-49-3 | 83   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

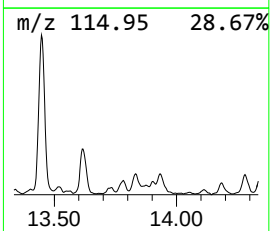
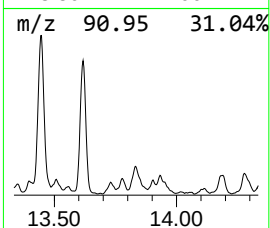
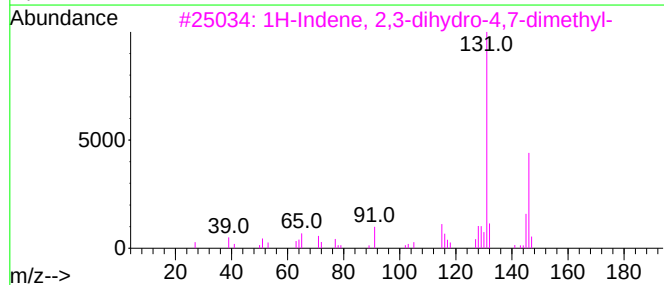
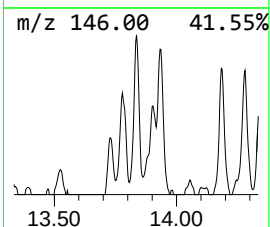
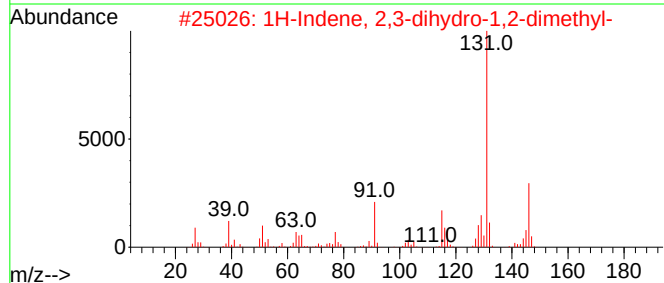
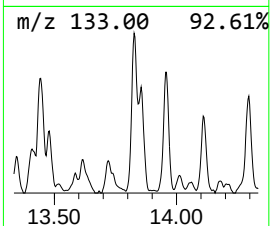
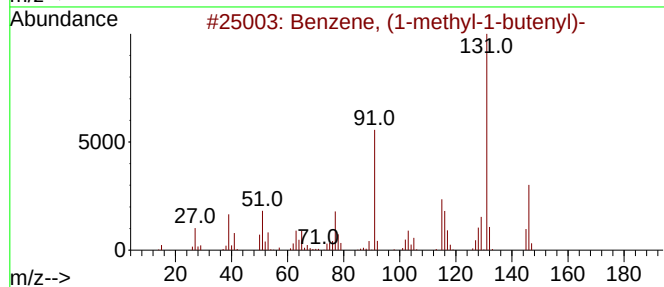
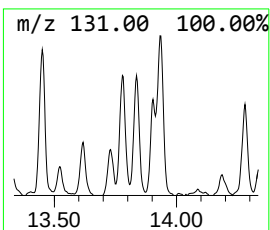
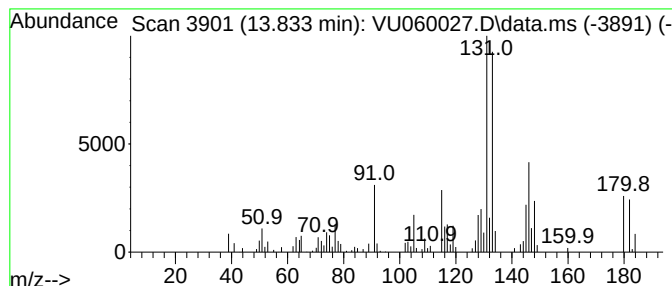
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 21 unknown-02 Concentration Rank 22

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 13.833 | 0.86 ug/L | 124331 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 46   |
| 2    |    |   | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 46   |
| 3    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 45   |
| 4    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 45   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-5,6-dimet... | 146 | C11H14  | 001075-22-5 | 45   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

