

Data Path : Z:\voasrv\HPCHEM1\MSVOA W\Data\VW121718\
 Data File : VW007544.D
 Acq On : 14 Dec 2018 19:01
 Operator : SY/AP
 Sample : J6347-05
 Misc : 2.82G/5ML/MSVOA W/SOIL
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 P001-SD002-0612-01

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W121118S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.958	475	501	530	rBV	721854	3053875	3.02%	0.919%
2	5.434	562	579	613	rBV	931081	3682987	3.65%	1.108%
3	5.940	647	662	684	rBV	2776516	8688900	8.61%	2.614%
4	6.982	821	833	852	rVB	2031731	6517859	6.46%	1.961%
5	7.952	972	992	1002	rBV3	447925	1633145	1.62%	0.491%
6	9.336	1204	1219	1236	rBV	961435	2210100	2.19%	0.665%
7	10.402	1375	1394	1399	rBV3	32907336	100971723	100.00%	30.380%
8	11.664	1591	1601	1606	rBV2	555319	1103018	1.09%	0.332%
9	11.744	1606	1614	1623	rVB	20504487	39117625	38.74%	11.769%
10	11.853	1623	1632	1635	rBV3	35685597	86681683	85.85%	26.080%
11	12.183	1673	1686	1694	rBV	29065747	52770339	52.26%	15.877%
12	12.615	1752	1757	1762	rVB	1350216	2158700	2.14%	0.649%
13	12.811	1781	1789	1793	rBV	726884	1186934	1.18%	0.357%
14	12.871	1793	1799	1802	rBV	2195027	3460318	3.43%	1.041%
15	12.902	1802	1804	1807	rVV	890262	1329684	1.32%	0.400%
16	12.945	1807	1811	1817	rVB2	1451673	2594227	2.57%	0.781%
17	13.109	1831	1838	1844	rBV	733587	1241953	1.23%	0.374%
18	13.256	1856	1862	1868	rBV	3692688	5819715	5.76%	1.751%
19	13.585	1908	1916	1922	rBV2	1612534	2973150	2.94%	0.895%
20	13.877	1959	1964	1969	rBV2	953162	1409174	1.40%	0.424%
21	14.079	1994	1997	2005	rVB3	574793	1515475	1.50%	0.456%
22	14.804	2110	2116	2121	rBV2	442605	1055777	1.05%	0.318%
23	15.371	2205	2209	2216	rVB	672738	1188893	1.18%	0.358%

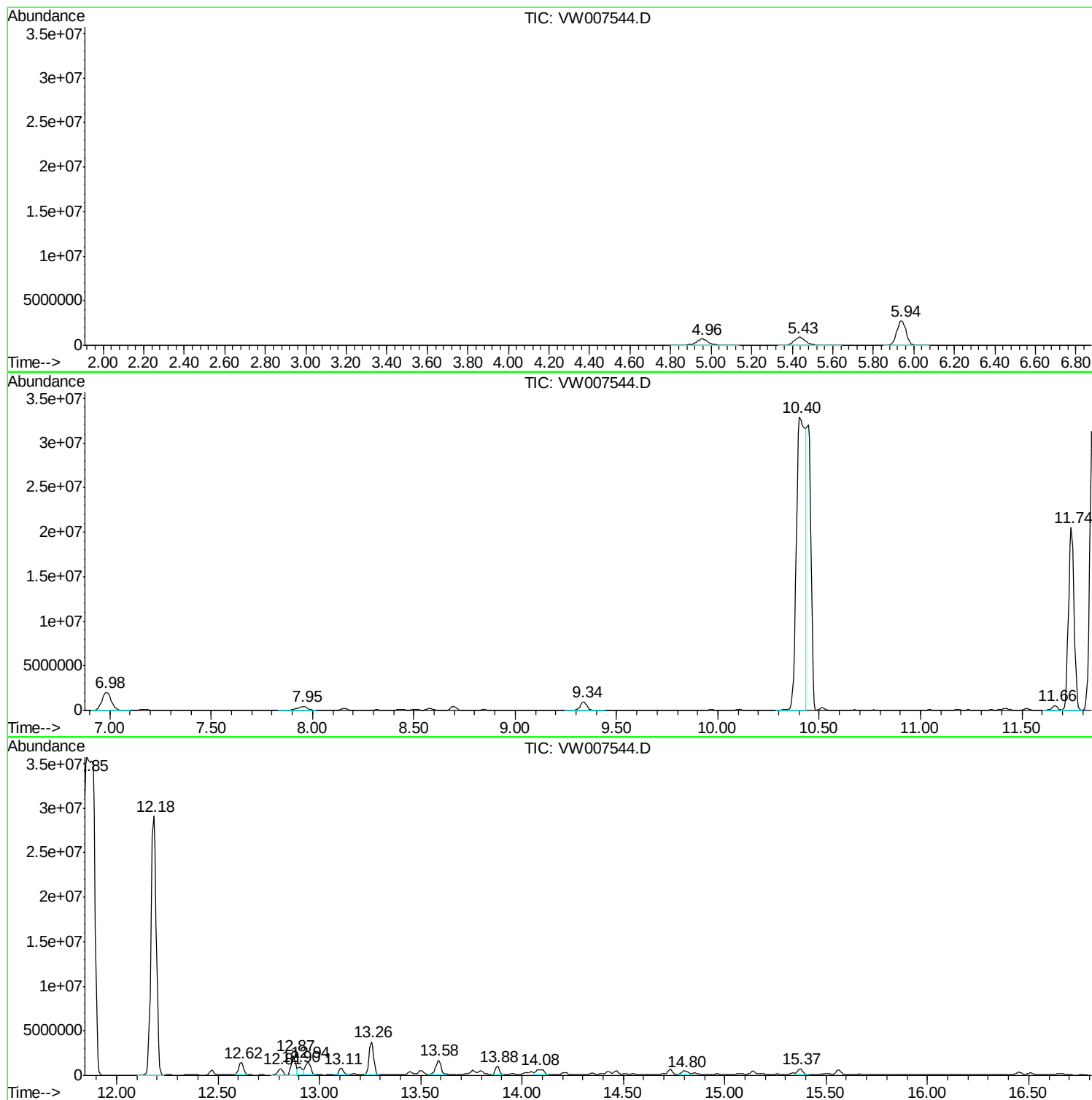
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Quant Title : SW846 8260

