

Data Path : Z:\VOASRV\HPCHEM1\MSVOA\_W\DATA\VW121718\  
 Data File : VW007593.D  
 Acq On : 15 Dec 2018 20:02  
 Operator : SY/AP  
 Sample : J6378-07  
 Misc : 2.70G/5ML/MSVOA\_W/SOIL  
 ALS Vial : 79 Sample Multiplier: 1

Instrument :  
 MSVOA\_W  
 ClientSampleId :  
 P001-SD010-0612-01

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA\_W\METHOD\82W121118S.M

Title : SW846 8260

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.952       | 474           | 500         | 525          | rBV      | 466595         | 1802515       | 19.51%          | 7.320%        |
| 2         | 5.434       | 561           | 579         | 604          | rBV      | 555439         | 1986982       | 21.50%          | 8.069%        |
| 3         | 5.934       | 646           | 661         | 680          | rBV      | 1097664        | 3349985       | 36.25%          | 13.604%       |
| 4         | 6.982       | 820           | 833         | 850          | rVB      | 750824         | 2271889       | 24.59%          | 9.226%        |
| 5         | 7.952       | 970           | 992         | 1000         | rBV3     | 164343         | 551284        | 5.97%           | 2.239%        |
| 6         | 8.153       | 1016          | 1025        | 1032         | rBV      | 62438          | 143813        | 1.56%           | 0.584%        |
| 7         | 8.433       | 1061          | 1071        | 1077         | rVV3     | 46792          | 113172        | 1.22%           | 0.460%        |
| 8         | 8.506       | 1077          | 1083        | 1088         | rVV      | 44183          | 98976         | 1.07%           | 0.402%        |
| 9         | 8.573       | 1088          | 1094        | 1106         | rVV2     | 70782          | 159230        | 1.72%           | 0.647%        |
| 10        | 8.842       | 1130          | 1138        | 1147         | rBB      | 52253          | 104143        | 1.13%           | 0.423%        |
| 11        | 9.335       | 1206          | 1219        | 1233         | rBB      | 288163         | 603627        | 6.53%           | 2.451%        |
| 12        | 10.323      | 1373          | 1381        | 1396         | rBV      | 93243          | 186935        | 2.02%           | 0.759%        |
| 13        | 11.658      | 1588          | 1600        | 1607         | rBV2     | 374089         | 703539        | 7.61%           | 2.857%        |
| 14        | 11.835      | 1621          | 1629        | 1648         | rVB      | 5580566        | 9240922       | 100.00%         | 37.527%       |
| 15        | 12.615      | 1750          | 1757        | 1766         | rVB      | 285381         | 484696        | 5.25%           | 1.968%        |
| 16        | 12.871      | 1793          | 1799        | 1804         | rBV      | 151064         | 237457        | 2.57%           | 0.964%        |
| 17        | 12.945      | 1804          | 1811        | 1818         | rVV      | 126629         | 200390        | 2.17%           | 0.814%        |
| 18        | 13.255      | 1856          | 1862        | 1868         | rBV      | 379963         | 582037        | 6.30%           | 2.364%        |
| 19        | 13.499      | 1897          | 1902        | 1908         | rVV      | 109732         | 172027        | 1.86%           | 0.699%        |
| 20        | 13.585      | 1908          | 1916        | 1926         | rVB2     | 873434         | 1538992       | 16.65%          | 6.250%        |
| 21        | 13.792      | 1947          | 1950        | 1959         | rVB3     | 46978          | 92445         | 1.00%           | 0.375%        |

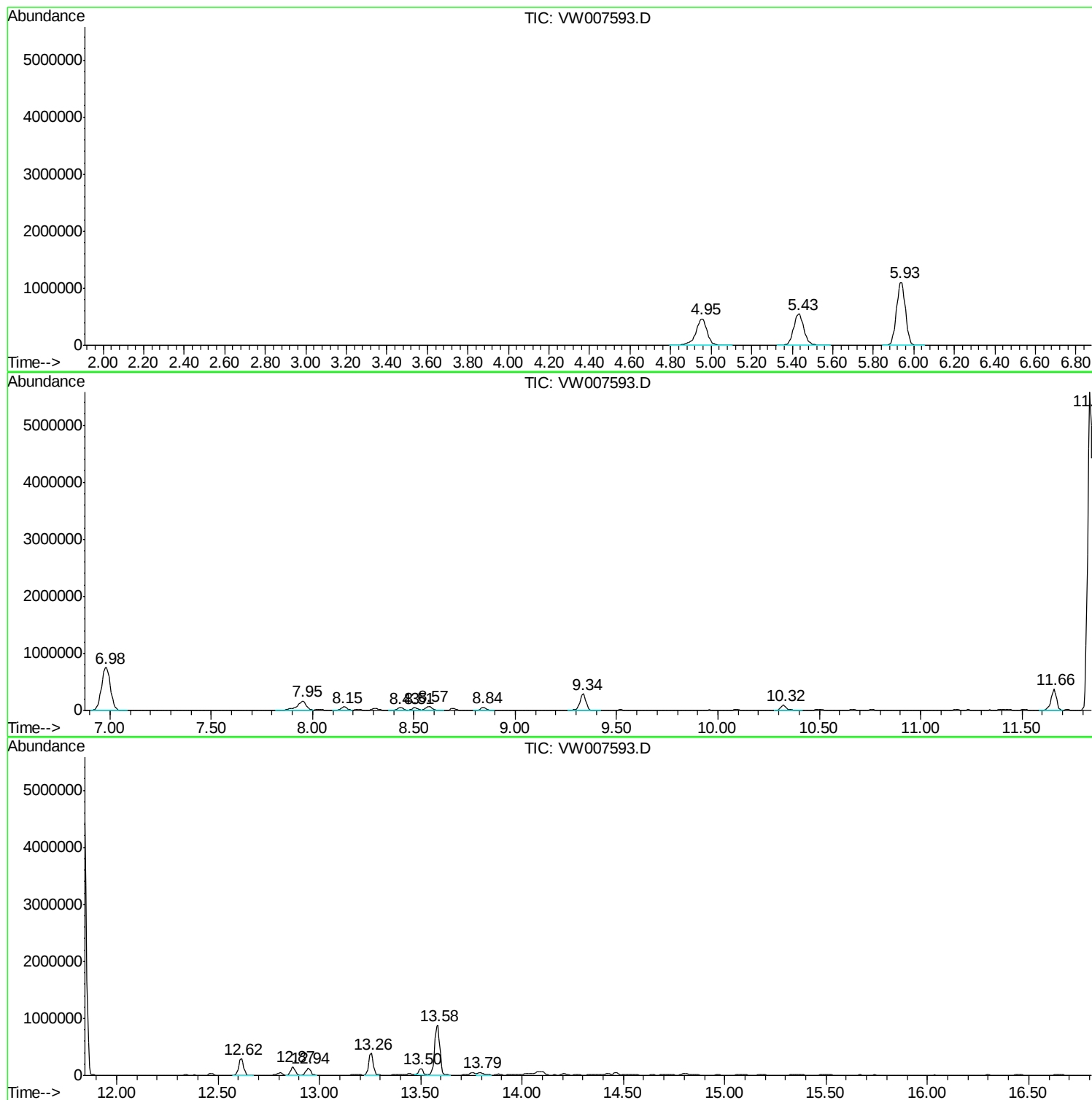
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ALS Vial : 79 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
P001-SD010-0612-01