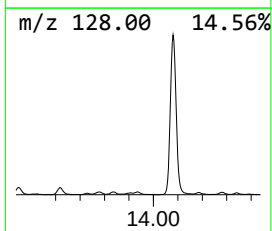
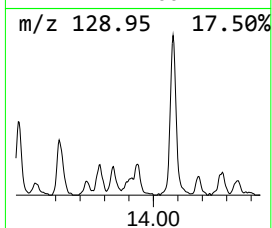
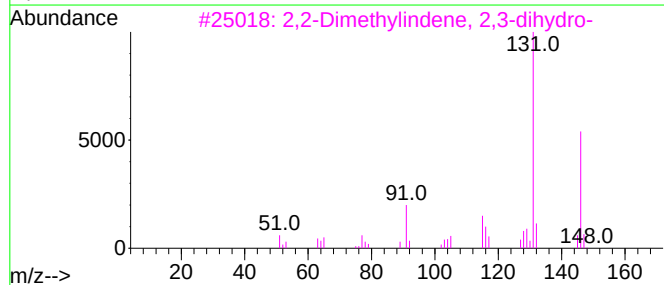
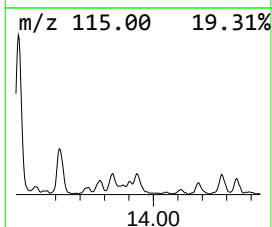
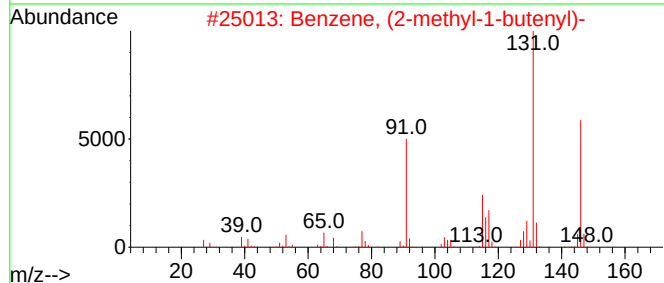
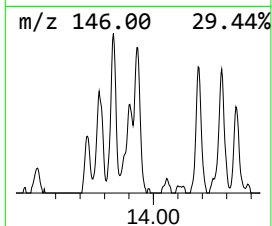
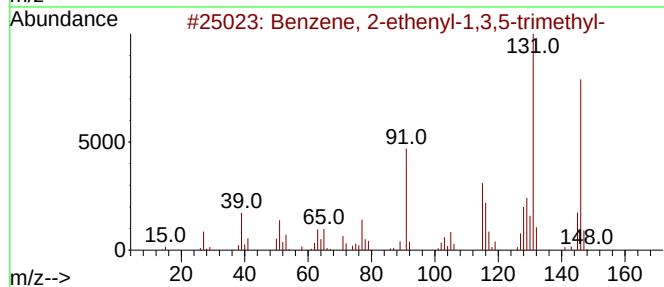
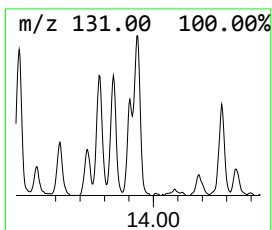
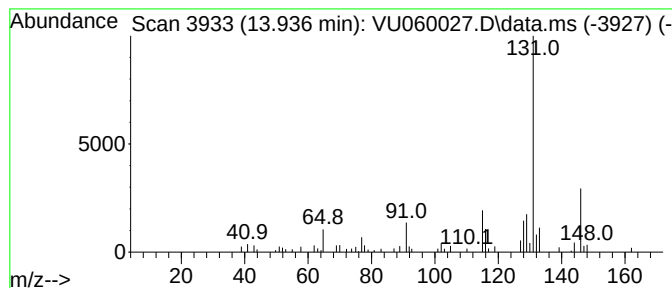
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 22 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 26

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 13.936 | 0.63 ug/L | 91326 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 | C11H14  | 000769-25-5 | 87   |
| 2    |    |   | Benzene, (2-methyl-1-butenyl)-      | 146 | C11H14  | 056253-64-6 | 83   |
| 3    |    |   | 2,2-Dimethylindene, 2,3-dihydro-    | 146 | C11H14  | 020836-11-7 | 83   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14  | 017059-48-2 | 81   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 80   |





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

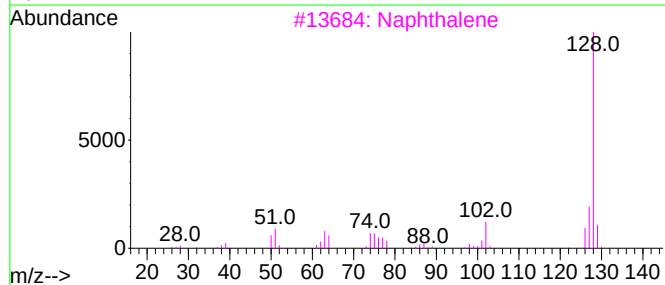
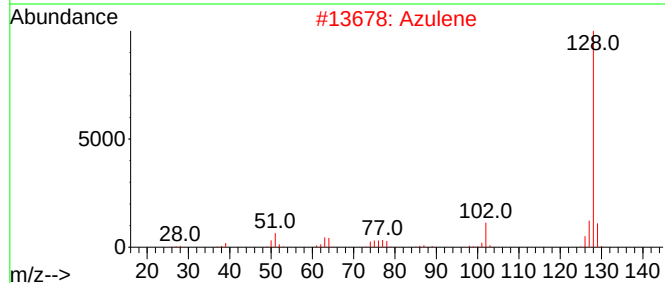
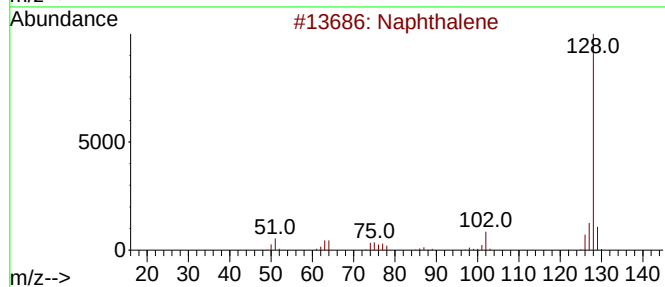
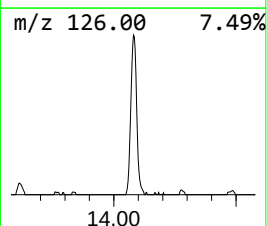
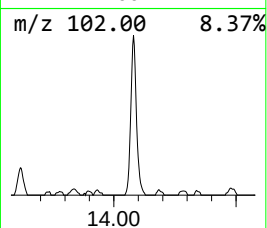
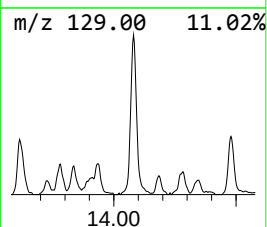
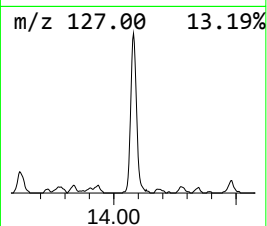
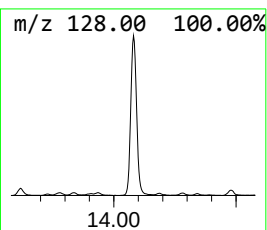
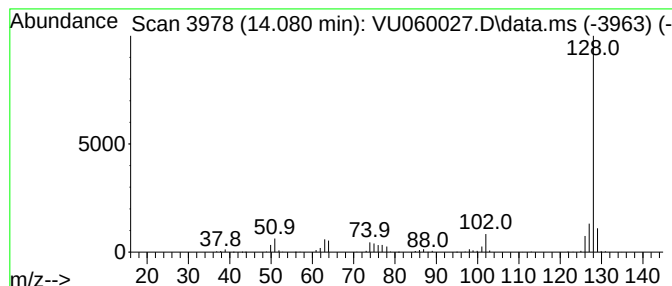
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 23 Azulene Concentration Rank 5

