

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061622\
 Data File : VD073623.D
 Acq On : 16 Jun 2022 17:58
 Operator : VA/SY
 Sample : N3351-03
 Misc : 5.06G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 40588

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_D\Method\82D060922S.M

Title : SW846 8260

Signal : TIC: VD073623.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.950	3	5	14	rVB	113166	124004	8.75%	1.720%
2	2.314	62	67	71	rBV2	29154	50880	3.59%	0.706%
3	4.961	506	517	524	rBV3	14922	48053	3.39%	0.666%
4	5.985	680	691	702	rBV5	12191	36121	2.55%	0.501%
5	7.908	1008	1018	1023	rBV	233734	547999	38.68%	7.600%
6	7.973	1023	1029	1042	rVB	340711	760141	53.66%	10.542%
7	8.326	1080	1089	1101	rBV	212166	467690	33.01%	6.486%
8	8.855	1171	1179	1191	rBV	540199	1084472	76.55%	15.040%
9	10.332	1422	1430	1438	rBV	793896	1416600	100.00%	19.646%
10	11.637	1643	1652	1663	rBV	600610	1057415	74.64%	14.665%
11	12.620	1812	1819	1831	rBV2	697700	1130410	79.80%	15.677%
12	13.561	1972	1979	1986	rBV	219535	350248	24.72%	4.857%
13	14.084	2062	2068	2074	rVB	86709	136651	9.65%	1.895%