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_W\METHOD\82W121118S.M  
Quant Title : SW846 8260

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



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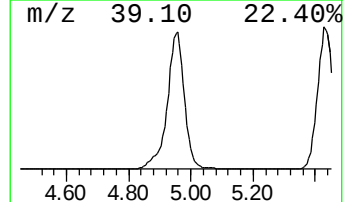
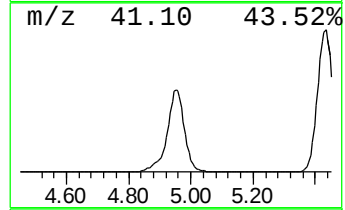
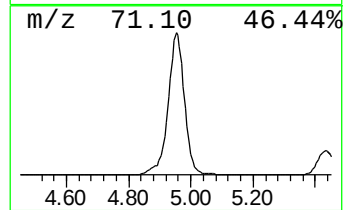
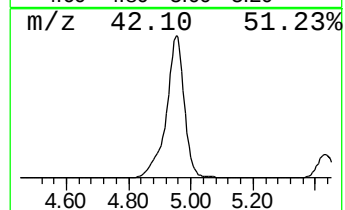
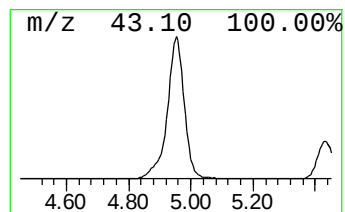
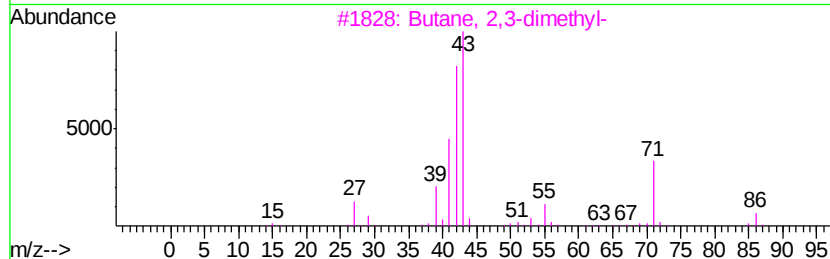
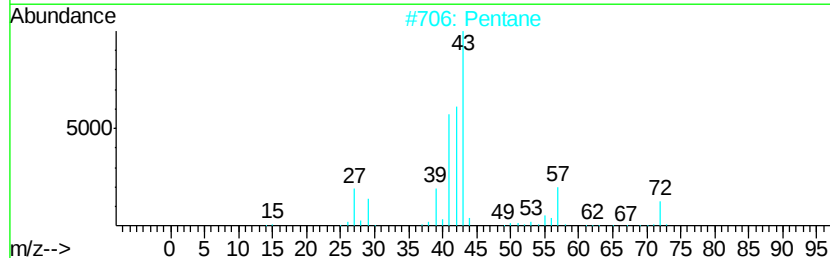
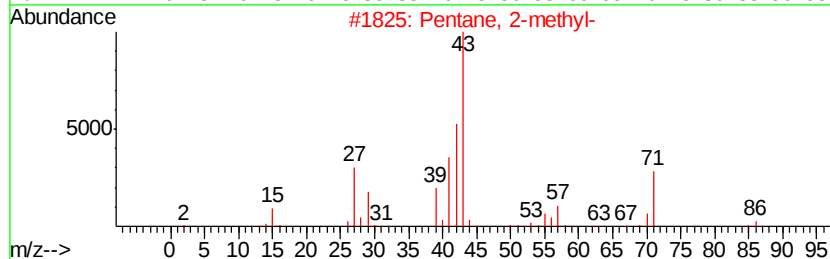
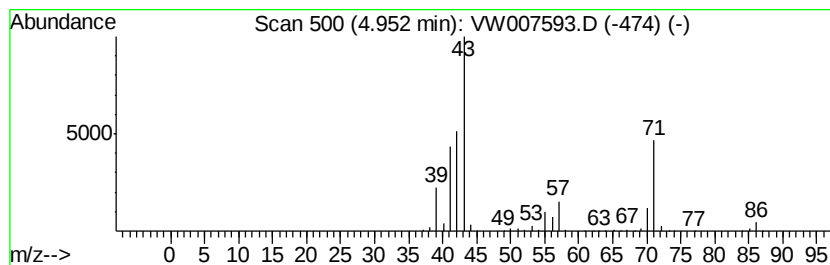
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_W\METHOD\82W121118S.M  
Quant Title : SW846 8260

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

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Peak Number 1 Pentane, 2-methyl- Concentration Rank 4

| R.T. | EstConc     | Area    | Relative to ISTD   | R.T. |
|------|-------------|---------|--------------------|------|
| 4.95 | 163.48 ug/l | 1802520 | Pentafluorobenzene | 7.95 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentane, 2-methyl-        | 86  | C6H14   | 000107-83-5 | 91   |
| 2    |    | Pentane                   | 72  | C5H12   | 000109-66-0 | 49   |
| 3    |    | Butane, 2,3-dimethyl-     | 86  | C6H14   | 000079-29-8 | 49   |
| 4    |    | 1-Butanol, 2,3-dimethyl-  | 102 | C6H14O  | 019550-30-2 | 42   |
| 5    |    | Pentane, 2,3,4-trimethyl- | 114 | C8H18   | 000565-75-3 | 37   |



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Instrument :  
MSVOA\_W  
ClientSampleId :  
P001-SD010-0612-01