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 14.080 | 2.79 ug/L | 405532 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene  | 128 | C10H8   | 000091-20-3 | 95   |
| 2    |    |   | Azulene      | 128 | C10H8   | 000275-51-4 | 94   |
| 3    |    |   | Naphthalene  | 128 | C10H8   | 000091-20-3 | 94   |
| 4    |    |   | Azulene      | 128 | C10H8   | 000275-51-4 | 94   |
| 5    |    |   | Azulene      | 128 | C10H8   | 000275-51-4 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

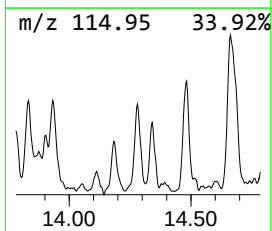
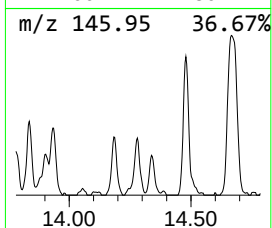
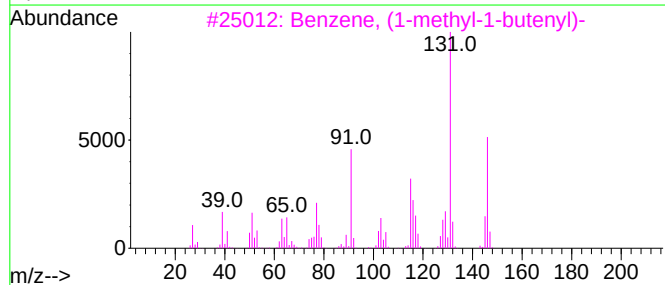
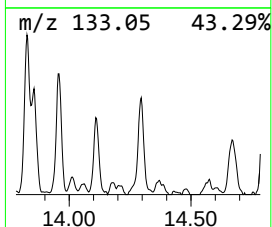
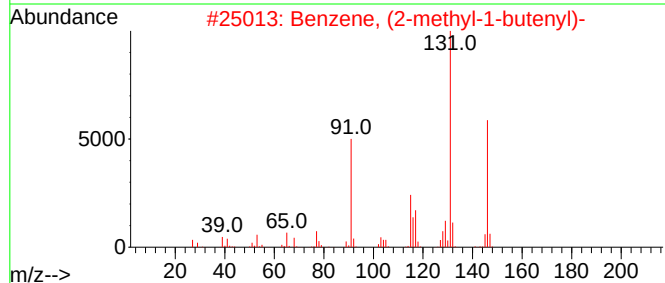
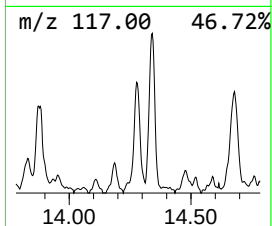
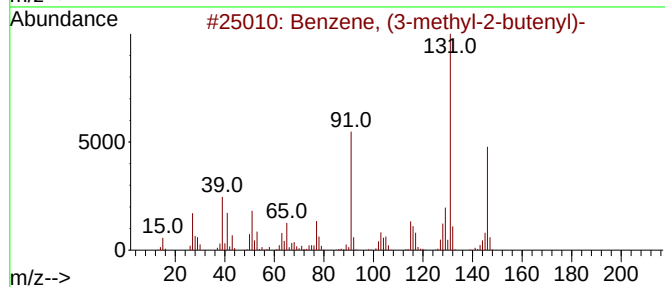
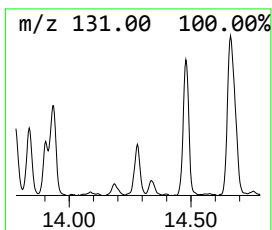
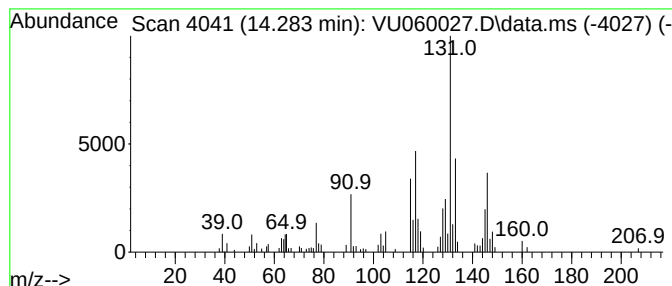
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Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 24 Benzene, (3-methyl-2-butenyl)- Concentration Rank 25

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 14.283 | 0.63 ug/L | 91483 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 55   |
| 2    |    |   | Benzene, (2-methyl-1-butenyl)-      | 146 | C11H14  | 056253-64-6 | 50   |
| 3    |    |   | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 49   |
| 4    |    |   | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 | C11H14  | 000769-25-5 | 49   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 46   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

