

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102519\
 Data File : VX013232.D
 Acq On : 25 Oct 2019 15:40
 Operator : JC/SP
 Sample : K5546-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GW810-8.0-13.0-102419

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_X\METHOD\82X102419W.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	1.070	26	30	43	rBV	332406	413039	66.51%	12.566%
2	1.167	43	46	56	rVB	29405	30363	4.89%	0.924%
3	1.265	58	62	71	rBV	14544	20732	3.34%	0.631%
4	1.381	78	81	87	rVB2	10069	15412	2.48%	0.469%
5	2.436	247	254	260	rBV	5054	8407	1.35%	0.256%
6	2.594	268	280	289	rBV	19716	36423	5.87%	1.108%
7	5.490	743	755	769	rBV2	77363	218930	35.25%	6.660%
8	5.648	769	781	794	rBV	82334	233274	37.56%	7.097%
9	6.057	836	848	862	rBV	88881	233185	37.55%	7.094%
10	6.849	969	978	992	rBV	117791	274408	44.19%	8.348%
11	8.709	1276	1283	1292	rBV	404066	620996	100.00%	18.893%
12	8.782	1292	1295	1304	rVB	5174	8474	1.36%	0.258%
13	10.111	1507	1513	1522	rBV	250645	345631	55.66%	10.515%
14	11.135	1675	1681	1690	rBV	303944	391730	63.08%	11.918%
15	11.361	1713	1718	1723	rBV3	5478	7364	1.19%	0.224%
16	12.074	1829	1835	1846	rBV	356304	428623	69.02%	13.040%

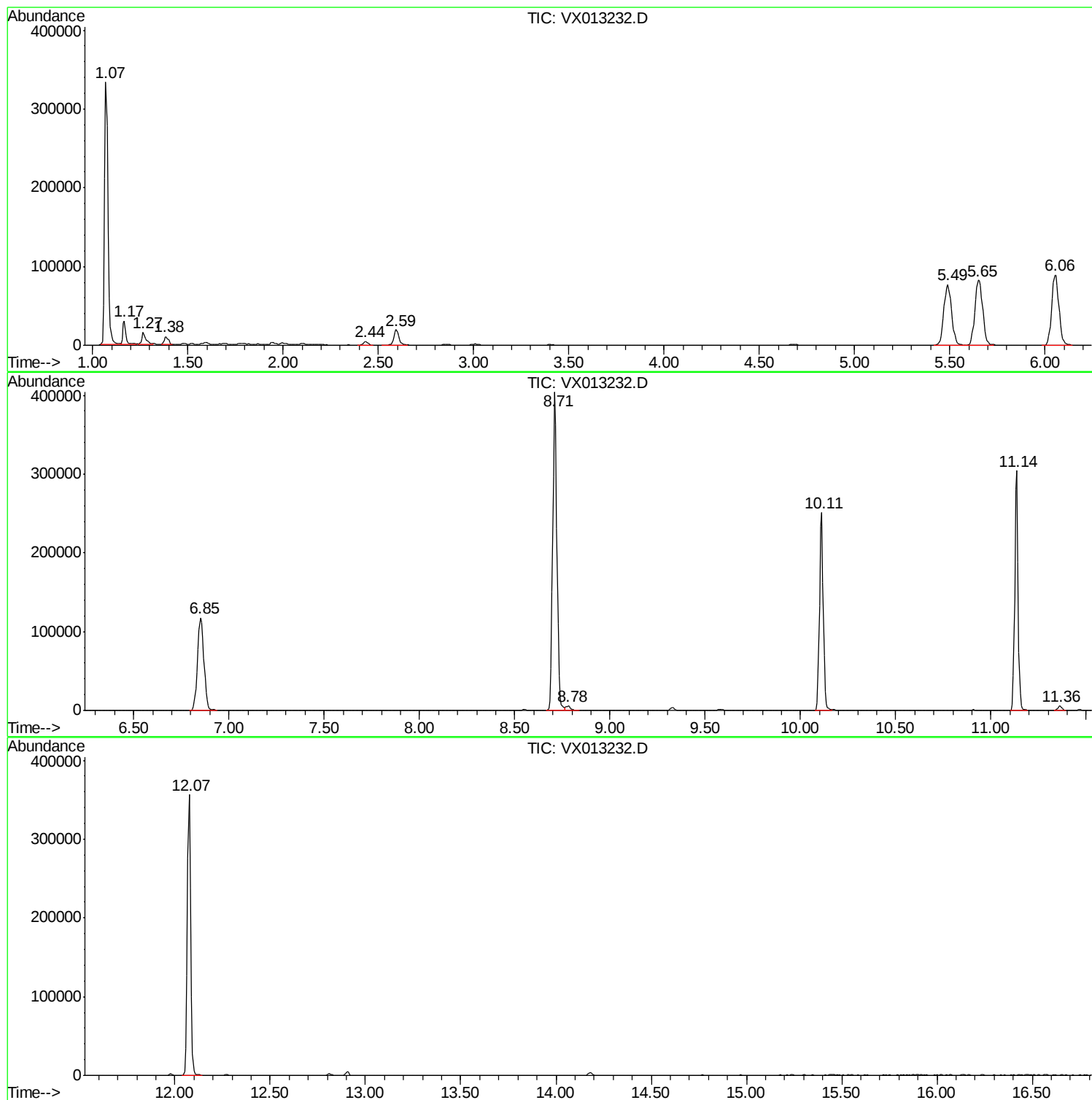
Sum of corrected areas: 3286991

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X102419W.M
Quant Title : SW846 8260