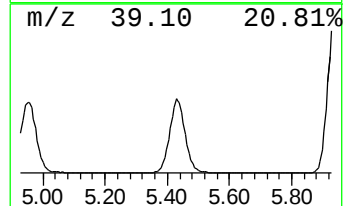
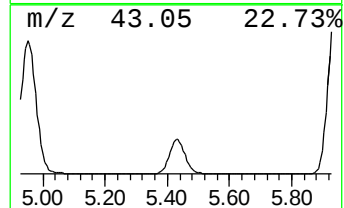
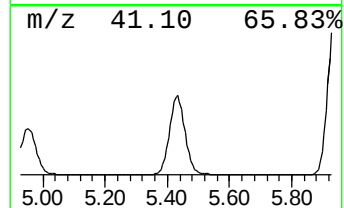
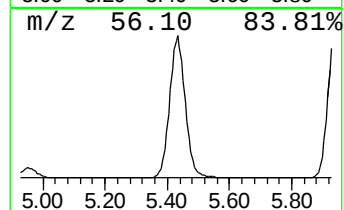
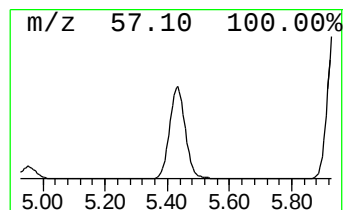
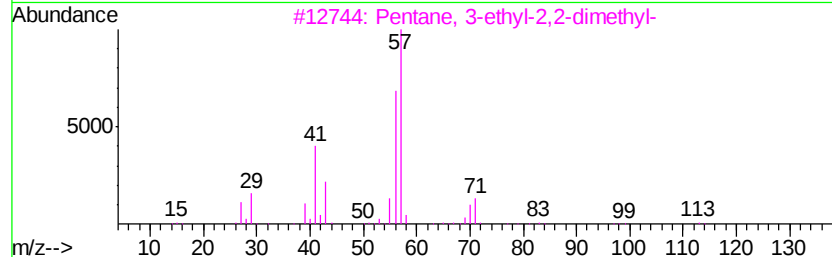
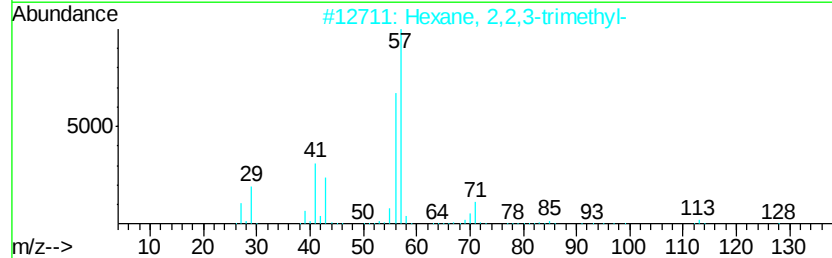
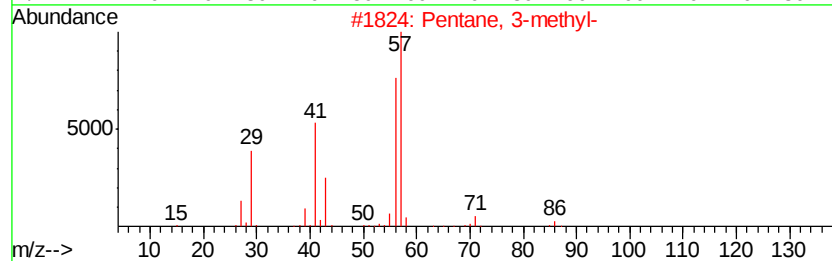
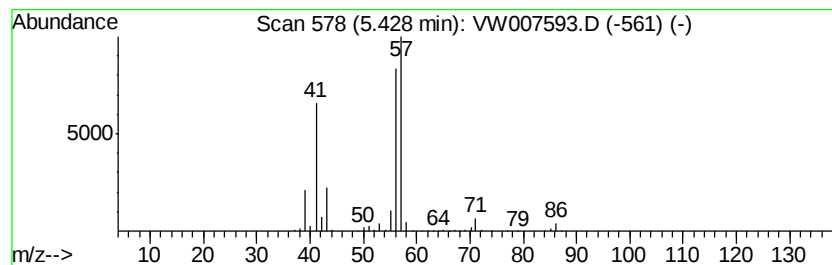
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_W\METHOD\82W121118S.M  
Quant Title : SW846 8260

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

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Peak Number 2 Pentane, 3-methyl- Concentration Rank 3

| R.T. | EstConc     | Area    | Relative to ISTD   | R.T. |
|------|-------------|---------|--------------------|------|
| 5.43 | 180.21 ug/l | 1986980 | Pentafluorobenzene | 7.95 |

| Hit# | of | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Pentane, 3-methyl-                 | 86  | C6H14    | 000096-14-0 | 90   |
| 2    |    | Hexane, 2,2,3-trimethyl-           | 128 | C9H20    | 016747-25-4 | 64   |
| 3    |    | Pentane, 3-ethyl-2,2-dimethyl-     | 128 | C9H20    | 016747-32-3 | 43   |
| 4    |    | 3,4-Dimethyldihydrofuran-2,5-dione | 128 | C6H8O3   | 007475-92-5 | 40   |
| 5    |    | Sulphuric acid dibutyl ester       | 210 | C8H18O4S | 000625-22-9 | 40   |



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MSVOA\_W  
ClientSampleId :  
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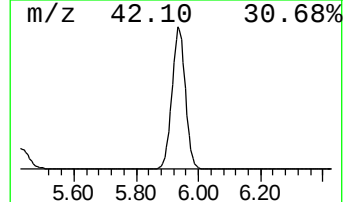
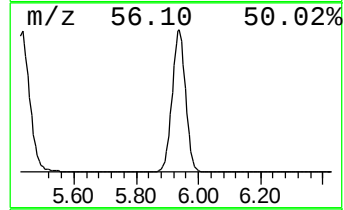
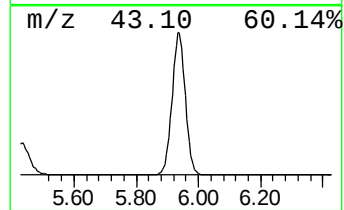
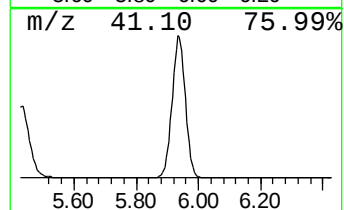
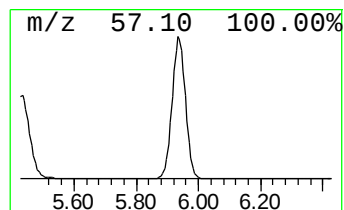
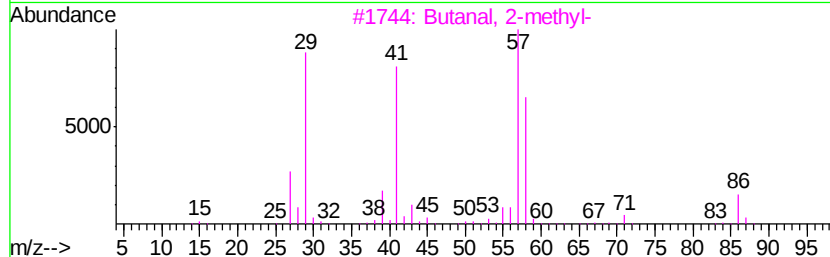
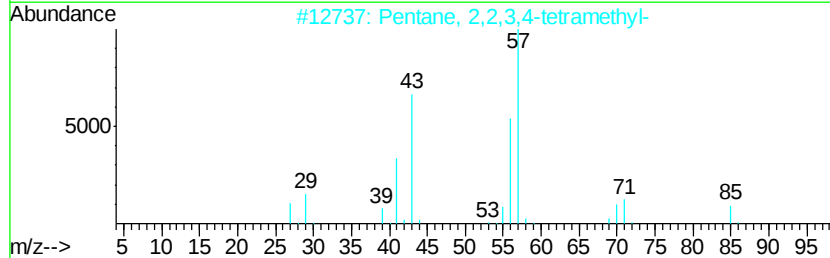
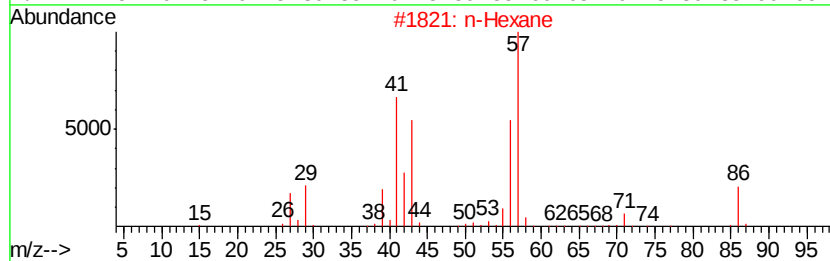
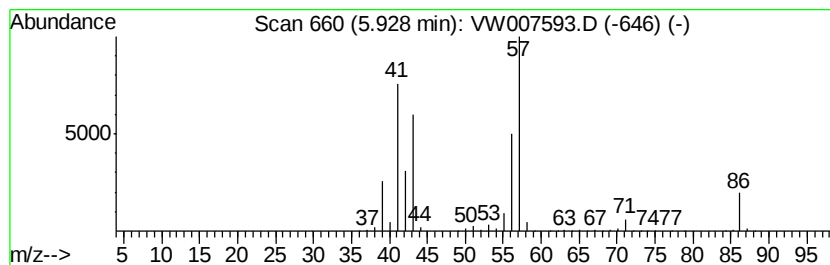
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_W\METHOD\82W121118S.M  
Quant Title : SW846 8260