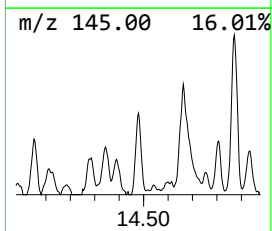
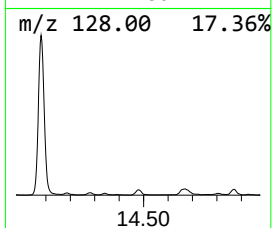
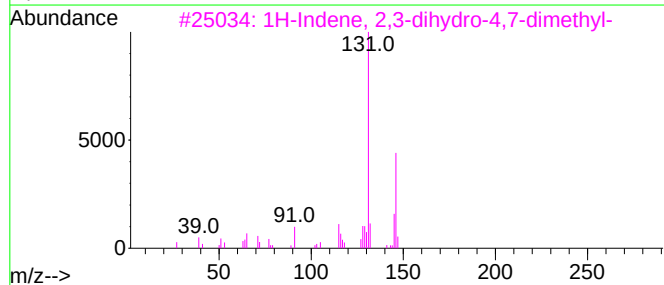
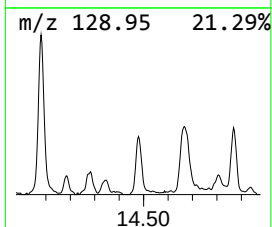
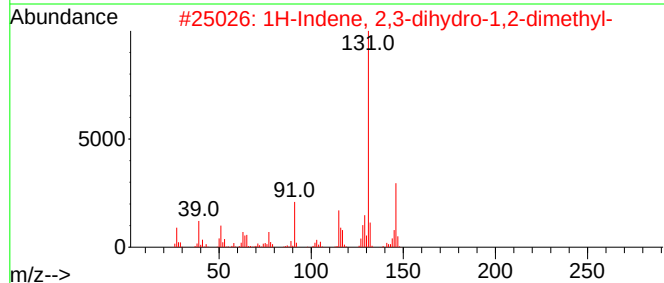
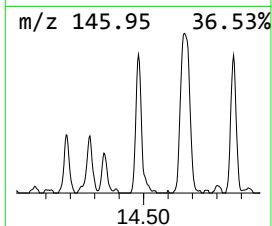
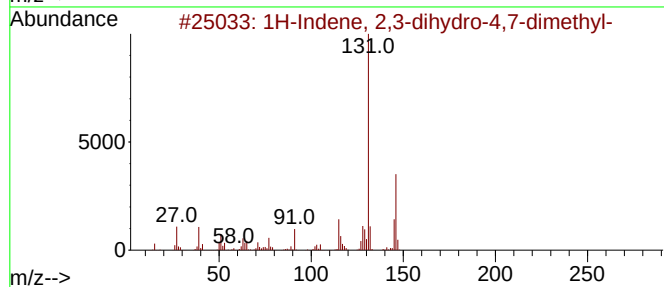
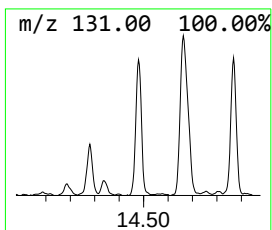
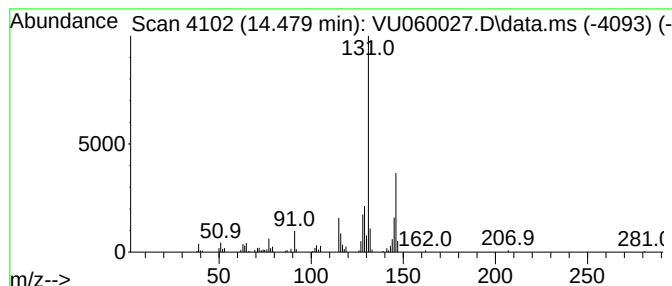
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 25 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 23

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 14.479 | 0.85 ug/L | 122752 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 94   |
| 2    |    |   | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 94   |
| 3    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 93   |
| 4    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 91   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 91   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