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



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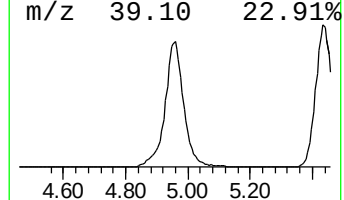
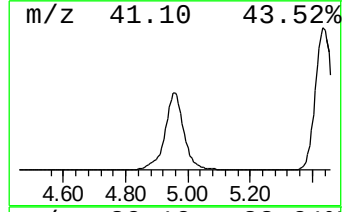
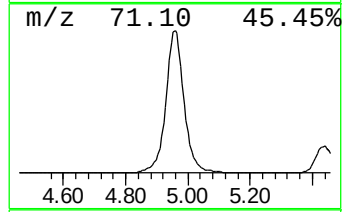
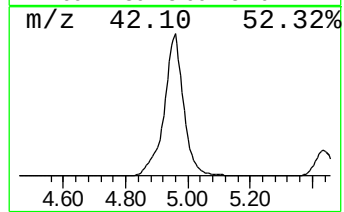
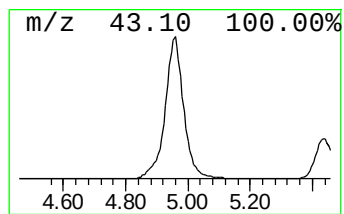
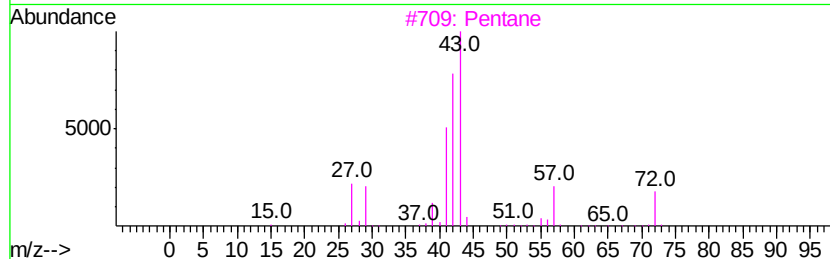
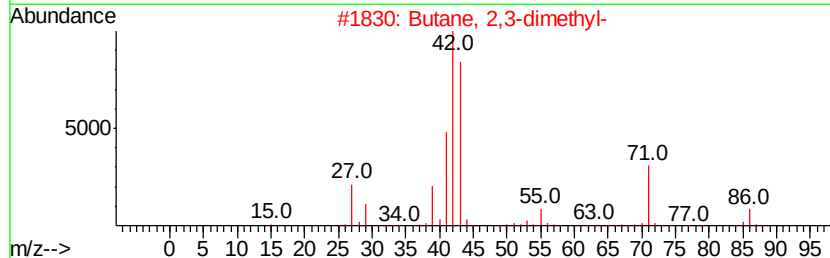
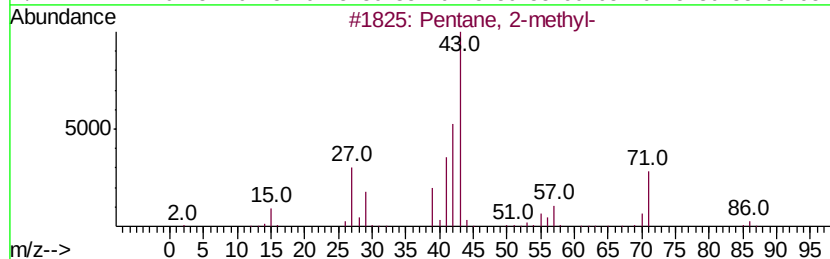
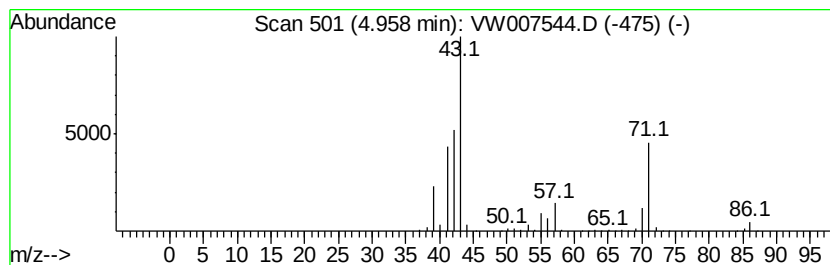
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W121118S.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	93.50 ug/l	3053880	Pentafluorobenzene	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	49
3		Pentane	72	C5H12	000109-66-0	43
4		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	42
5		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38