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

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Peak Number 3 n-Hexane Concentration Rank 1

| R.T. | EstConc     | Area    | Relative to ISTD   | R.T. |
|------|-------------|---------|--------------------|------|
| 5.93 | 303.84 ug/l | 3349990 | Pentafluorobenzene | 7.95 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | n-Hexane                      | 86  | C6H14   | 000110-54-3 | 94   |
| 2    |    | Pentane, 2,2,3,4-tetramethyl- | 128 | C9H20   | 001186-53-4 | 38   |
| 3    |    | Butanal, 2-methyl-            | 86  | C5H10O  | 000096-17-3 | 38   |
| 4    |    | Propanal, 2,2-dimethyl-       | 86  | C5H10O  | 000630-19-3 | 38   |
| 5    |    | 1-Pentene, 4-methyl-          | 84  | C6H12   | 000691-37-2 | 37   |



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Instrument :  
MSVOA\_W  
ClientSampleId :  
P001-SD010-0612-01

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_W\METHOD\82W121118S.M  
Quant Title : SW846 8260

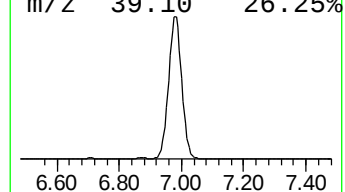
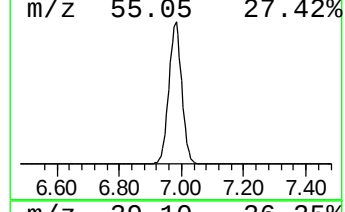
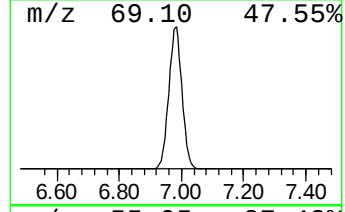
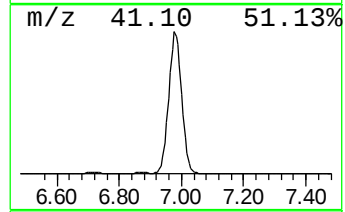
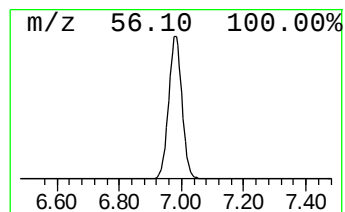
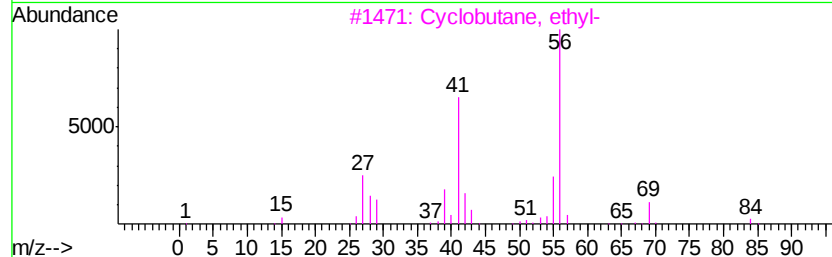
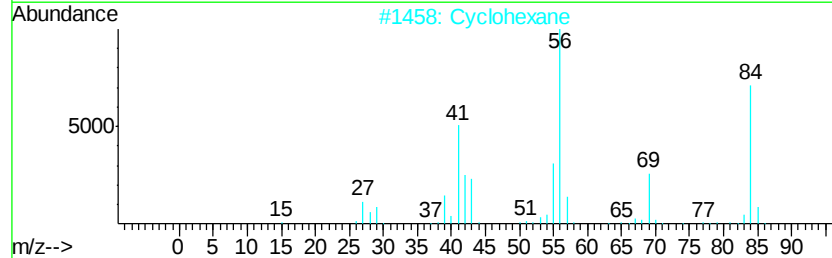
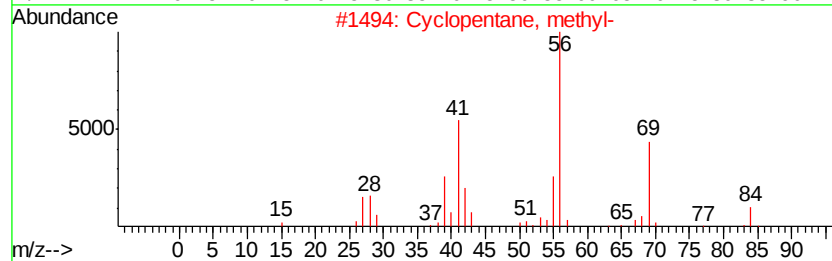
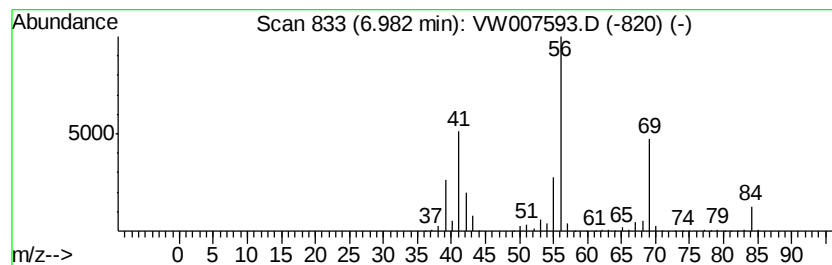
TIC Library : C:\DATABASE\NIST11.L  
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\*\*\*\*\*  
Peak Number 4 Cyclopentane, methyl- Concentration Rank 2

| R.T. | EstConc     | Area    | Relative to ISTD   | R.T. |
|------|-------------|---------|--------------------|------|
| 6.98 | 206.05 ug/l | 2271890 | Pentafluorobenzene | 7.95 |

| Hit# of | 5 | Tentative ID            | MW | MolForm | CAS#        | Qual |
|---------|---|-------------------------|----|---------|-------------|------|
| 1       |   | Cyclopentane, methyl-   | 84 | C6H12   | 000096-37-7 | 94   |
| 2       |   | Cyclohexane             | 84 | C6H12   | 000110-82-7 | 78   |
| 3       |   | Cyclobutane, ethyl-     | 84 | C6H12   | 004806-61-5 | 72   |
| 4       |   | 1H-Tetrazole, 5-methyl- | 84 | C2H4N4  | 004076-36-2 | 72   |
| 5       |   | Cyclobutane             | 56 | C4H8    | 000287-23-0 | 53   |



