

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102620\
 Data File : VU040960.D
 Acq On : 26 Oct 2020 22:48
 Operator : SY/MD
 Sample : L4492-16
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 C0AP4

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

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_U\METHOD\SOMUTR102620WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.366	3	8	22	rVB	162031	156709	1.17%	1.010%
2	1.604	71	82	94	rBV3	96645	144650	1.08%	0.932%
3	4.665	1013	1034	1081	rBV2	36047	148176	1.11%	0.955%
4	5.742	1346	1369	1392	rBV2	103069	271257	2.03%	1.748%
5	6.263	1517	1531	1544	rBV	87573	159180	1.19%	1.026%
6	7.906	2029	2042	2055	rBV	116934	196362	1.47%	1.265%
7	8.562	2229	2246	2261	rBV	8027679	13392904	100.00%	86.288%
8	8.639	2261	2270	2302	rVB	233864	483009	3.61%	3.112%
9	9.420	2501	2513	2534	rBV	130015	213357	1.59%	1.375%
10	11.806	3240	3255	3270	rBV	112013	176682	1.32%	1.138%
11	12.189	3362	3374	3387	rBV	112267	178869	1.34%	1.152%

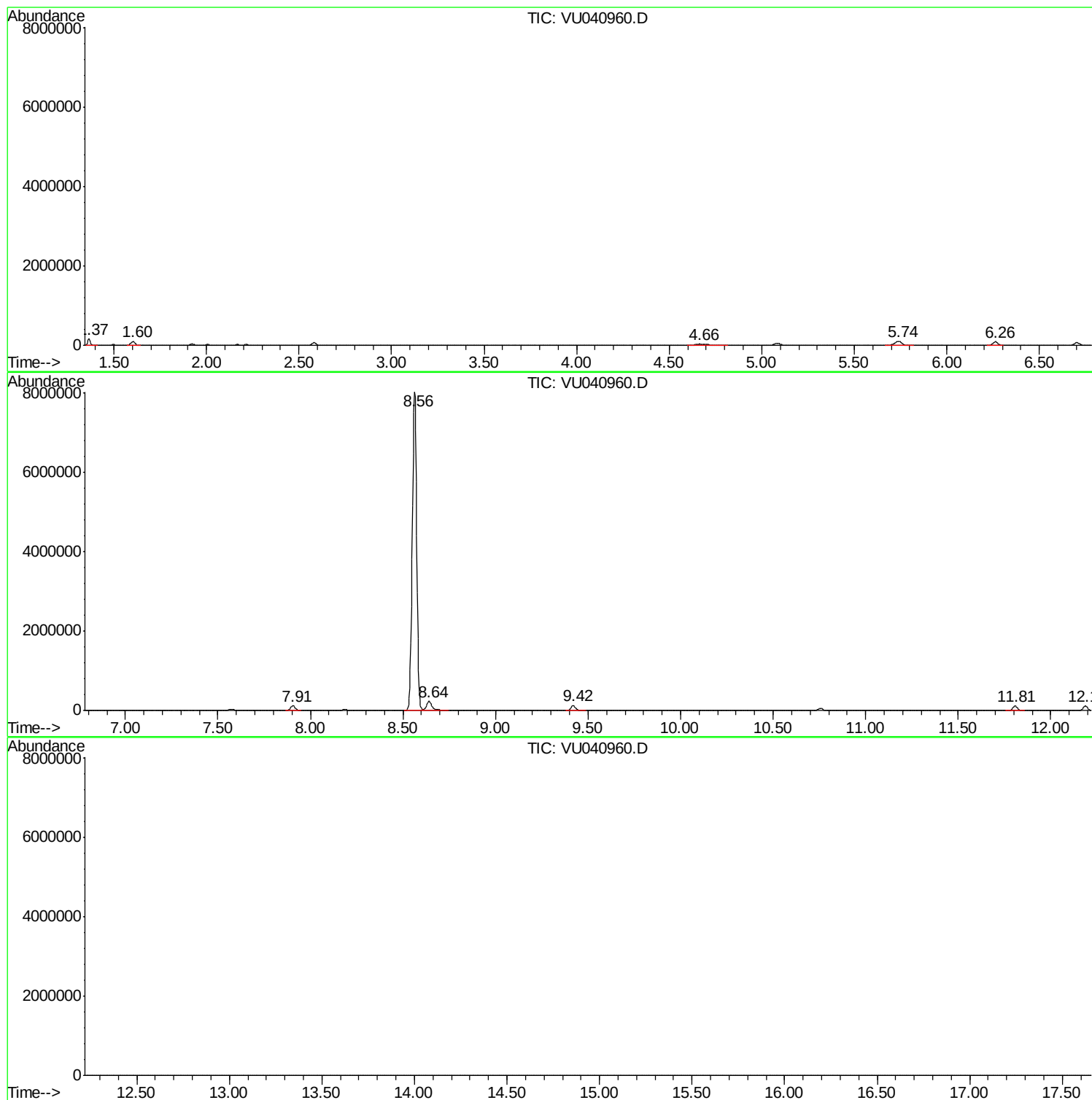
Sum of corrected areas: 15521155

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR102620WMA.M
Quant Title : TRACE VOA SOM01.0