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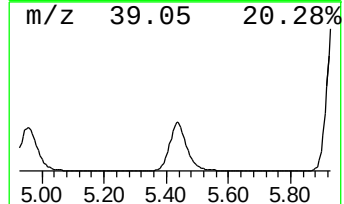
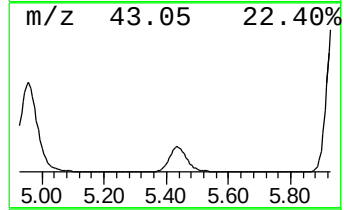
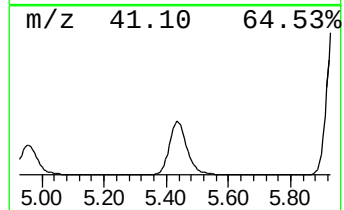
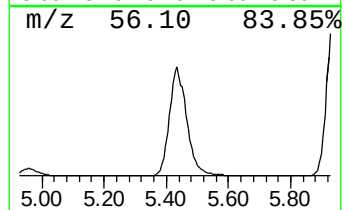
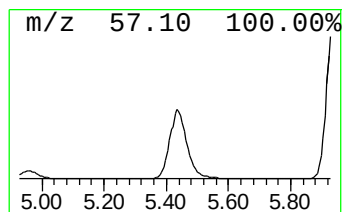
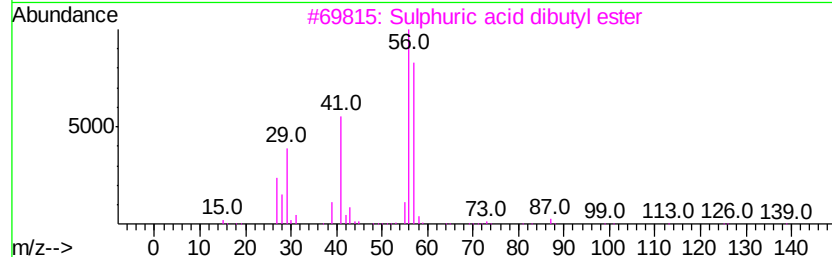
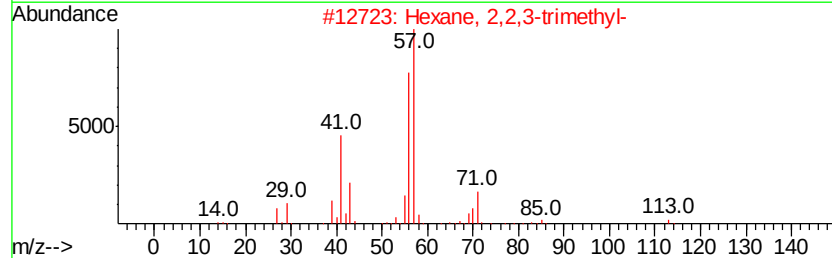
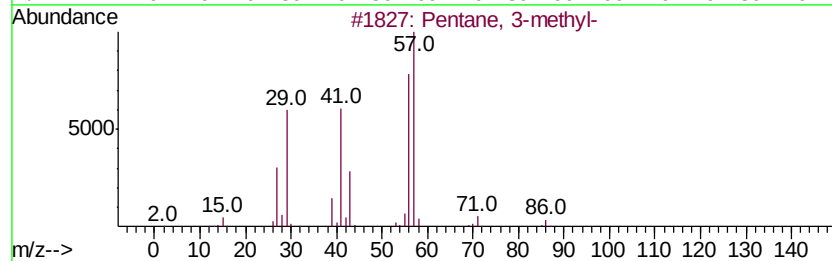
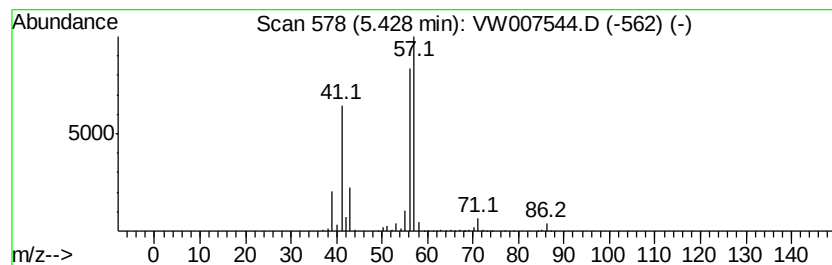
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W121118S.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.43	112.76 ug/l	3682990	Pentafluorobenzene	7.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	90
2			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
3			Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40
4			Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	40
5			n-Hexane	86	C6H14	000110-54-3	38



