

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN083121\
 Data File : VN068373.D
 Acq On : 31 Aug 2021 16:49
 Operator : JC/MD
 Sample : IBLK
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IBLK

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_N\methods\82N083021W.M

Title : SW846 8260

Signal : TIC: VN068373.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.234	77	91	110	rBV	30148	66370	2.94%	0.511%
2	2.741	269	280	282	rBV7	19010	25040	1.11%	0.193%
3	4.151	783	806	831	rBV	112584	318733	14.12%	2.456%
4	6.289	1585	1603	1626	rBV4	9050	28885	1.28%	0.223%
5	6.579	1692	1711	1742	rBV7	30407	103681	4.59%	0.799%
6	8.024	2226	2250	2262	rBV	317491	768266	34.02%	5.920%
7	8.088	2262	2274	2309	rVB2	546436	1244094	55.10%	9.587%
8	8.440	2382	2405	2434	rBV2	292779	674673	29.88%	5.199%
9	8.971	2584	2603	2641	rBV2	765430	1605265	71.09%	12.370%
10	10.443	3134	3152	3182	rBV	1202483	2258034	100.00%	17.400%
11	11.141	3381	3412	3436	rBV	109950	302932	13.42%	2.334%
12	11.747	3612	3638	3668	rBV	1142358	2194810	97.20%	16.913%
13	12.734	3989	4006	4028	rBV	919398	1603079	70.99%	12.353%
14	13.678	4341	4358	4379	rBV	1039932	1783516	78.99%	13.743%