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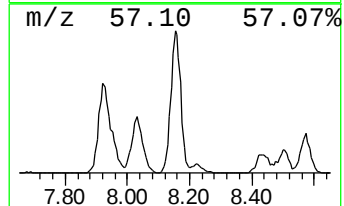
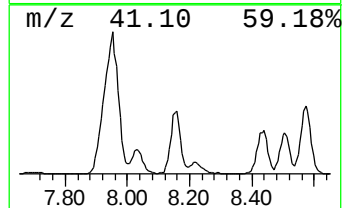
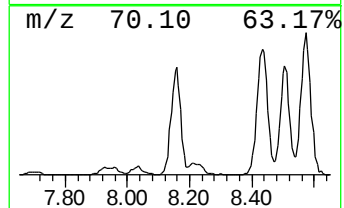
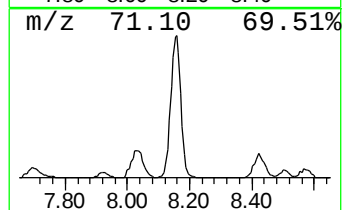
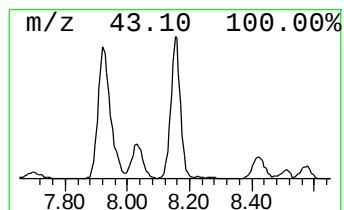
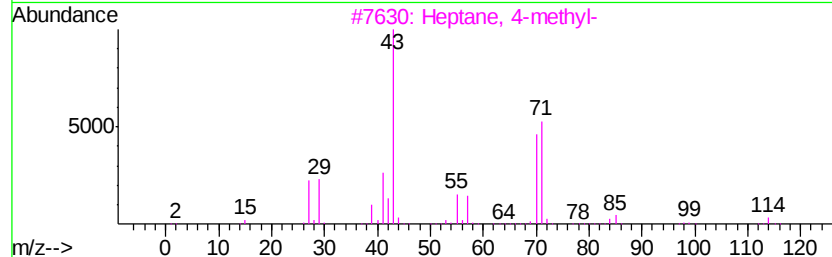
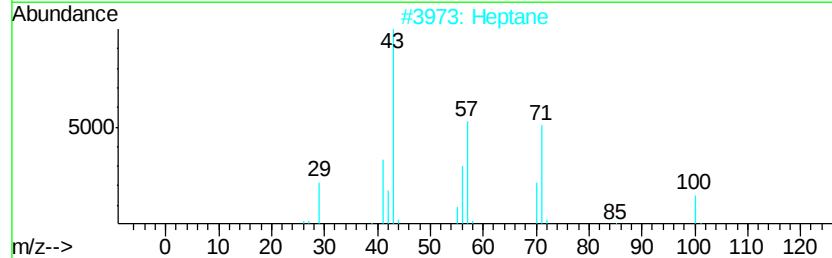
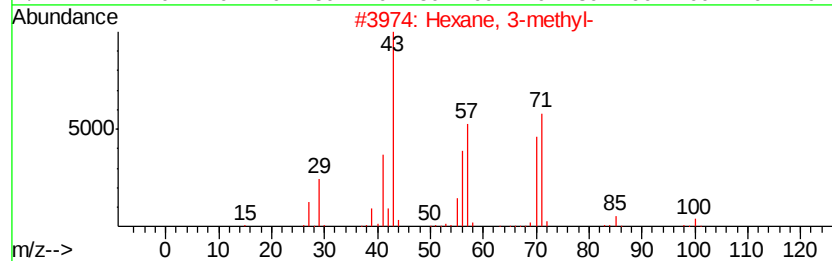
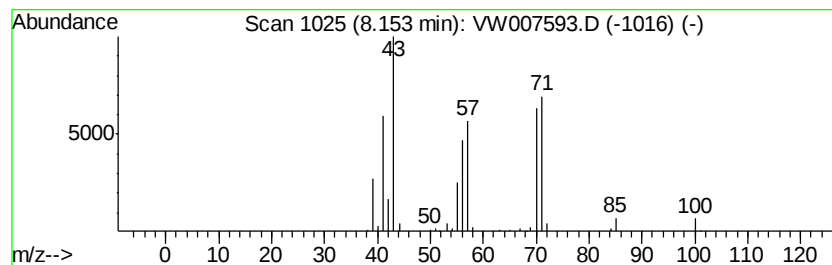
Instrument :  
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Hexane, 3-methyl- Concentration Rank 8

| R.T.    | EstConc                   | Area         | Relative to ISTD   | R.T.        |      |      |
|---------|---------------------------|--------------|--------------------|-------------|------|------|
| 8.15    | 13.04 ug/l                | 143813       | Pentafluorobenzene | 7.95        |      |      |
| Hit# of | 5                         | Tentative ID | MW                 | MolForm     | CAS# | Qual |
| 1       | Hexane, 3-methyl-         | 100          | C7H16              | 000589-34-4 | 91   |      |
| 2       | Heptane                   | 100          | C7H16              | 000142-82-5 | 80   |      |
| 3       | Heptane, 4-methyl-        | 114          | C8H18              | 000589-53-7 | 59   |      |
| 4       | Pentane, 2,3,3-trimethyl- | 114          | C8H18              | 000560-21-4 | 53   |      |
| 5       | Hexane, 2,3-dimethyl-     | 114          | C8H18              | 000584-94-1 | 53   |      |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_W\DATA\VW121718\  
Data File : VW007593.D  
Acq On : 15 Dec 2018 20:02  
Operator : SY/AP  
Sample : J6378-07  
Misc : 2.70G/5ML/MSVOA\_W/SOIL  
ALS Vial : 79 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
P001-SD010-0612-01

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_W\METHOD\82W121118S.M  
Quant Title : SW846 8260