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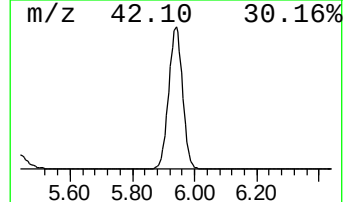
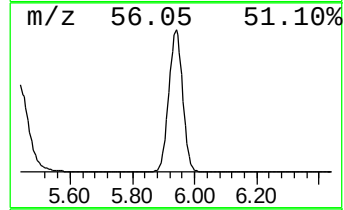
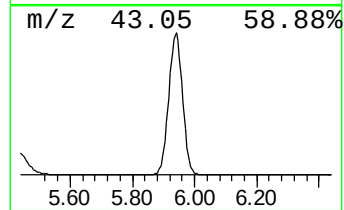
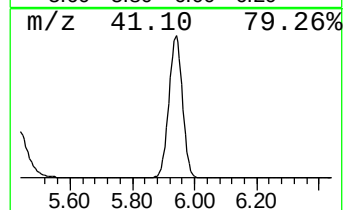
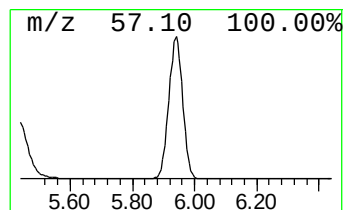
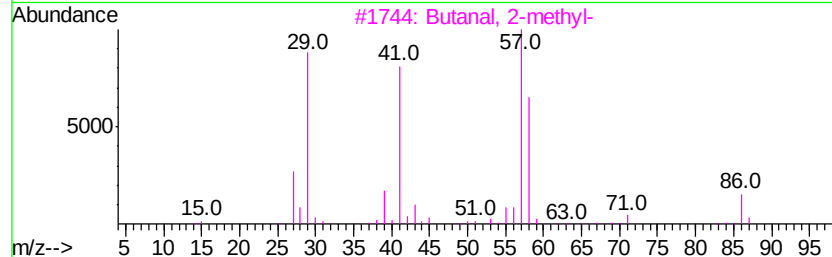
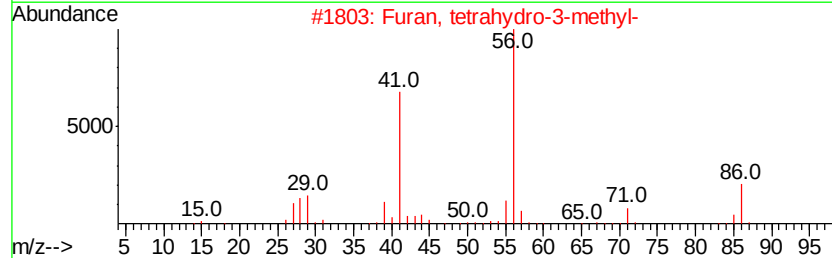
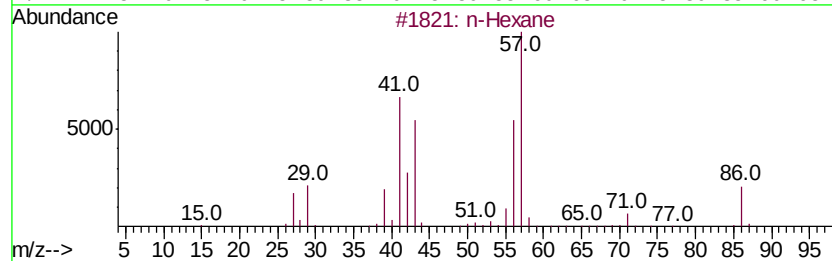
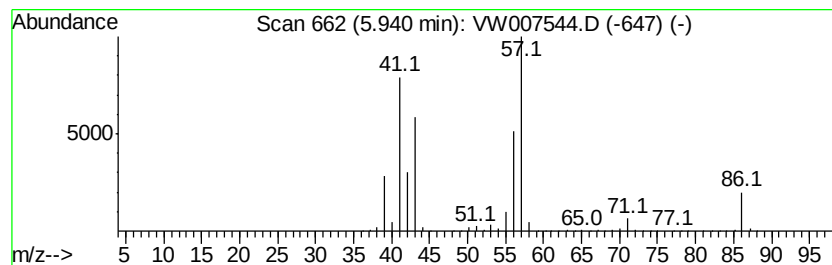
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.94	266.02 ug/l	8688900	Pentafluorobenzene	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	94
2		Furan, tetrahydro-3-methyl-	86	C5H10O	013423-15-9	50
3		Butanal, 2-methyl-	86	C5H10O	000096-17-3	43
4		Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	43
5		Pentane, 3-methyl-	86	C6H14	000096-14-0	38



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Misc : 2.82G/5ML/MSVOA W/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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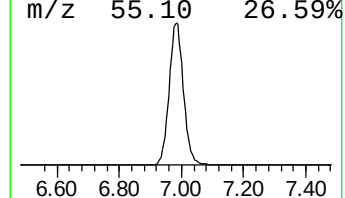
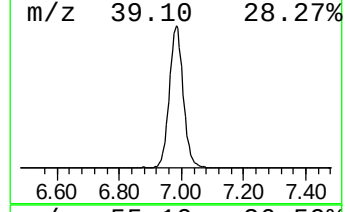
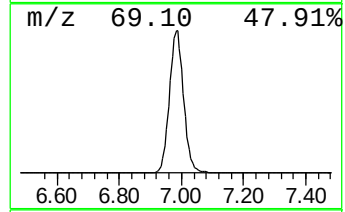
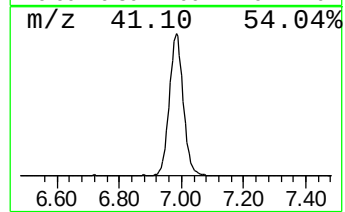
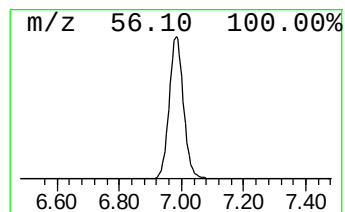
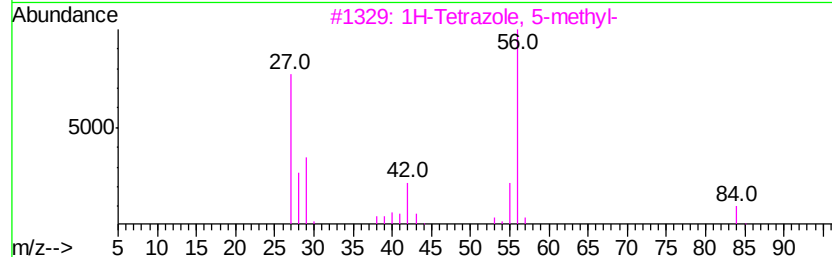
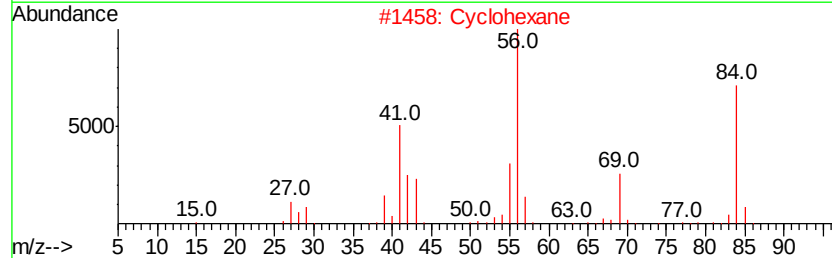
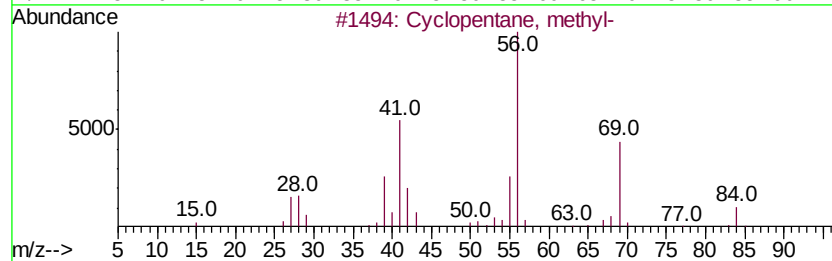
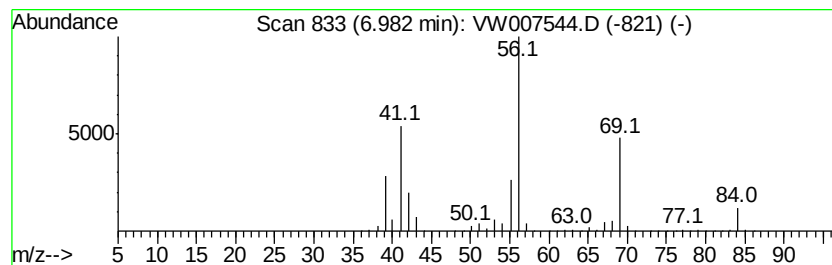
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Quant Title : SW846 8260