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



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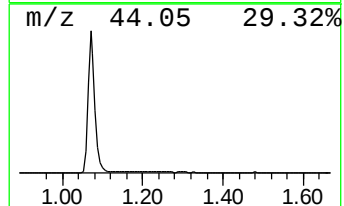
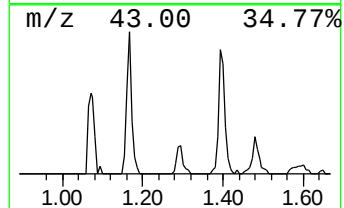
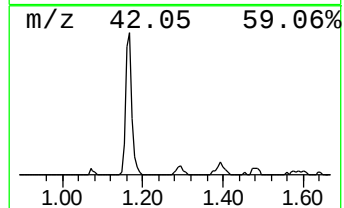
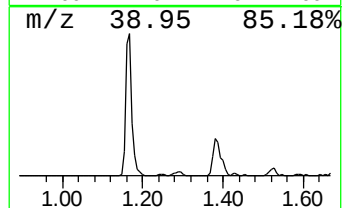
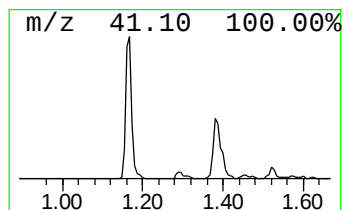
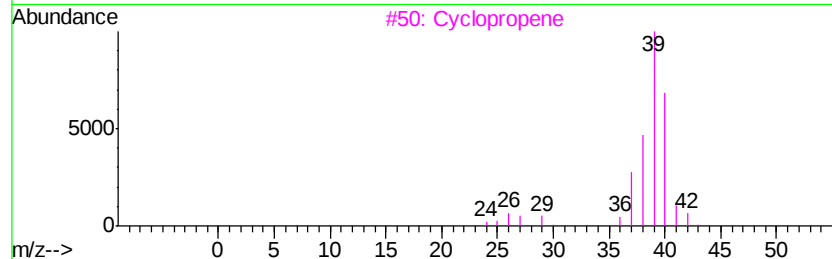
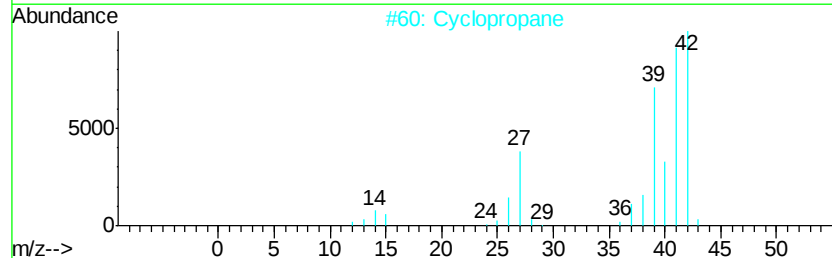
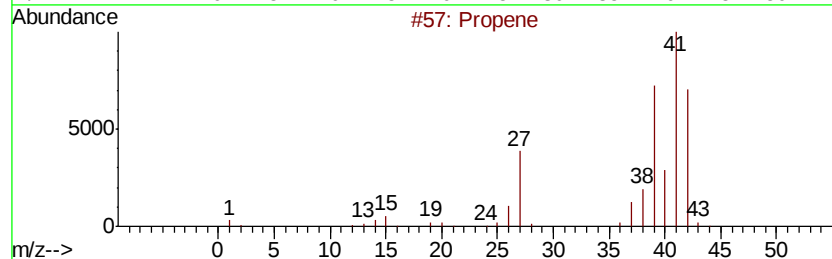
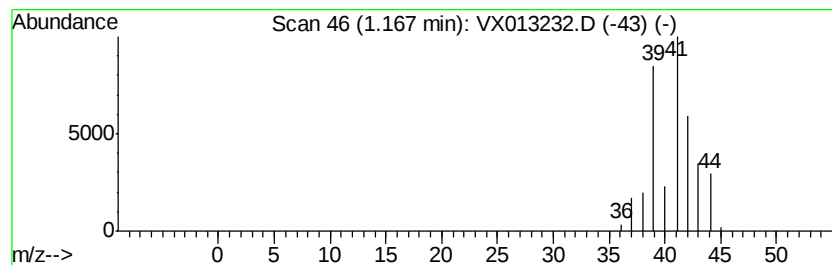
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Peak Number 2 Propene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.17	6.51 ug/l	30363	Pentafluorobenzene	5.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene		42	C3H6	000115-07-1	83
2	Cyclopropane		42	C3H6	000075-19-4	40
3	Cyclopropene		40	C3H4	002781-85-3	5
4	Acetonitrile		41	C2H3N	000075-05-8	3
5	2-Oxetanone, 4-methyl-		86	C4H6O2	003068-88-0	2