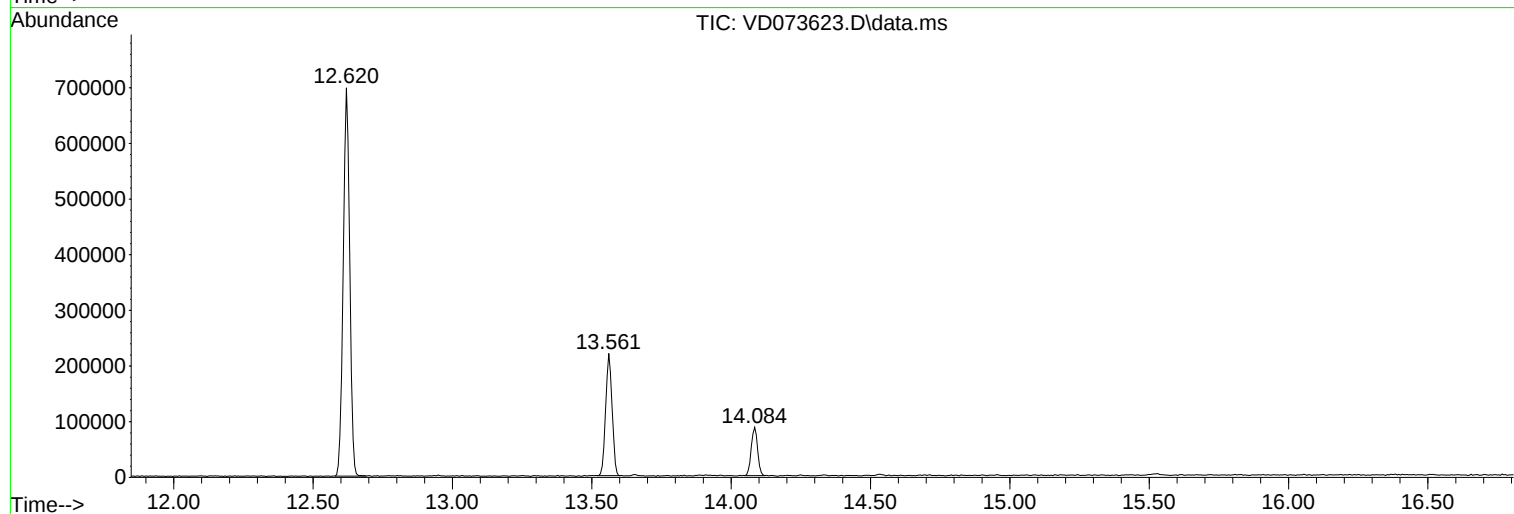
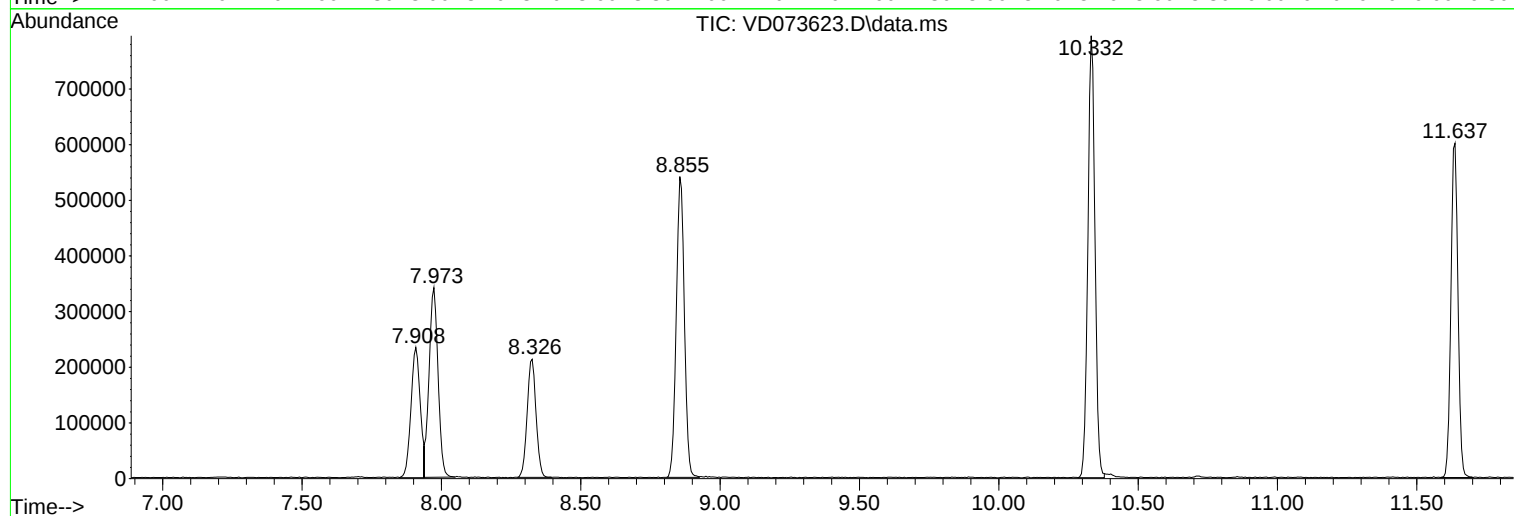
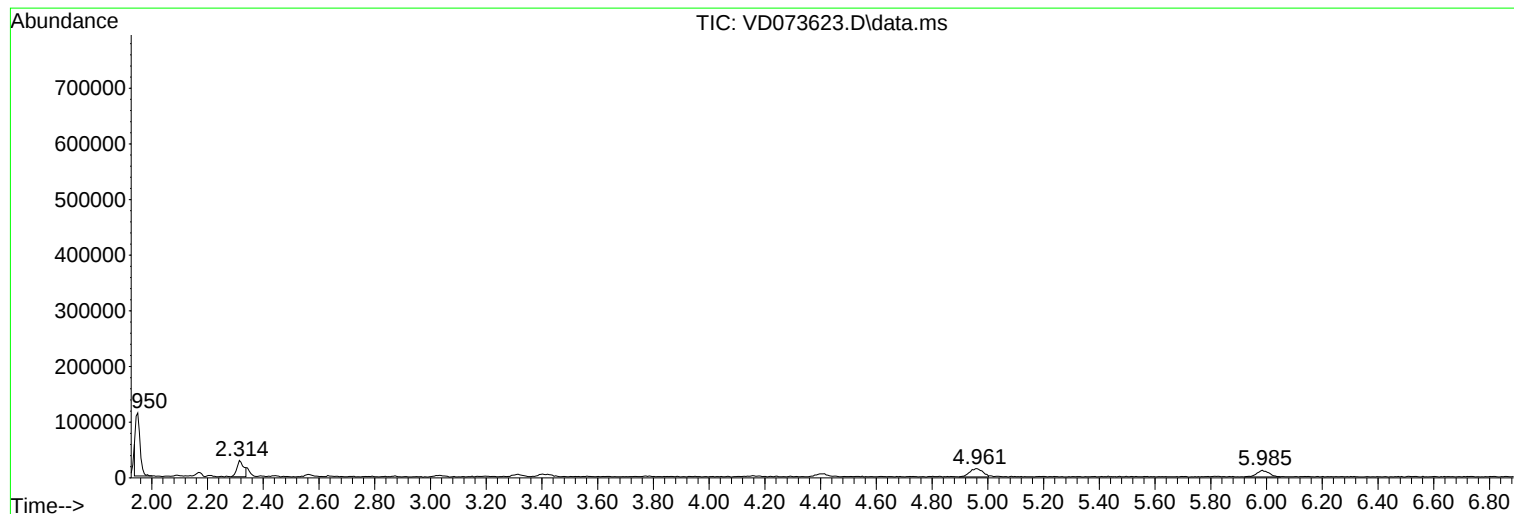
Sum of corrected areas: 7210684