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

Peak Number 4 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.98	199.55 ug/l	6517860	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		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	59
5		Cyclobutane	56	C4H8	000287-23-0	53



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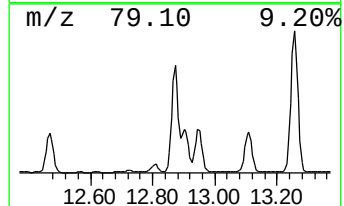
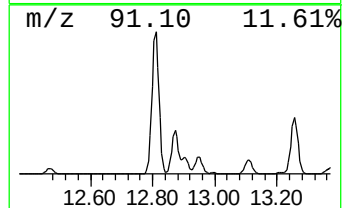
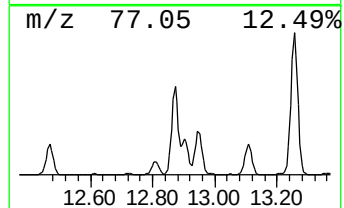
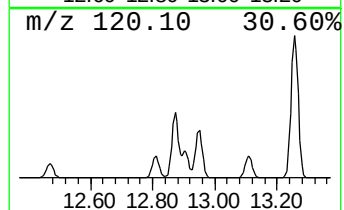
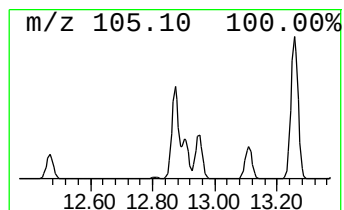
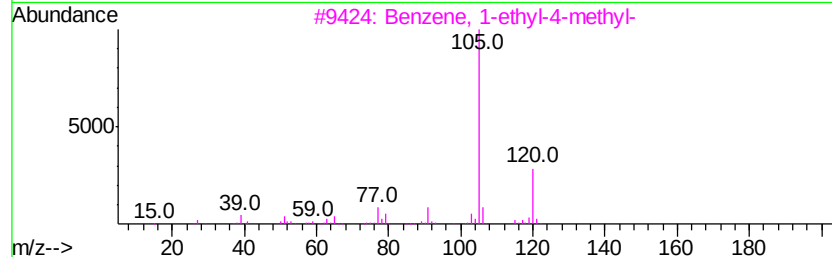
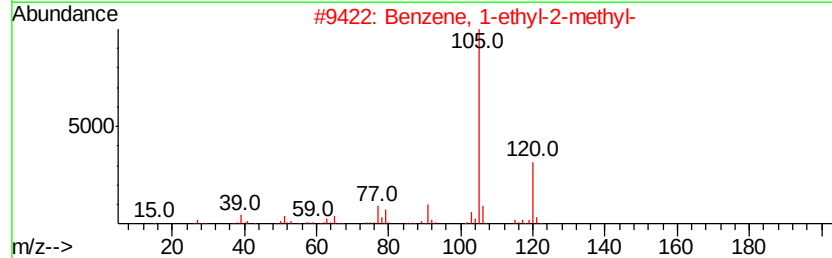
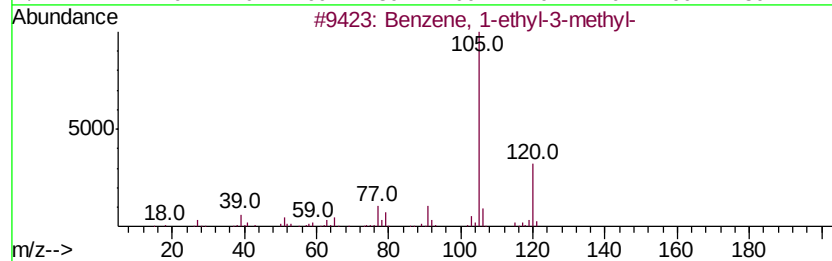
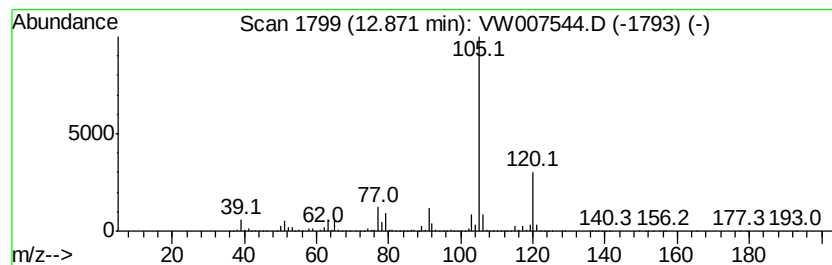
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	58.19 ug/l	3460320	1,4-Dichlorobenzene-d4	13.56
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-2-methyl-	120 C9H12	000611-14-3 95
3		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 91
4		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 91
5		Mesitylene	120 C9H12	000108-67-8 91