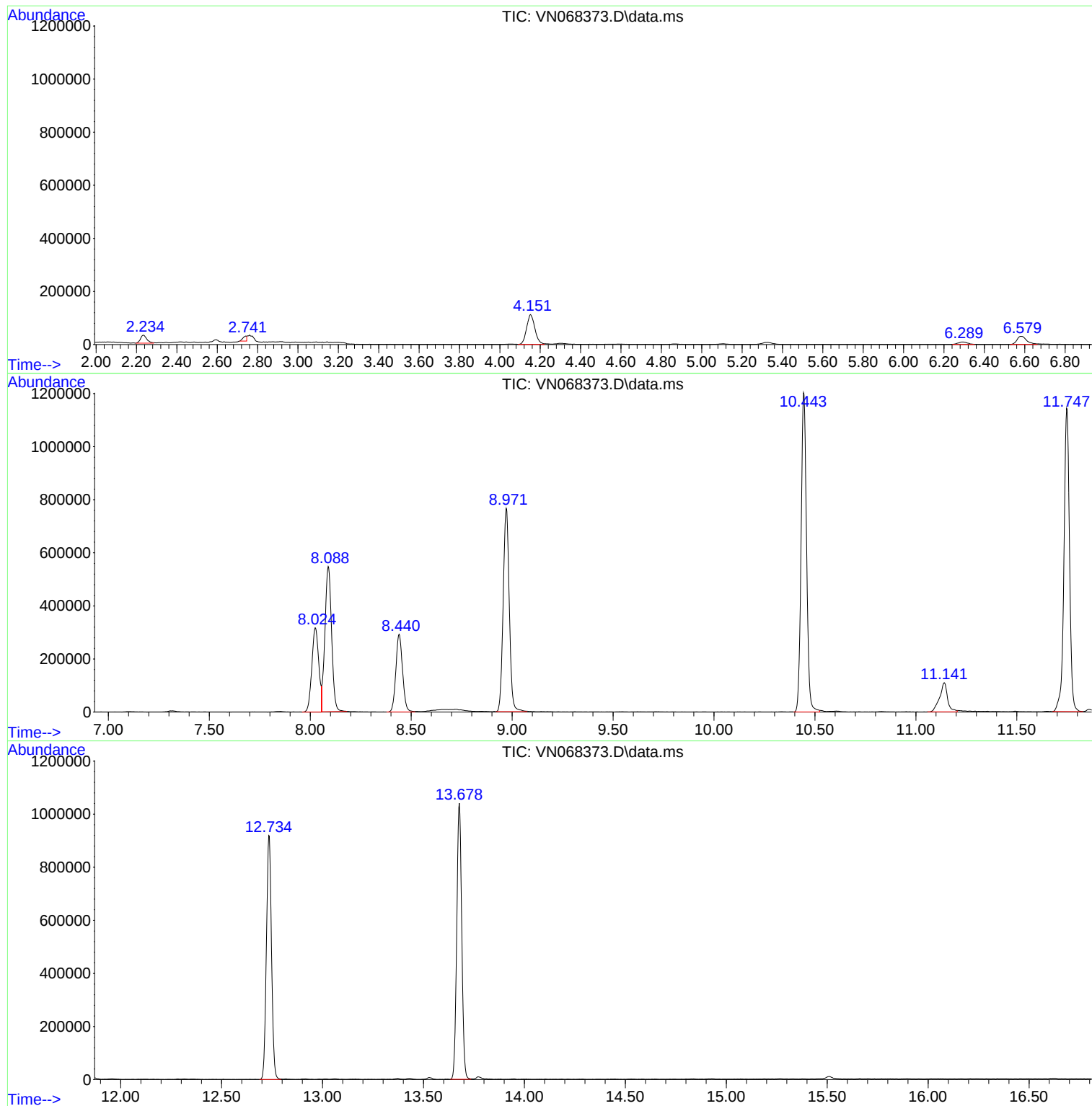
Sum of corrected areas: 12977378

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Quant Title : SW846 8260

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



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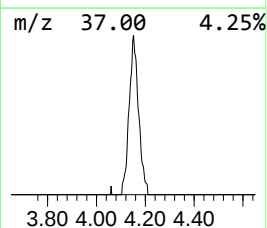
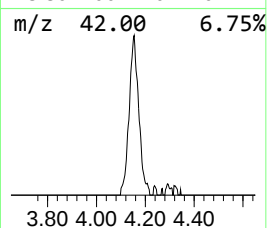
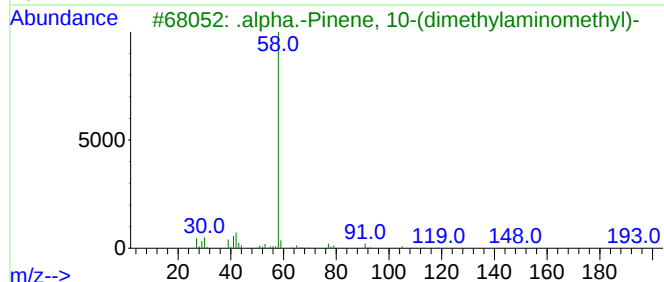
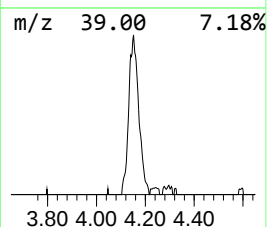
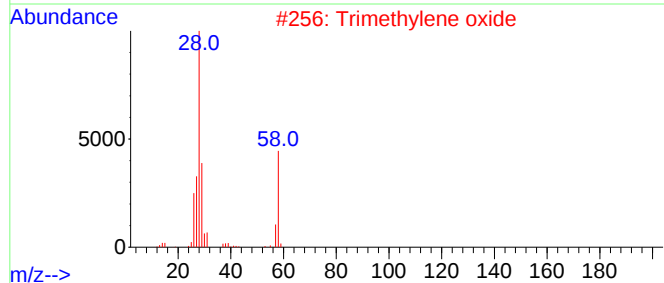
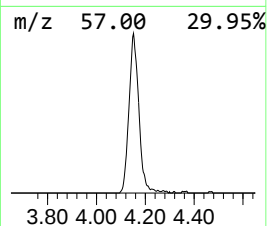
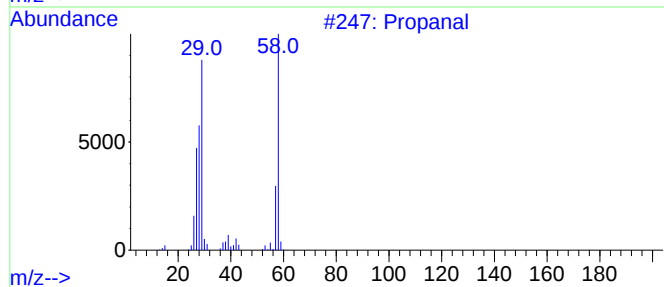
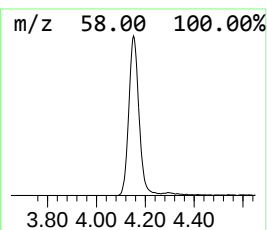
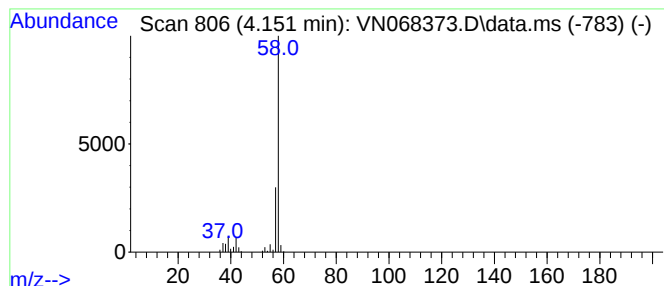
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Peak Number 1 Propanal Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.151	12.81 ug/l	318733	Pentafluorobenzene	8.088