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D060922S.M
Quant Title : SW846 8260

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



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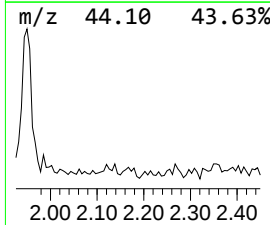
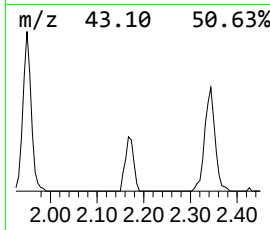
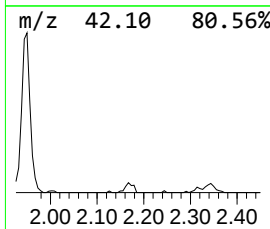
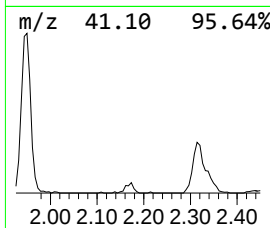
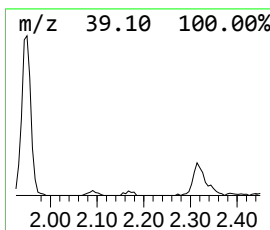
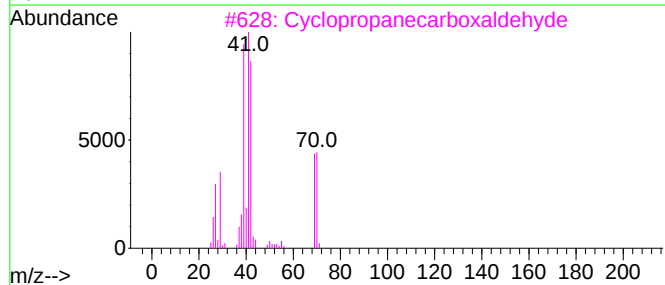
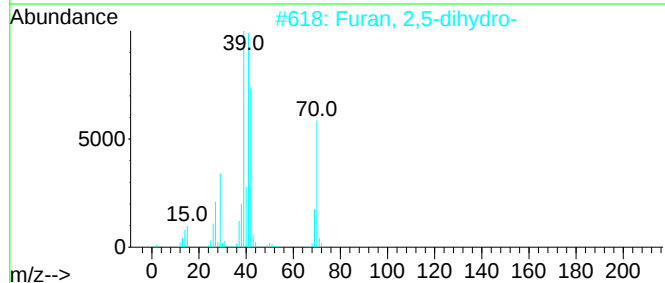
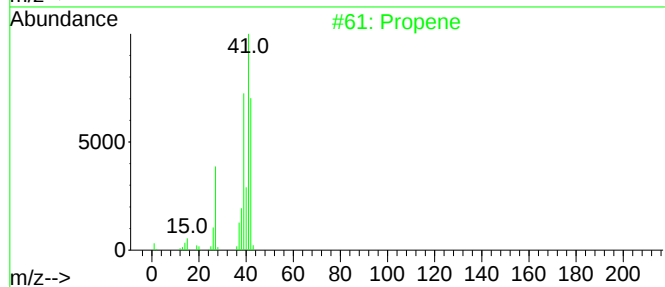
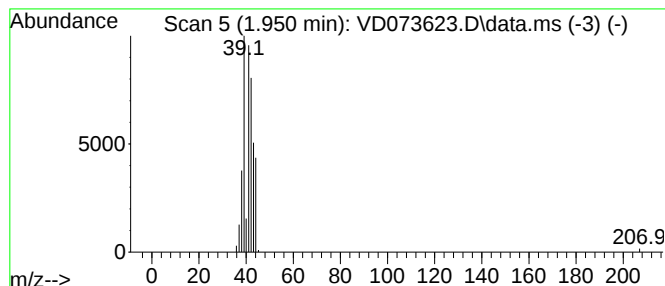
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D060922S.M
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Peak Number 1 Propene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.950	8.16 ug/l	124004	Pentafluorobenzene	7.973

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propene	42	C3H6	000115-07-1	64
2			Furan, 2,5-dihydro-	70	C4H6O	001708-29-8	39
3			Cyclopropanecarboxaldehyde	70	C4H6O	001489-69-6	39
4			N-Methyl-7-azabicyclo(2,2,1)hept...	109	C7H11N	055590-26-6	38
5			Cyclopropane	42	C3H6	000075-19-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propene	1.950	8.2	ug/l	124004	1	7.973	760141	50.0