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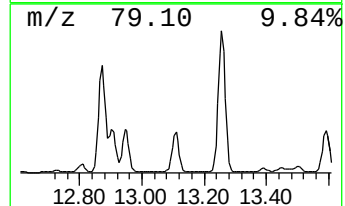
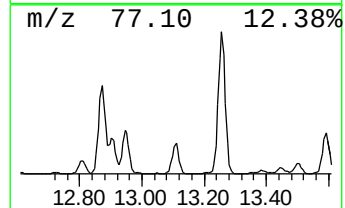
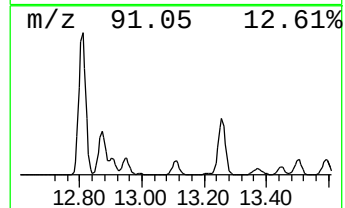
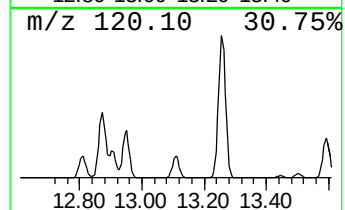
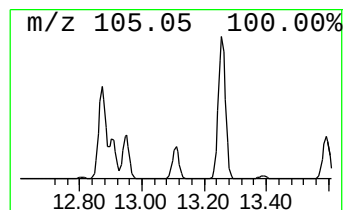
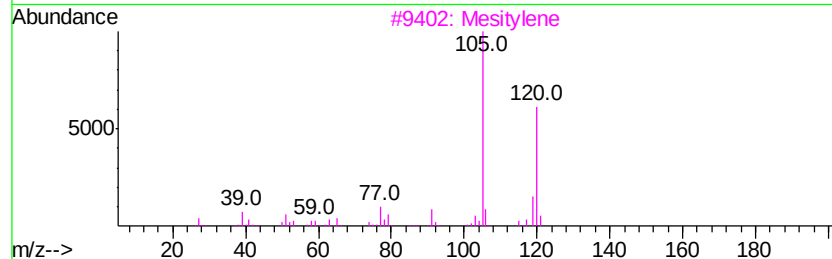
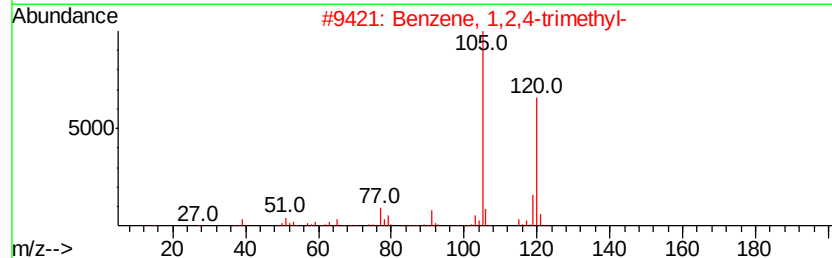
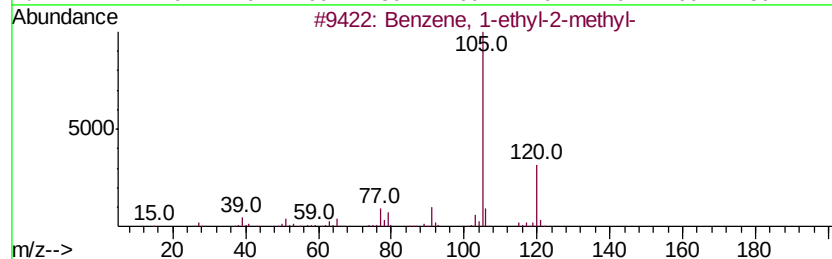
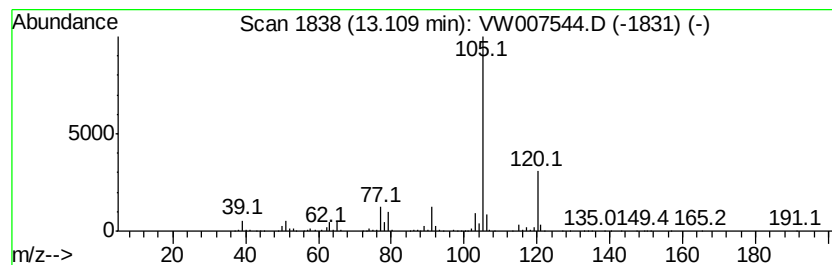
Instrument :
MSVOA_W
ClientSampleId :
P001-SD002-0612-01

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W121118S.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.11	20.89 ug/l	1241950	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 94
2		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 91
3		Mesitylene	120 C9H12	000108-67-8 91
4		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 91
5		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 91



Data Path : Z:\voasrv\HPCHEM1\MSVOA W\Data\VW121718\
Data File : VW007544.D
Acq On : 14 Dec 2018 19:01
Operator : SY/AP
Sample : J6347-05
Misc : 2.82G/5ML/MSVOA W/SOIL
ALS Vial : 21 Sample Multiplier: 1

