

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU040221\
 Data File : VU042908.D
 Acq On : 02 Apr 2021 20:12
 Operator : SY/MD
 Sample : M1903-09 10X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 31031

Integration Parameters: RTEINT.P

Integrator: RTE

Smoother: 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_U\Method\82U032421W.M

Title : SW846 8260

Signal : TIC: VU042908.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.549	351	376	394	rBV	1990057	3315112	100.00%	45.833%
2	2.629	395	401	433	rVB	71823	158952	4.79%	2.198%
3	4.063	831	847	876	rBV2	49513	156545	4.72%	2.164%
4	5.298	1216	1231	1245	rBV	102715	217644	6.57%	3.009%
5	5.382	1245	1257	1275	rVB	165796	333173	10.05%	4.606%
6	5.710	1344	1359	1377	rBV	116544	235951	7.12%	3.262%
7	6.256	1515	1529	1554	rBV	238404	448920	13.54%	6.207%
8	7.906	2029	2042	2064	rBV	368302	634030	19.13%	8.766%
9	9.423	2501	2514	2527	rBV	335677	555377	16.75%	7.678%
10	9.497	2527	2537	2558	rVB4	15427	41672	1.26%	0.576%
11	10.639	2880	2892	2905	rBV	281592	453736	13.69%	6.273%
12	10.713	2905	2915	2927	rVB	47386	80607	2.43%	1.114%
13	11.819	3248	3259	3277	rVB	349306	561385	16.93%	7.761%
14	12.069	3326	3337	3356	rBV3	19013	39910	1.20%	0.552%