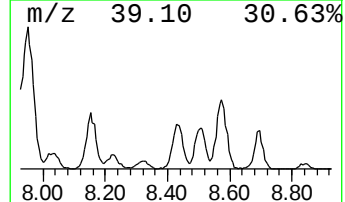
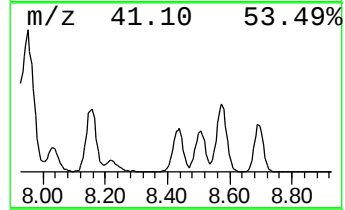
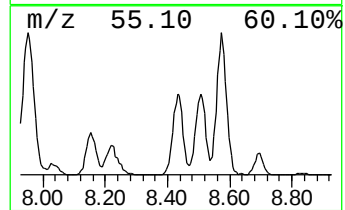
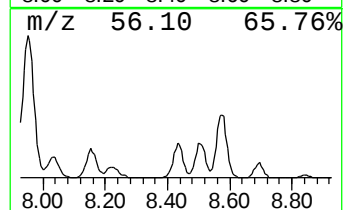
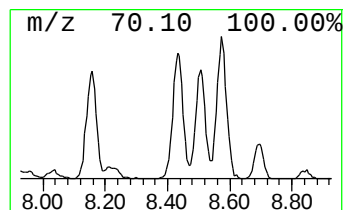
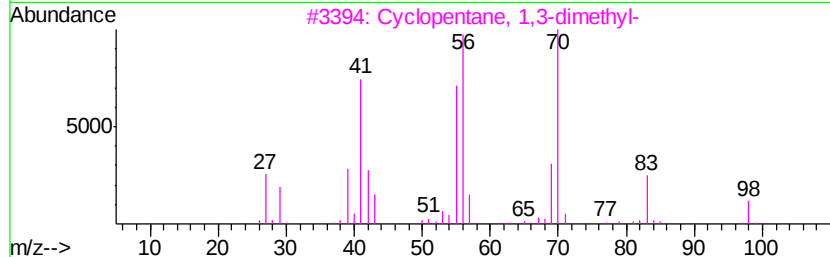
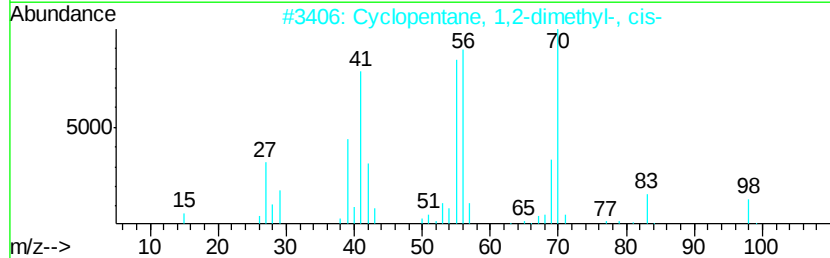
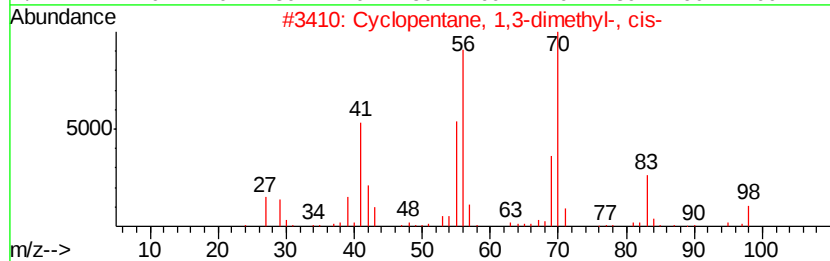
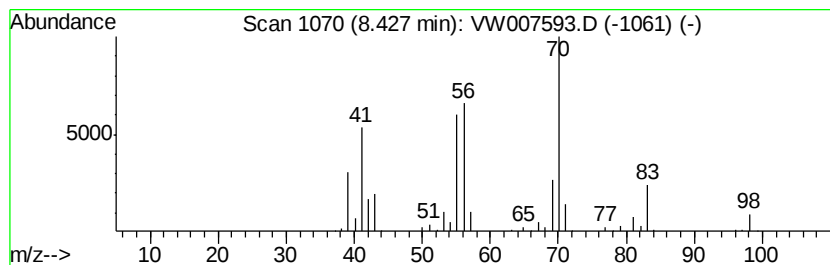
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Cyclopentane, 1,3-dimethyl-... Concentration Rank 6

| R.T. | EstConc    | Area   | Relative to ISTD    | R.T. |
|------|------------|--------|---------------------|------|
| 8.43 | 54.33 ug/l | 113172 | 1,4-Difluorobenzene | 8.84 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopentane, 1,3-dimethyl-, cis-   | 98  | C7H14   | 002532-58-3 | 91   |
| 2    |    | Cyclopentane, 1,2-dimethyl-, cis-   | 98  | C7H14   | 001192-18-3 | 90   |
| 3    |    | Cyclopentane, 1,3-dimethyl-         | 98  | C7H14   | 002453-00-1 | 90   |
| 4    |    | Cyclopentane, 1,3-dimethyl-, trans- | 98  | C7H14   | 001759-58-6 | 87   |
| 5    |    | 2H-Pyran-2,6(3H)-dione, dihydro-... | 142 | C7H10O3 | 004160-82-1 | 64   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_W\DATA\VW121718\  
Data File : VW007593.D  
Acq On : 15 Dec 2018 20:02  
Operator : SY/AP  
Sample : J6378-07  
Misc : 2.70G/5ML/MSVOA\_W/SOIL  
ALS Vial : 79 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
P001-SD010-0612-01

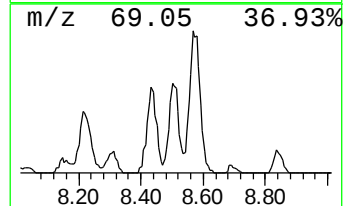
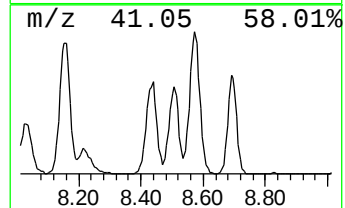
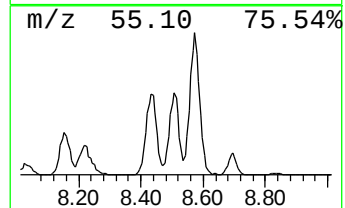
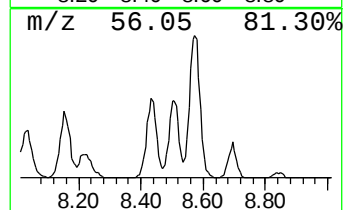
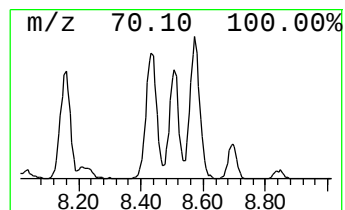
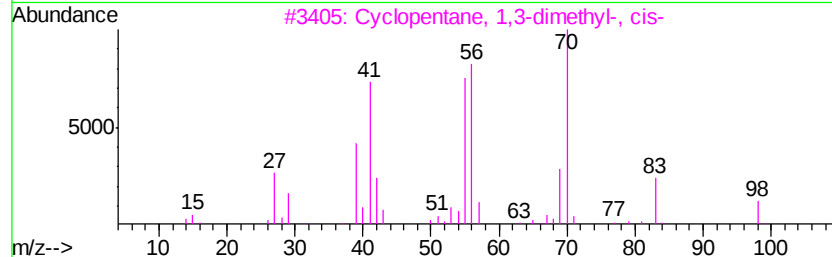
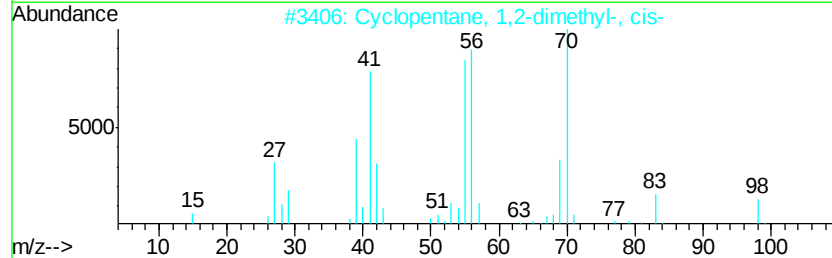
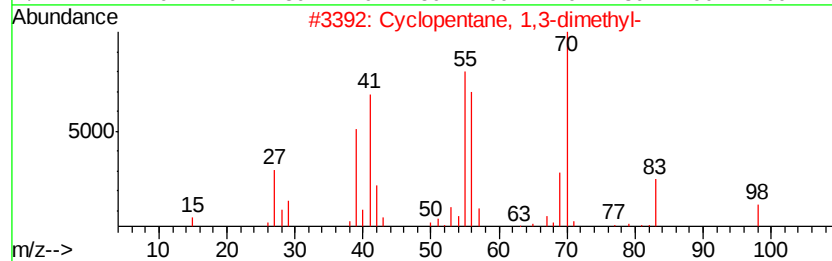
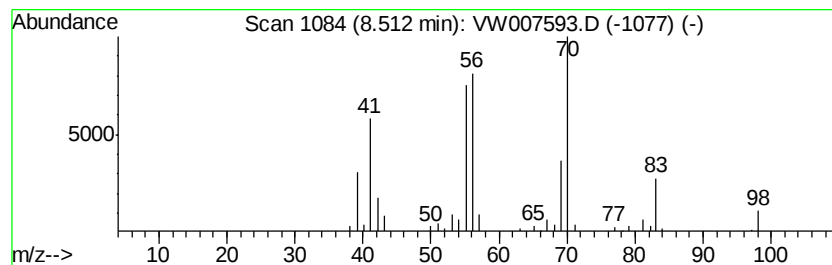
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Quant Title : SW846 8260