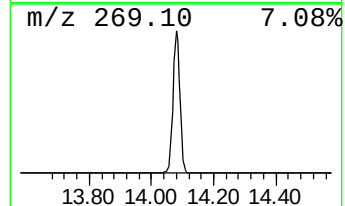
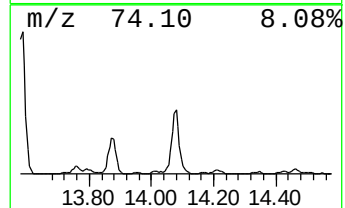
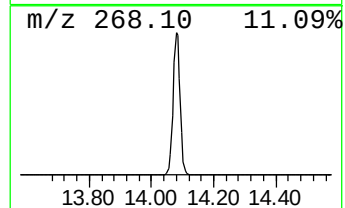
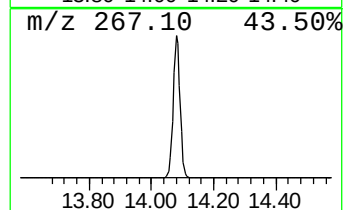
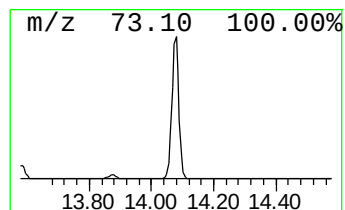
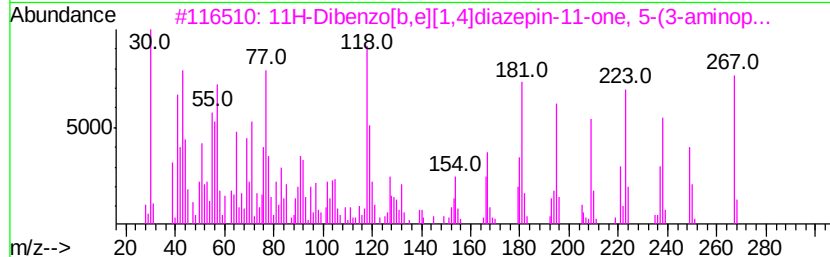
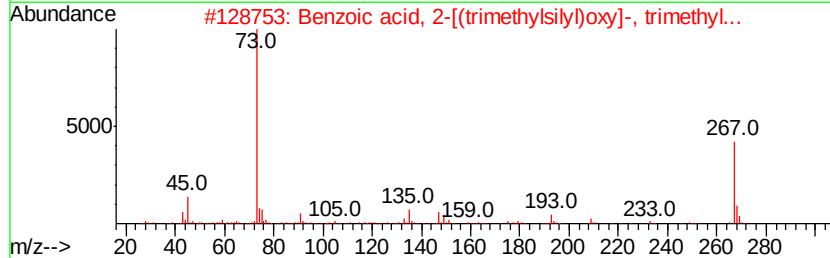
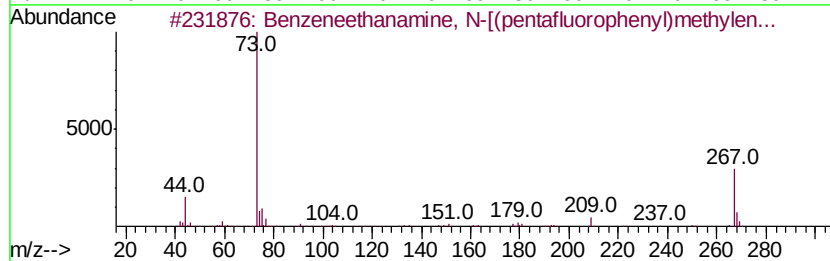
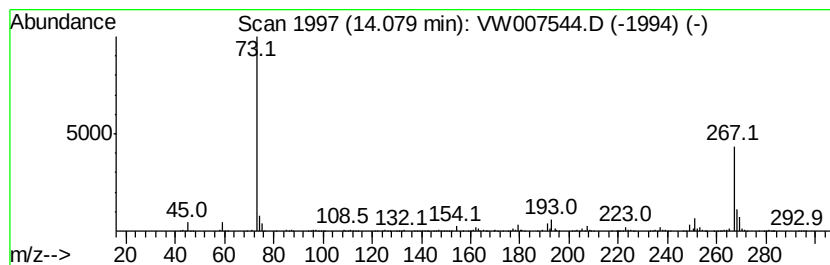
Instrument :
MSVOA_W
ClientSampleId :
P001-SD002-0612-01

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W121118S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzeneethanamine, N-[(pent... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.08	25.49 ug/l	1515480	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzeneethanamine, N-[(pentafluoro...]	475 C21H26F5N02Si2	055429-85-1 64
2		Benzoic acid, 2-[(trimethylsilyl)...]	282 C13H22O3Si2	003789-85-3 50
3		11H-Dibenzo[b,e][1,4]diazepin-11...	267 C16H17N3O	013450-73-2 43
4		Trimethylsilyloxvethane, 2-(3-tr...	282 C14H26O2Si2	1000365-49-0 25
5		4-Methylcatechol, bis(trimethyls...	268 C13H24O2Si2	1000332-79-9 22



Data Path : Z:\voasrv\HPCHEM1\MSVOA W\Data\VW121718\
Data File : VW007544.D
Acq On : 14 Dec 2018 19:01
Operator : SY/AP
Sample : J6347-05
Misc : 2.82G/5ML/MSVOA W/SOIL
ALS Vial : 21 Sample Multiplier: 1