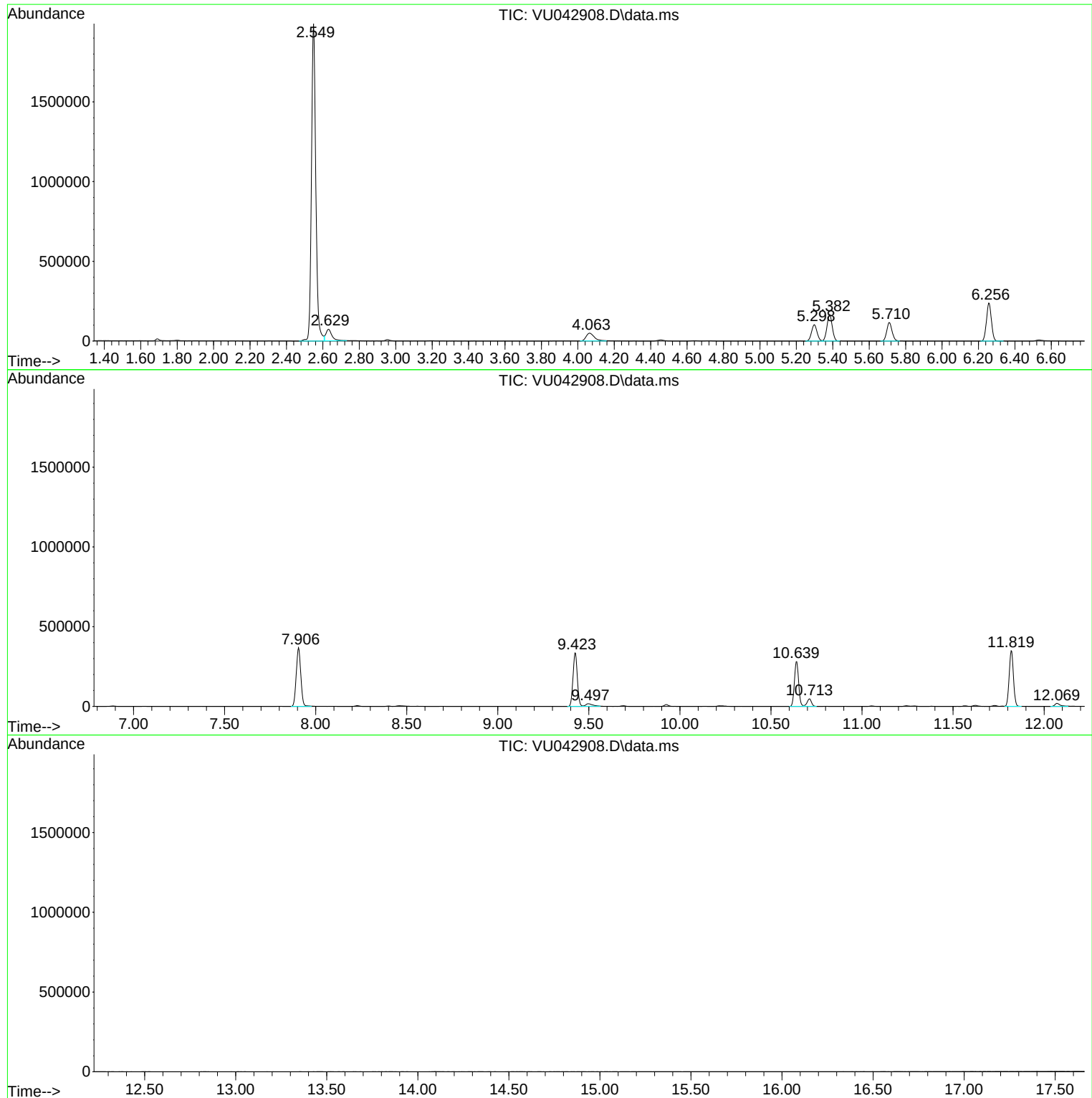
Sum of corrected areas: 7233014

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Quant Title : SW846 8260

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



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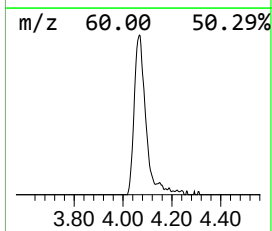
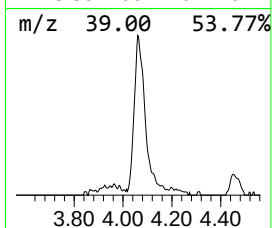
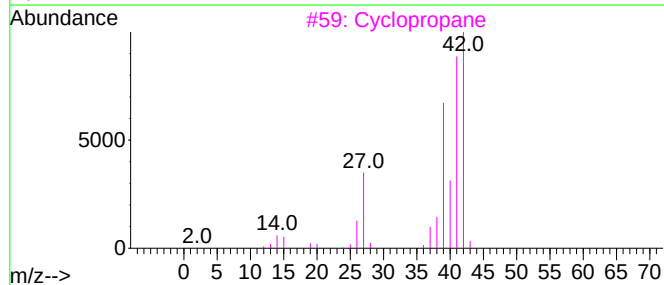
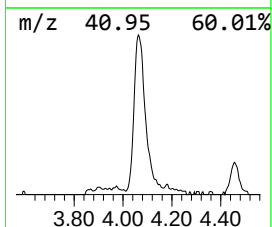
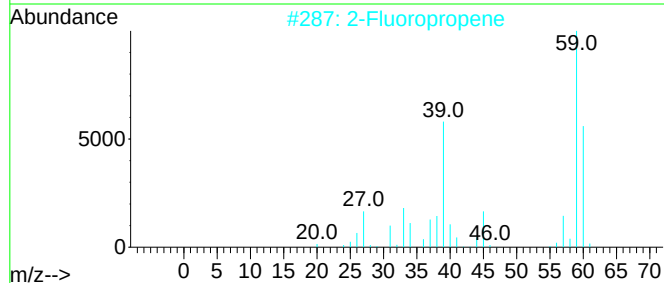
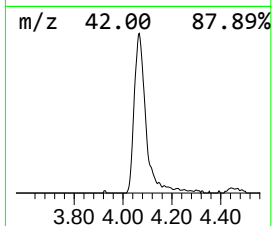
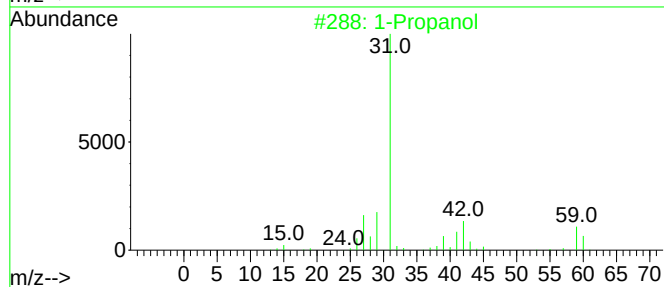
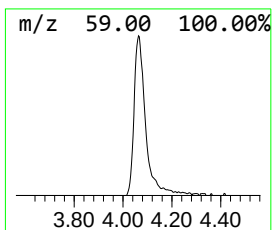
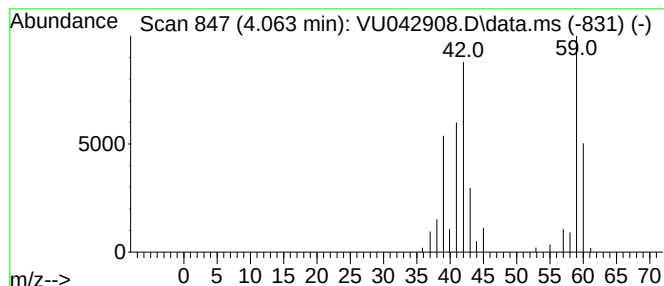
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Peak Number 1 1-Propanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.063	23.49 ug/l	156545	Pentafluorobenzene	5.385

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propanol			60	C3H8O	000071-23-8	52
2	2-Fluoropropene			60	C3H5F	001184-60-7	49
3	Cyclopropane			42	C3H6	000075-19-4	38
4	Allyl fluoride			60	C3H5F	000818-92-8	35
5	Propene			42	C3H6	000115-07-1	23