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

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Peak Number 7 Cyclopentane, 1,3-dimethyl- Concentration Rank 7

| R.T. | EstConc    | Area  | Relative to ISTD    | R.T. |
|------|------------|-------|---------------------|------|
| 8.51 | 47.52 ug/l | 98976 | 1,4-Difluorobenzene | 8.84 |

| Hit# | of | Tentative ID                        | MW | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|----|---------|-------------|------|
| 1    | 5  | Cyclopentane, 1,3-dimethyl-         | 98 | C7H14   | 002453-00-1 | 91   |
| 2    |    | Cyclopentane, 1,2-dimethyl-, cis-   | 98 | C7H14   | 001192-18-3 | 90   |
| 3    |    | Cyclopentane, 1,3-dimethyl-, cis-   | 98 | C7H14   | 002532-58-3 | 90   |
| 4    |    | Cyclopentane, 1,3-dimethyl-, trans- | 98 | C7H14   | 001759-58-6 | 76   |
| 5    |    | Cyclopentane, 1,2-dimethyl-         | 98 | C7H14   | 002452-99-5 | 58   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_W\DATA\VW121718\  
Data File : VW007593.D  
Acq On : 15 Dec 2018 20:02  
Operator : SY/AP  
Sample : J6378-07  
Misc : 2.70G/5ML/MSVOA\_W/SOIL  
ALS Vial : 79 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
P001-SD010-0612-01

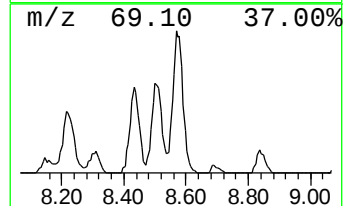
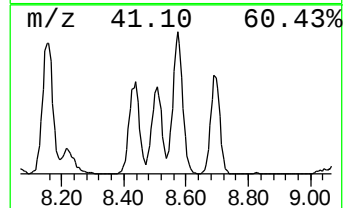
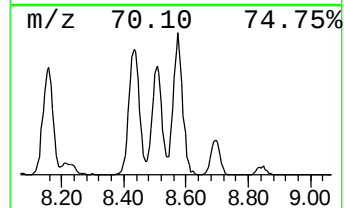
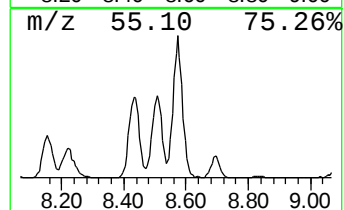
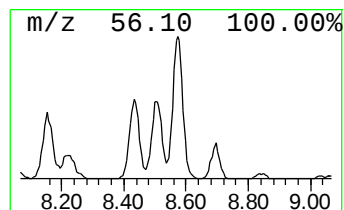
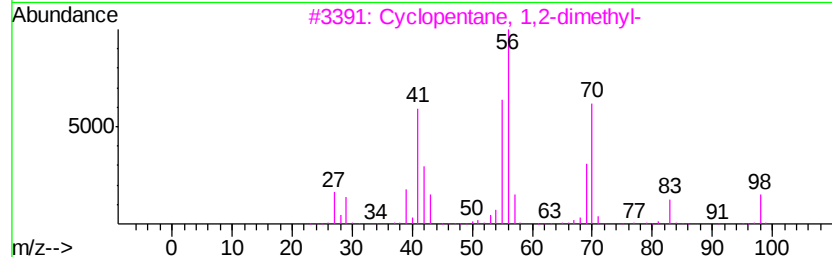
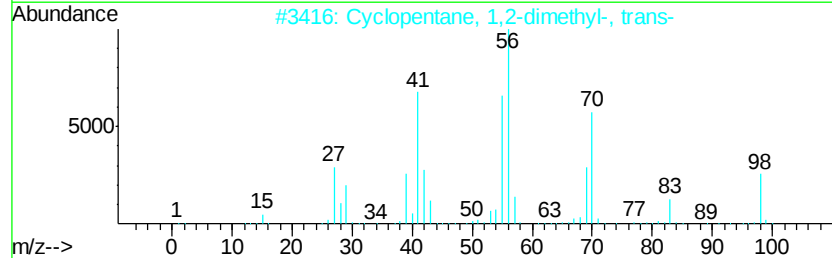
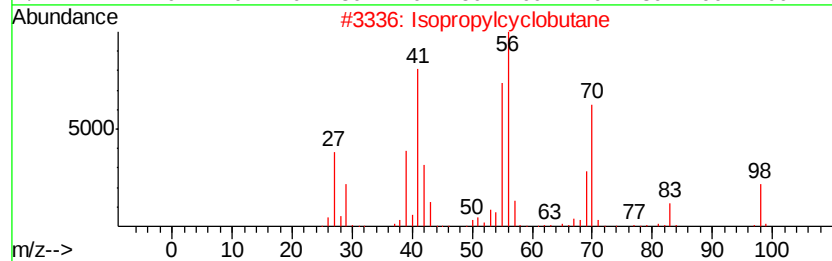
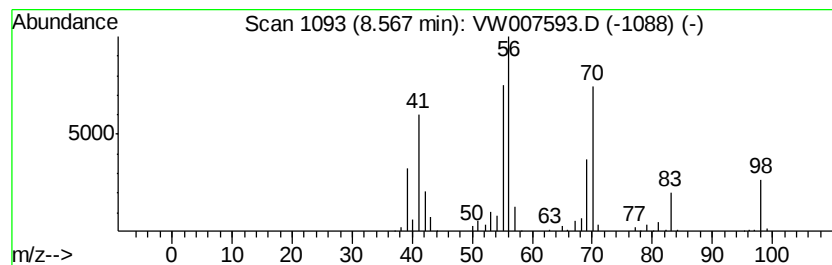
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Quant Title : SW846 8260