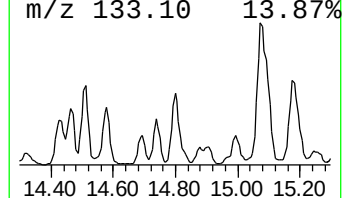
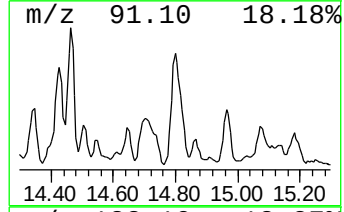
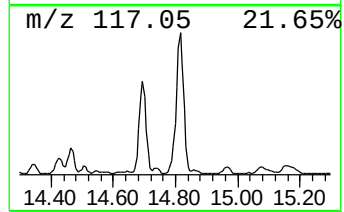
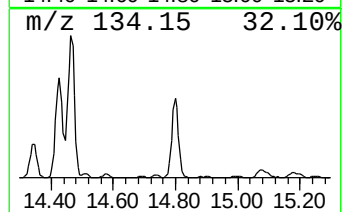
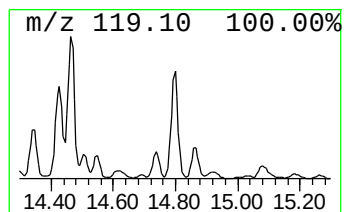
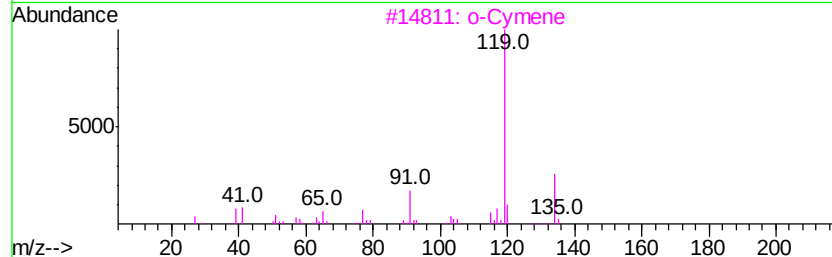
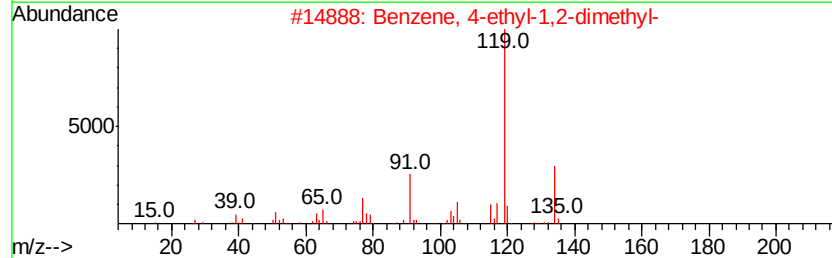
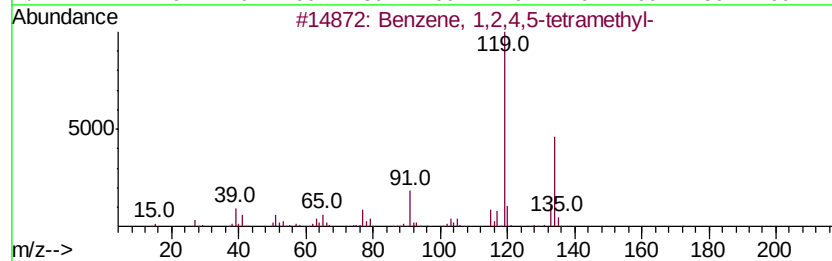
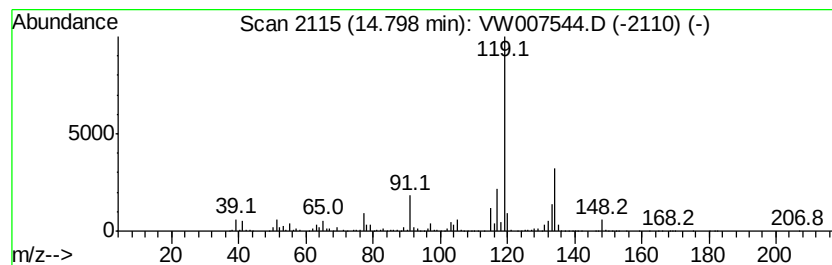
Instrument :
MSVOA_W
ClientSampleId :
P001-SD002-0612-01

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W121118S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	17.76 ug/l	1055780	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 64
2		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 60
3		o-Cymene	134 C10H14	000527-84-4 60
4		p-Cymene	134 C10H14	000099-87-6 60
5		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW121718\
Data File : VW007544.D
Acq On : 14 Dec 2018 19:01
Operator : SY/AP
Sample : J6347-05
Misc : 2.82G/5ML/MSVOA_W/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
P001-SD002-0612-01

Quant Method : Z:\VOASRV\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-	4.96	93.5	ug/l	3053880	1	7.95	1633150	50.0
Pentane, 3-methyl-	5.43	112.8	ug/l	3682990	1	7.95	1633150	50.0
n-Hexane	5.94	266.0	ug/l	8688900	1	7.95	1633150	50.0
Cyclopentane, met...	6.98	199.6	ug/l	6517860	1	7.95	1633150	50.0
Benzene, 1-ethyl-...	12.87	58.2	ug/l	3460320	4	13.56	2973150	50.0
Benzene, 1-ethyl-...	13.11	20.9	ug/l	1241950	4	13.56	2973150	50.0
Benzeneethanamine...	14.08	25.5	ug/l	1515480	4	13.56	2973150	50.0
Benzene, 1,2,4,5-...	14.80	17.8	ug/l	1055780	4	13.56	2973150	50.0