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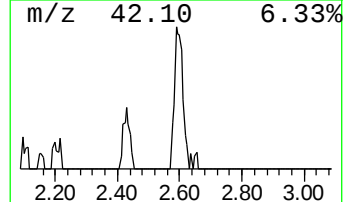
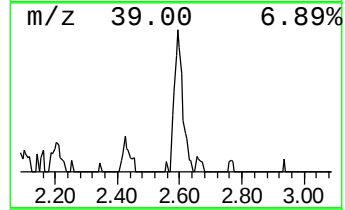
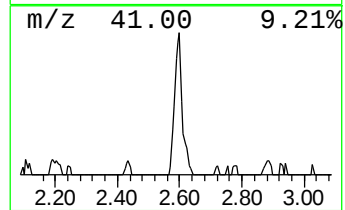
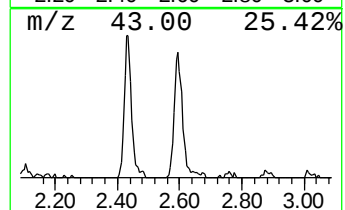
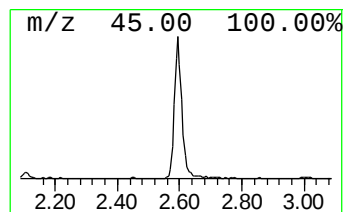
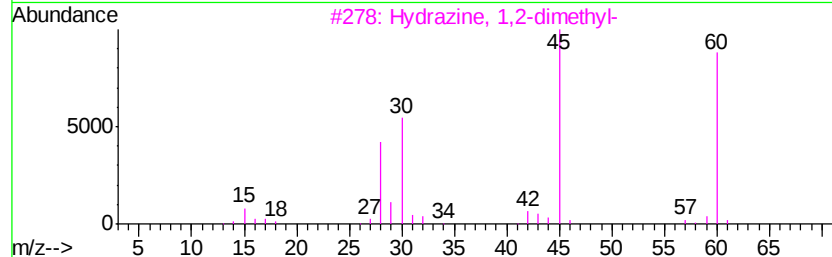
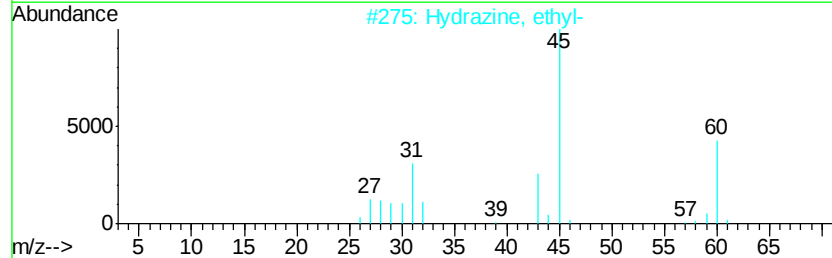
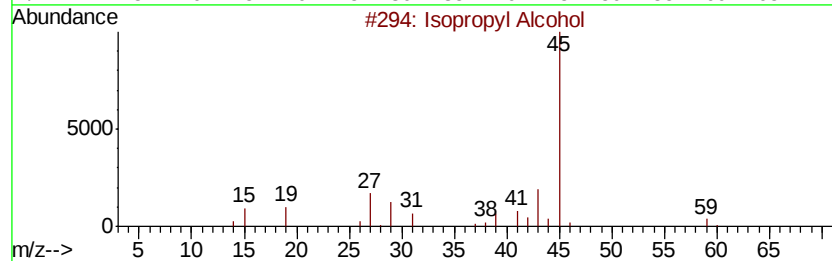
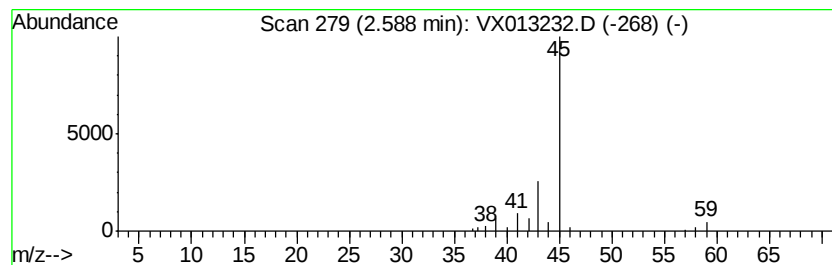
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Peak Number 3 Isopropyl Alcohol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.59	7.81 ug/l	36423	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropyl Alcohol	60	C3H8O	000067-63-0	78
2		Hydrazine, ethyl-	60	C2H8N2	000624-80-6	40
3		Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	9
4		Ethylamine	45	C2H7N	000075-04-7	5
5		2-Propanone, 1-methoxy-	88	C4H8O2	005878-19-3	4



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propene	1.17	6.5	ug/l	30363	1	5.65	233274	50.0
Isopropyl Alcohol	2.59	7.8	ug/l	36423	1	5.65	233274	50.0