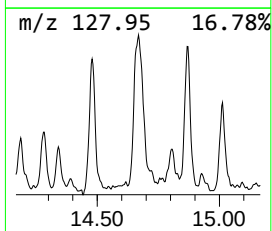
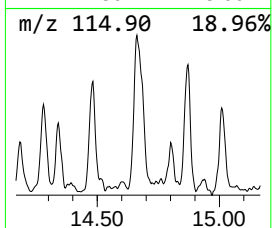
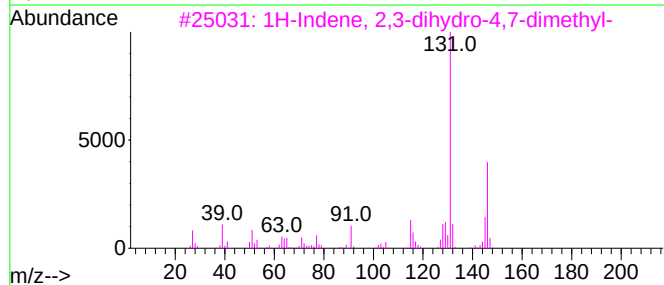
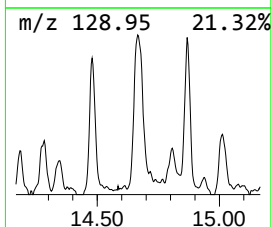
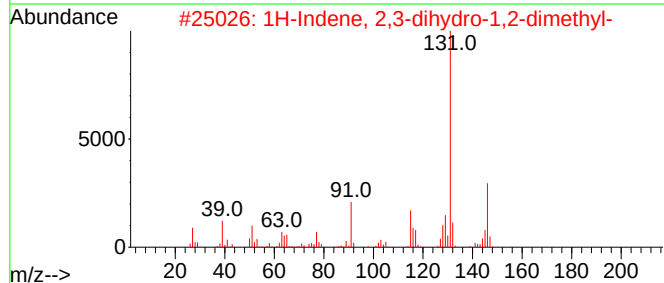
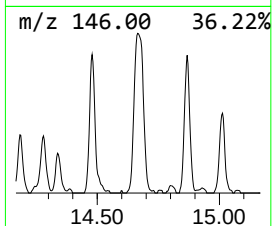
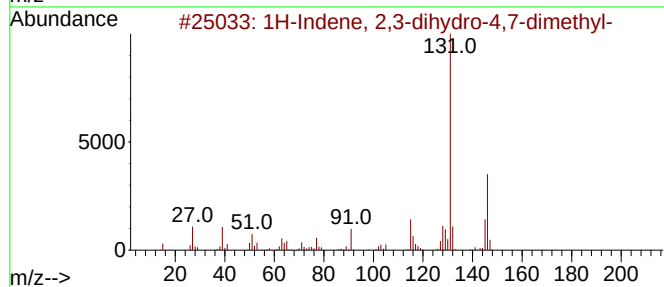
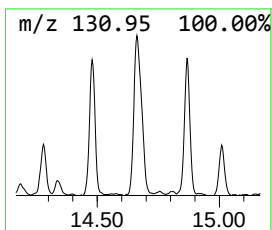
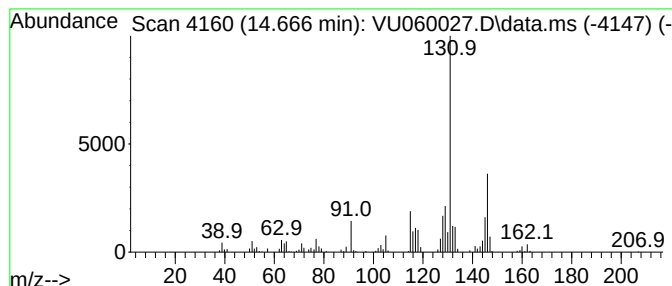
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 12

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 14.666 | 1.76 ug/L | 256245 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 94   |
| 2    |    |   | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 93   |
| 3    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 89   |
| 4    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 87   |
| 5    |    |   | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 87   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

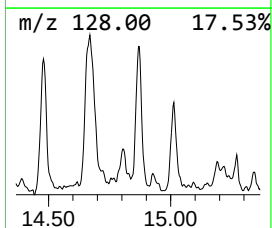
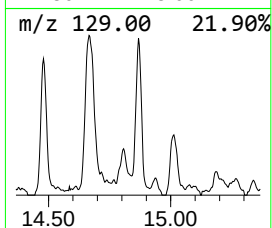
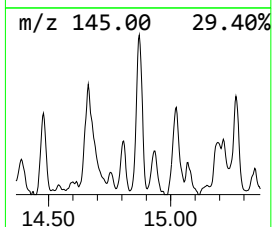
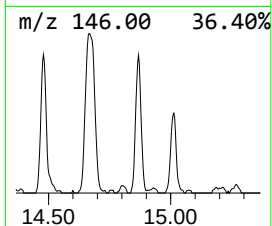
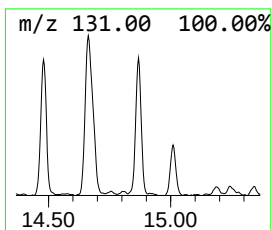
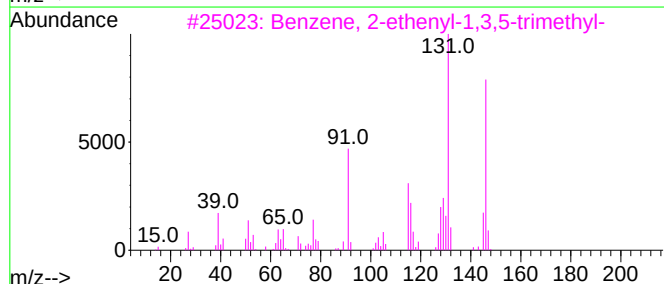
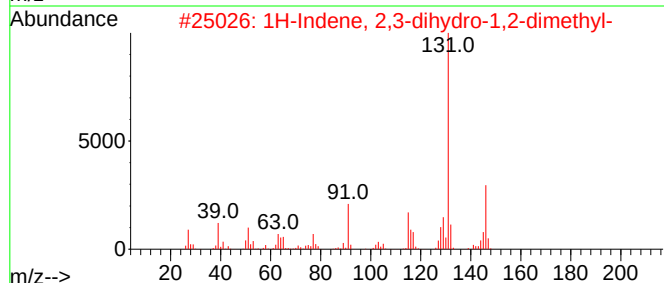
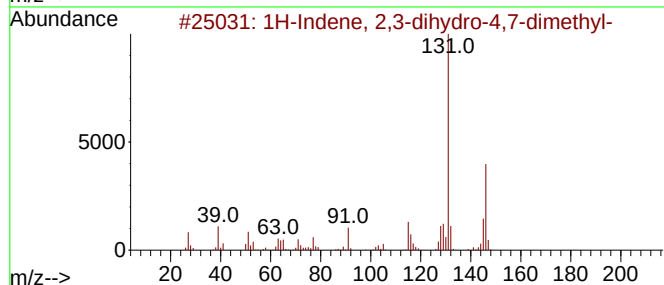
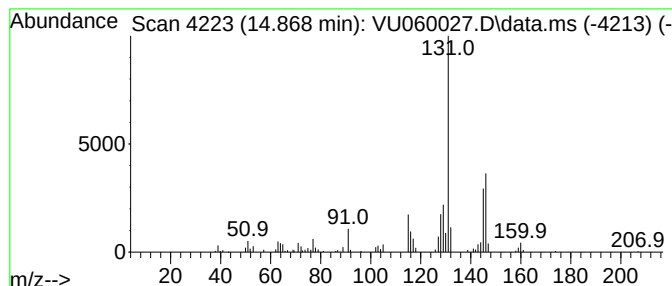
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 27 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 19

| R.T.   | EstConc   | Area   | Relative to ISTD       | R.T.   |
|--------|-----------|--------|------------------------|--------|
| 14.868 | 0.93 ug/L | 134940 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 87   |
| 2    |    |   | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 83   |
| 3    |    |   | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 | C11H14  | 000769-25-5 | 81   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 81   |
| 5    |    |   | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 81   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