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

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Peak Number 8 Isopropylcyclobutane Concentration Rank 5

| R.T. | EstConc    | Area   | Relative to ISTD    | R.T. |
|------|------------|--------|---------------------|------|
| 8.57 | 76.45 ug/l | 159230 | 1,4-Difluorobenzene | 8.84 |

| Hit# | of | Tentative ID                        | MW | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|----|---------|-------------|------|
| 1    | 5  | Isopropylcyclobutane                | 98 | C7H14   | 000872-56-0 | 91   |
| 2    |    | Cyclopentane, 1,2-dimethyl-, trans- | 98 | C7H14   | 000822-50-4 | 91   |
| 3    |    | Cyclopentane, 1,2-dimethyl-         | 98 | C7H14   | 002452-99-5 | 87   |
| 4    |    | Cyclopentane, 1,3-dimethyl-         | 98 | C7H14   | 002453-00-1 | 72   |
| 5    |    | Cyclopentane, 1,3-dimethyl-, cis-   | 98 | C7H14   | 002532-58-3 | 72   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_W\DATA\VW121718\  
Data File : VW007593.D  
Acq On : 15 Dec 2018 20:02  
Operator : SY/AP  
Sample : J6378-07  
Misc : 2.70G/5ML/MSVOA\_W/SOIL  
ALS Vial : 79 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
P001-SD010-0612-01

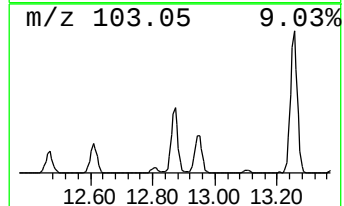
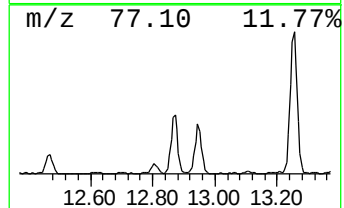
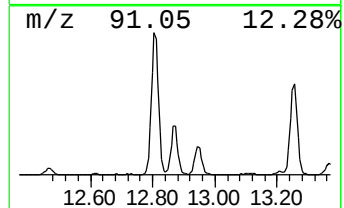
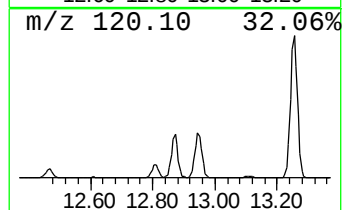
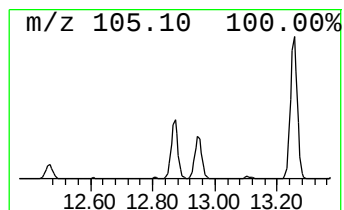
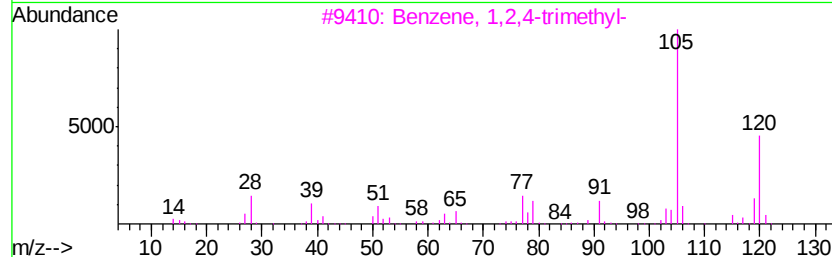
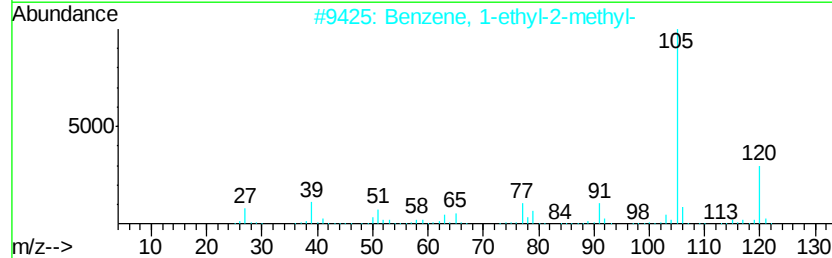
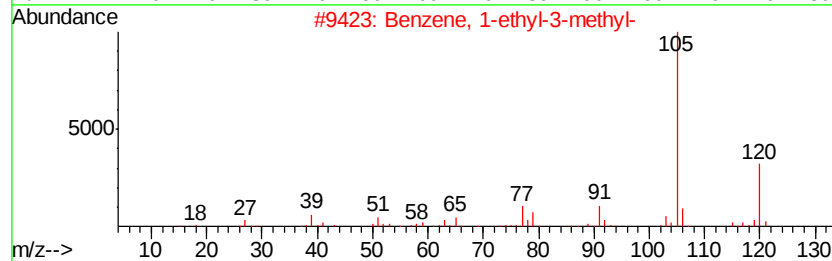
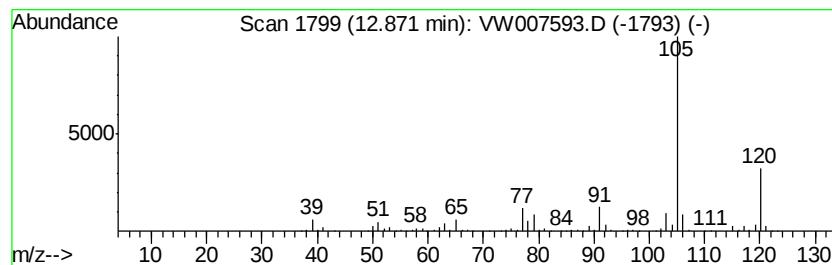
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Quant Title : SW846 8260

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

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Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 9

| R.T.  | EstConc   | Area   | Relative to ISTD       | R.T.  |
|-------|-----------|--------|------------------------|-------|
| 12.87 | 7.71 ug/l | 237457 | 1,4-Dichlorobenzene-d4 | 13.56 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 97   |
| 2    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 3    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 91   |
| 4    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 5    |    | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA\_W\DATA\VW121718\  
Data File : VW007593.D  
Acq On : 15 Dec 2018 20:02  
Operator : SY/AP  
Sample : J6378-07  
Misc : 2.70G/5ML/MSVOA\_W/SOIL  
ALS Vial : 79 Sample Multiplier: 1

Instrument :  
MSVOA\_W  
ClientSampleId :  
P001-SD010-0612-01

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA\_W\METHOD\82W121118S.M  
Quant Title : SW846 8260

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Pentane, 2-methyl-   | 4.95  | 163.5   | ug/l  | 1802520  | 1                     | 7.95  | 551284  | 50.0 |
| Pentane, 3-methyl-   | 5.43  | 180.2   | ug/l  | 1986980  | 1                     | 7.95  | 551284  | 50.0 |
| n-Hexane             | 5.93  | 303.8   | ug/l  | 3349990  | 1                     | 7.95  | 551284  | 50.0 |
| Cyclopentane, met... | 6.98  | 206.1   | ug/l  | 2271890  | 1                     | 7.95  | 551284  | 50.0 |
| Hexane, 3-methyl-    | 8.15  | 13.0    | ug/l  | 143813   | 1                     | 7.95  | 551284  | 50.0 |
| Cyclopentane, 1,3... | 8.43  | 54.3    | ug/l  | 113172   | 2                     | 8.84  | 104143  | 50.0 |
| Cyclopentane, 1,3... | 8.51  | 47.5    | ug/l  | 98976    | 2                     | 8.84  | 104143  | 50.0 |
| Isopropylcyclobutane | 8.57  | 76.5    | ug/l  | 159230   | 2                     | 8.84  | 104143  | 50.0 |
| Benzene, 1-ethyl-... | 12.87 | 7.7     | ug/l  | 237457   | 4                     | 13.56 | 1538990 | 50.0 |