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



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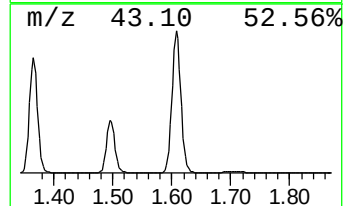
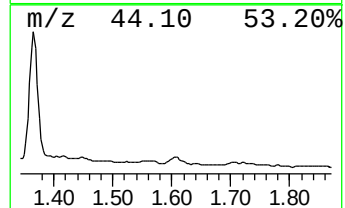
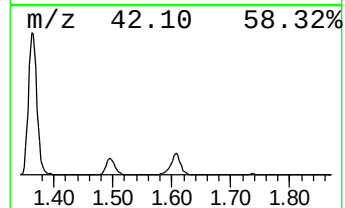
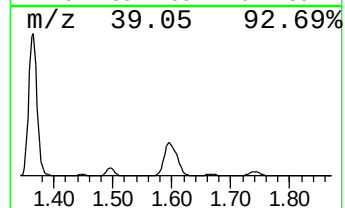
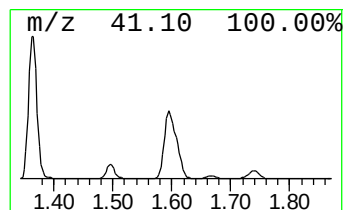
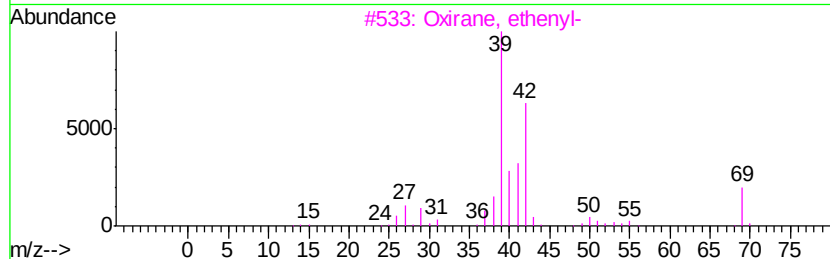
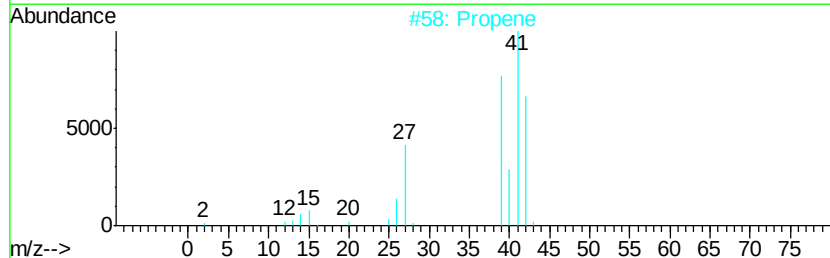
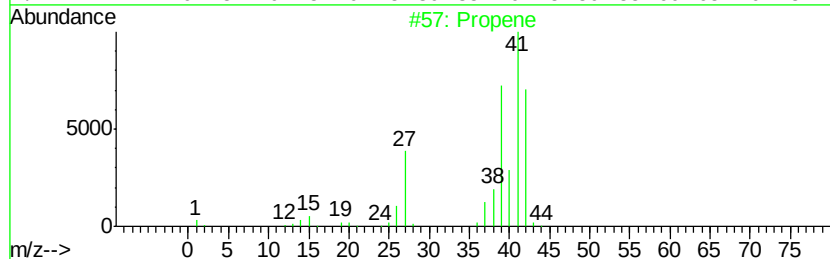
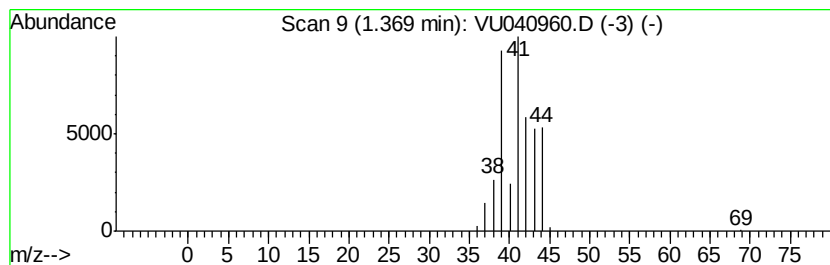
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR102620WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.37	4.92 ug/L	156709	1,4-Difluorobenzene	6.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propene	42	C3H6	000115-07-1	64
2			Propene	42	C3H6	000115-07-1	42
3			Oxirane, ethenyl-	70	C4H6O	000930-22-3	9
4			Furan	68	C4H4O	000110-00-9	9
5			2-Butenenitrile	67	C4H5N	004786-20-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	1.37	4.9	ug/L	156709	1	6.26	159180	5.0