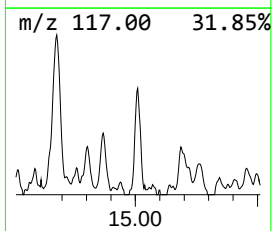
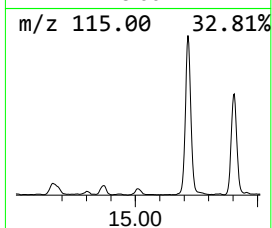
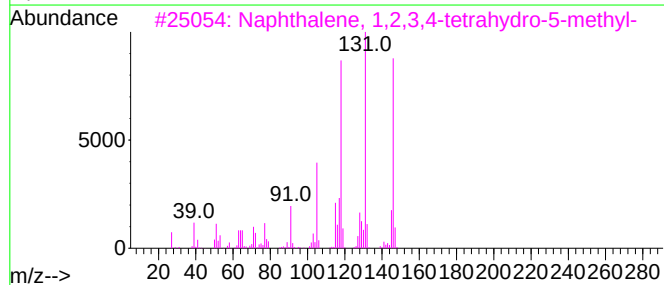
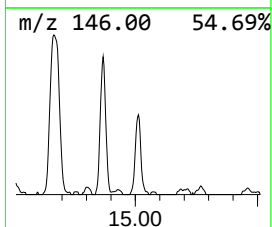
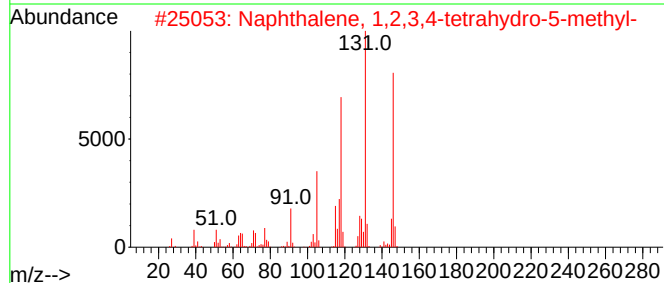
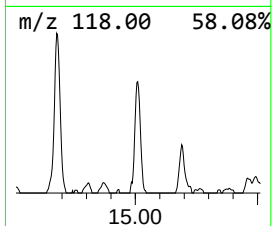
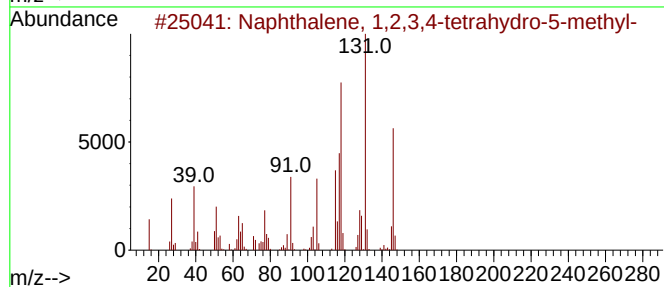
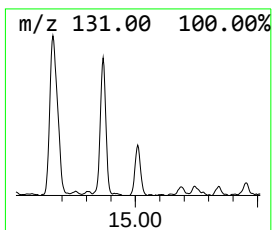
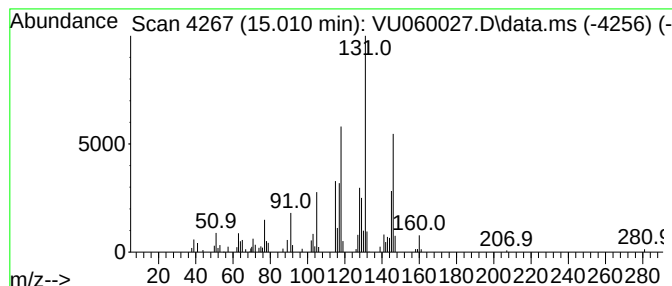
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 28 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 27

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 15.010 | 0.63 ug/L | 91051 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5 | 90   |
| 2    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5 | 81   |
| 3    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5 | 81   |
| 4    |    |   | 2-Propyn-1-ol, 3-(4-methylphenyl)-  | 146 | C10H10O | 016017-24-6 | 68   |
| 5    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 64   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