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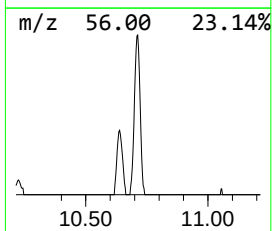
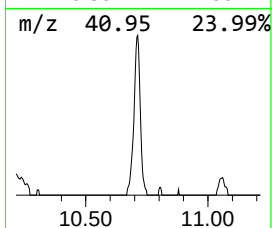
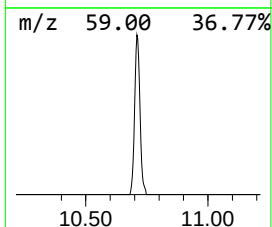
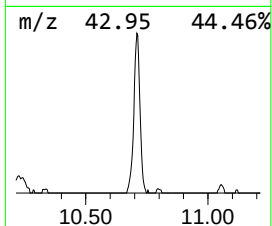
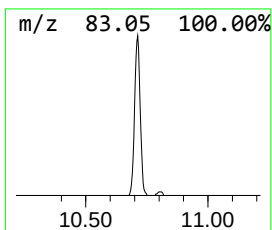
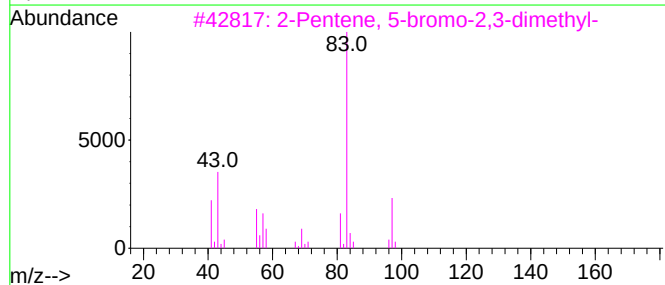
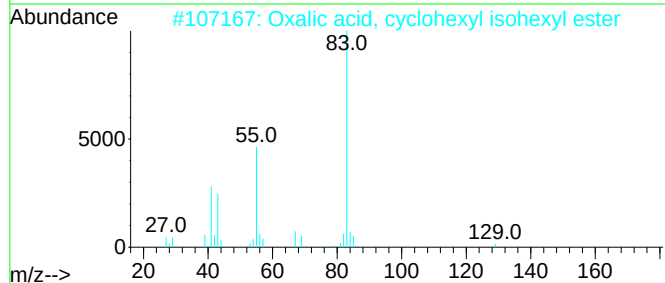
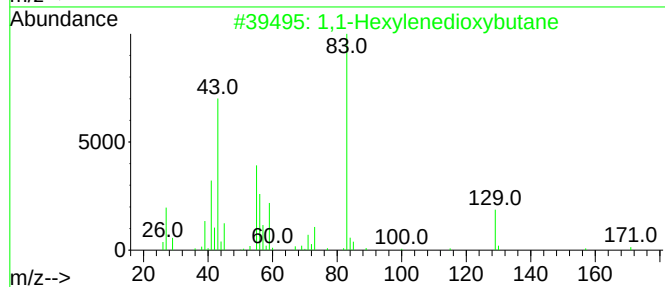
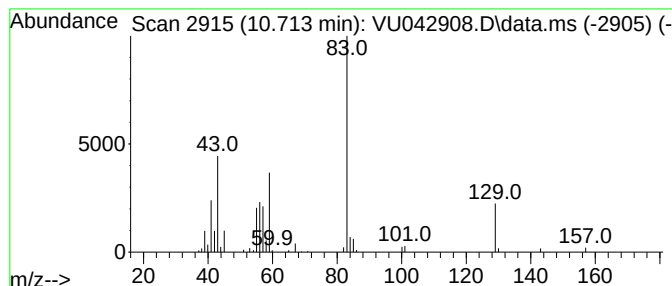
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Peak Number 2 1,1-Hexylenedioxybutane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.713	7.18 ug/l	80607	1,4-Dichlorobenzene-d4	11.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1-Hexylenedioxybutane	172	C10H20O2	1000249-77-4	78
2			Oxalic acid, cyclohexyl isohexyl...	256	C14H24O4	1000309-30-7	37
3			2-Pentene, 5-bromo-2,3-dimethyl-	176	C7H13Br	056312-52-8	25
4			3-Methyl-2-butenic acid, oct-3-...	210	C13H22O2	1000299-31-2	23
5			3-Methylbut-2-enoic acid, 4-nitr...	221	C11H11NO4	1000307-59-8	17



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Propanol	4.063	23.5	ug/l	156545	1	5.385	333173	50.0
1,1-Hexylenedio...	10.713	7.2	ug/l	80607	4	11.822	561385	50.0