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanal	58	C3H6O	000123-38-6	72
2			Trimethylene oxide	58	C3H6O	000503-30-0	9
3			.alpha.-Pinene, 10-(dimethylamin...	193	C13H23N	015063-97-5	9
4			Succindialdehyde	86	C4H6O2	000638-37-9	9
5			Oxirane, methyl-, (S)-	58	C3H6O	016088-62-3	7



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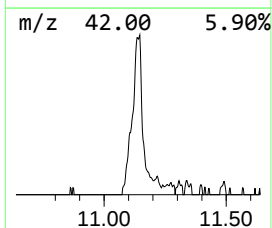
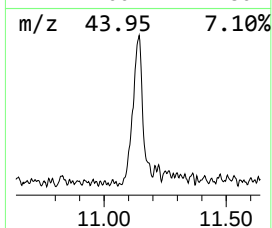
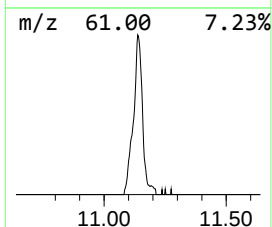
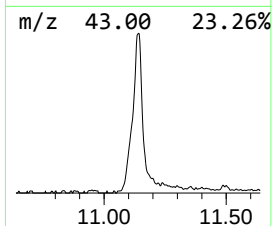
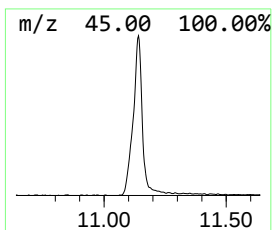
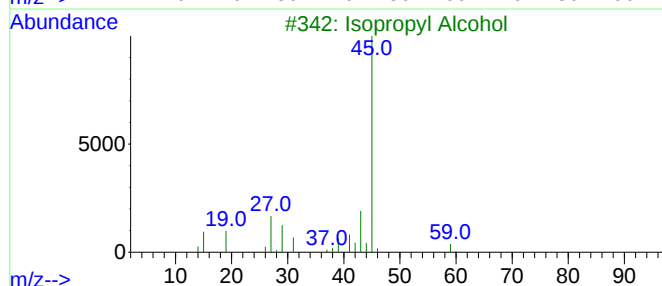
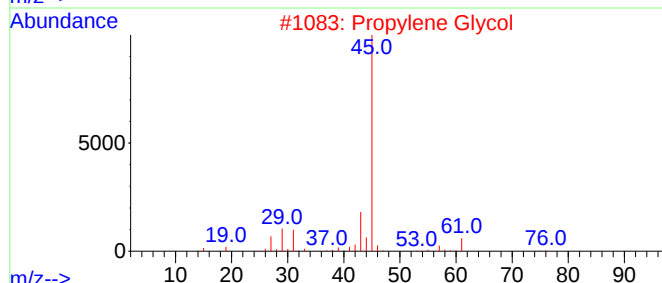
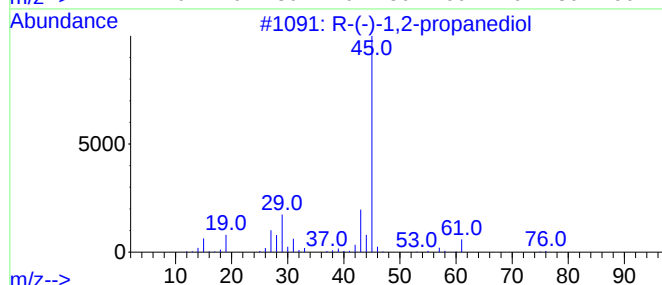
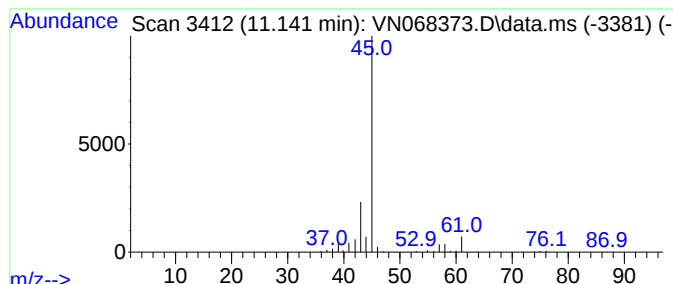
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Peak Number 2 R-(-)-1,2-propanediol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.140	6.90 ug/l	302932	Chlorobenzene-d5	11.749

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
2			Propylene Glycol	76	C3H8O2	000057-55-6	78
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	50
4			2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	40
5			2-Hexanol, 3-methyl-	116	C7H16O	002313-65-7	40



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propanal	4.151	12.8	ug/l	318733	1	8.088	1244090	50.0
R-(-)-1,2-propa...	11.140	6.9	ug/l	302932	3	11.749	2194810	50.0