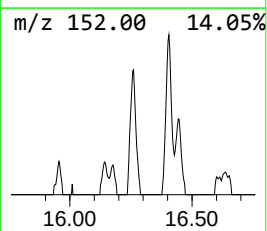
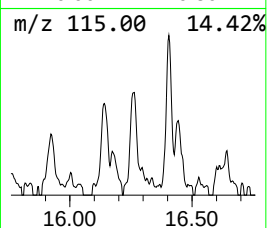
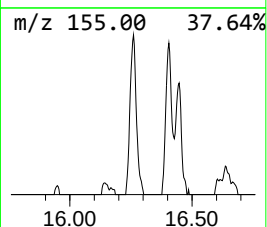
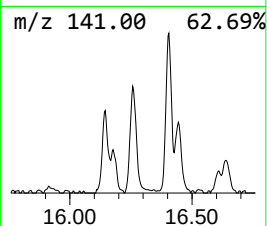
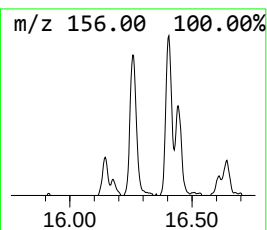
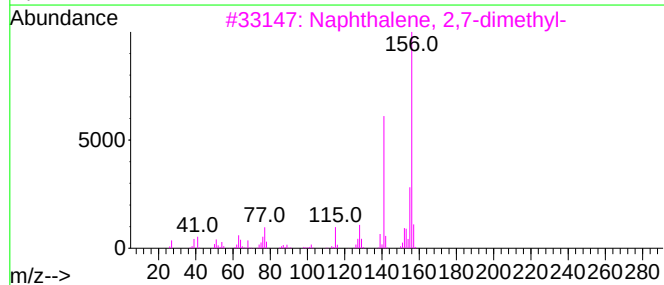
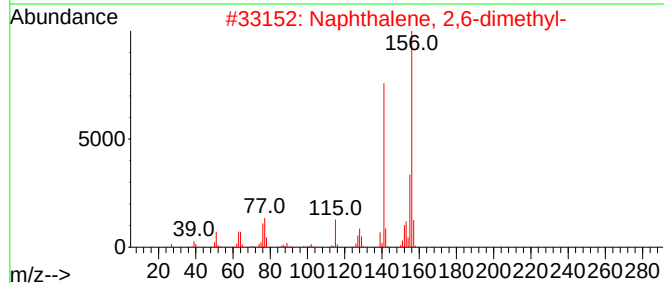
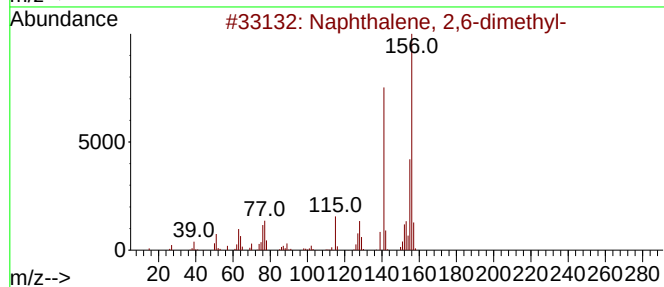
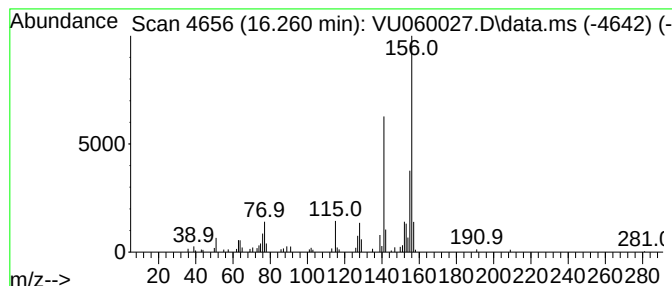
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.M  
Quant Title : TRACE VOA SFAM1.0

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

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Peak Number 31 Naphthalene, 2,6-dimethyl- Concentration Rank 29

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 16.260 | 0.55 ug/L | 80342 | 1,4-Dichlorobenzene-d4 | 11.807 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 2    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 3    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 4    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |
| 5    |    |   | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 97   |



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

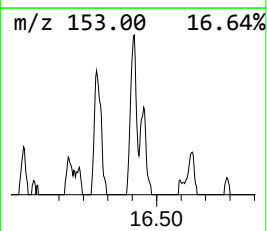
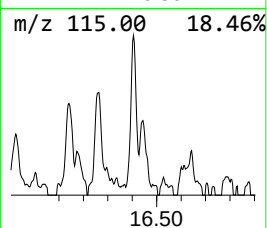
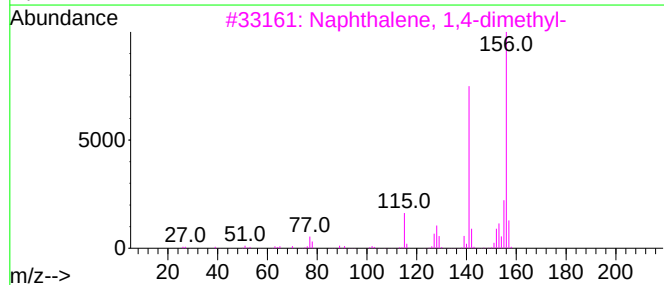
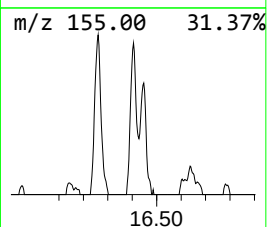
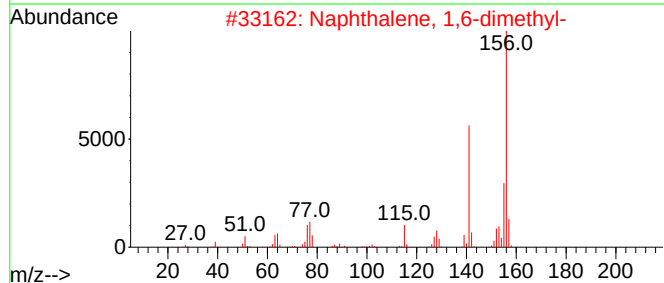
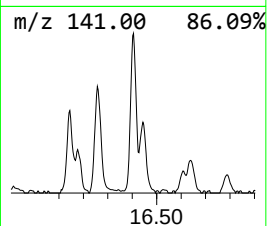
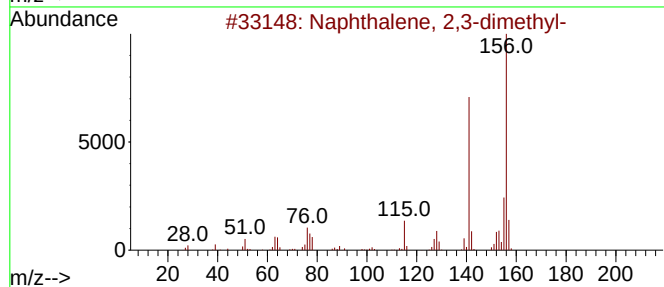
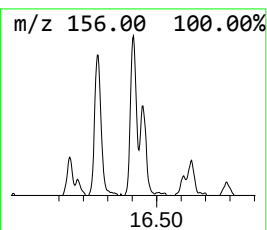
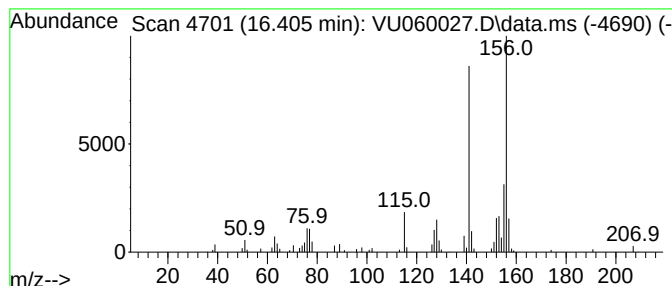
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Quant Title : TRACE VOA SFAM1.0

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

\*\*\*\*\*  
Peak Number 32 Naphthalene, 2,3-dimethyl- Concentration Rank 30

| R.T.   | EstConc   | Area  | Relative to ISTD       | R.T.   |
|--------|-----------|-------|------------------------|--------|
| 16.405 | 0.55 ug/L | 79887 | 1,4-Dichlorobenzene-d4 | 11.807 |

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU071824\  
Data File : VU060027.D  
Acq On : 18 Jul 2024 19:11  
Operator : MD/SY  
Sample : P3217-03DL 100X  
Misc : 25.0mL/MSVOA\_U/WATER  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
MSVOA\_U  
ClientSampleId :  
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR071824WMA.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, 1-ethy... | 10.991 | 1.8     | ug/L  | 260950   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 1-ethy... | 11.286 | 2.4     | ug/L  | 345527   | 3                     | 11.807 | 725980 | 5.0  |
| 7-Oxabicyclo[2...  | 11.634 | 0.6     | ug/L  | 88734    | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 1,2,3-... | 11.888 | 6.1     | ug/L  | 890725   | 3                     | 11.807 | 725980 | 5.0  |
| Indane             | 12.090 | 2.2     | ug/L  | 323573   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 1-meth... | 12.370 | 0.9     | ug/L  | 134868   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 2-ethy... | 12.460 | 1.0     | ug/L  | 147540   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 1,2,4,... | 12.489 | 0.9     | ug/L  | 137511   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 1-ethy... | 12.566 | 2.0     | ug/L  | 282649   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, (1-met... | 12.605 | 0.5     | ug/L  | 75465    | 3                     | 11.807 | 725980 | 5.0  |
| Indan, 1-methyl-   | 12.675 | 2.1     | ug/L  | 310950   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 1-ethy... | 12.868 | 1.1     | ug/L  | 155931   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 1,2,3,... | 12.965 | 1.7     | ug/L  | 249894   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 1,2,3,... | 13.019 | 2.3     | ug/L  | 335699   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, (2-met... | 13.286 | 0.9     | ug/L  | 125967   | 3                     | 11.807 | 725980 | 5.0  |
| 2,4-Dimethylsty... | 13.447 | 4.3     | ug/L  | 618794   | 3                     | 11.807 | 725980 | 5.0  |
| unknown-01         | 13.511 | 0.6     | ug/L  | 92280    | 3                     | 11.807 | 725980 | 5.0  |
| Naphthalene, 1,... | 13.617 | 1.6     | ug/L  | 239146   | 3                     | 11.807 | 725980 | 5.0  |
| Bicyclo[2.2.1]h... | 13.730 | 0.5     | ug/L  | 78265    | 3                     | 11.807 | 725980 | 5.0  |
| unknown-02         | 13.833 | 0.9     | ug/L  | 124331   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 2-ethe... | 13.936 | 0.6     | ug/L  | 91326    | 3                     | 11.807 | 725980 | 5.0  |
| Azulene            | 14.080 | 2.8     | ug/L  | 405532   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, (3-met... | 14.283 | 0.6     | ug/L  | 91483    | 3                     | 11.807 | 725980 | 5.0  |
| 1H-Indene, 2,3-... | 14.479 | 0.8     | ug/L  | 122752   | 3                     | 11.807 | 725980 | 5.0  |
| 1H-Indene, 2,3-... | 14.666 | 1.8     | ug/L  | 256245   | 3                     | 11.807 | 725980 | 5.0  |
| Benzene, 2-ethe... | 14.868 | 0.9     | ug/L  | 134940   | 3                     | 11.807 | 725980 | 5.0  |
| Naphthalene, 1,... | 15.010 | 0.6     | ug/L  | 91051    | 3                     | 11.807 | 725980 | 5.0  |
| Naphthalene, 2,... | 16.260 | 0.6     | ug/L  | 80342    | 3                     | 11.807 | 725980 | 5.0  |
| Naphthalene, 2,... | 16.405 | 0.6     | ug/L  | 79887    | 3                     | 11.807 | 725980 | 5.0  |