

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN102119\
 Data File : VN058877.D
 Acq On : 21 Oct 2019 11:05
 Operator : JC/SP
 Sample : K5452-07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FA19-MMW0422

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_N\METHODS\82N101819W.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	4.040	686	705	726	rBV	56340	162365	6.93%	1.324%
2	6.805	1547	1565	1580	rBV	8823	26187	1.12%	0.214%
3	6.911	1580	1598	1625	rVB2	67633	187190	7.99%	1.527%
4	7.574	1783	1804	1815	rBV	301777	756824	32.31%	6.173%
5	7.648	1815	1827	1848	rVB	537004	1285122	54.87%	10.482%
6	8.011	1922	1940	1965	rBV	370883	870300	37.16%	7.098%
7	8.570	2097	2114	2139	rBV	793889	1653730	70.61%	13.488%
8	10.085	2569	2585	2613	rBV	1278230	2342091	100.00%	19.102%
9	10.278	2631	2645	2648	rBV	30237	64188	2.74%	0.524%
10	10.628	2743	2754	2764	rBV3	37700	66042	2.82%	0.539%
11	11.406	2980	2996	3022	rBV	1083710	1887610	80.60%	15.396%
12	12.403	3289	3306	3320	rBV	823003	1398423	59.71%	11.406%
13	13.348	3587	3600	3624	rBV	867249	1446886	61.78%	11.801%
14	16.017	4424	4430	4443	rVB	62539	113819	4.86%	0.928%

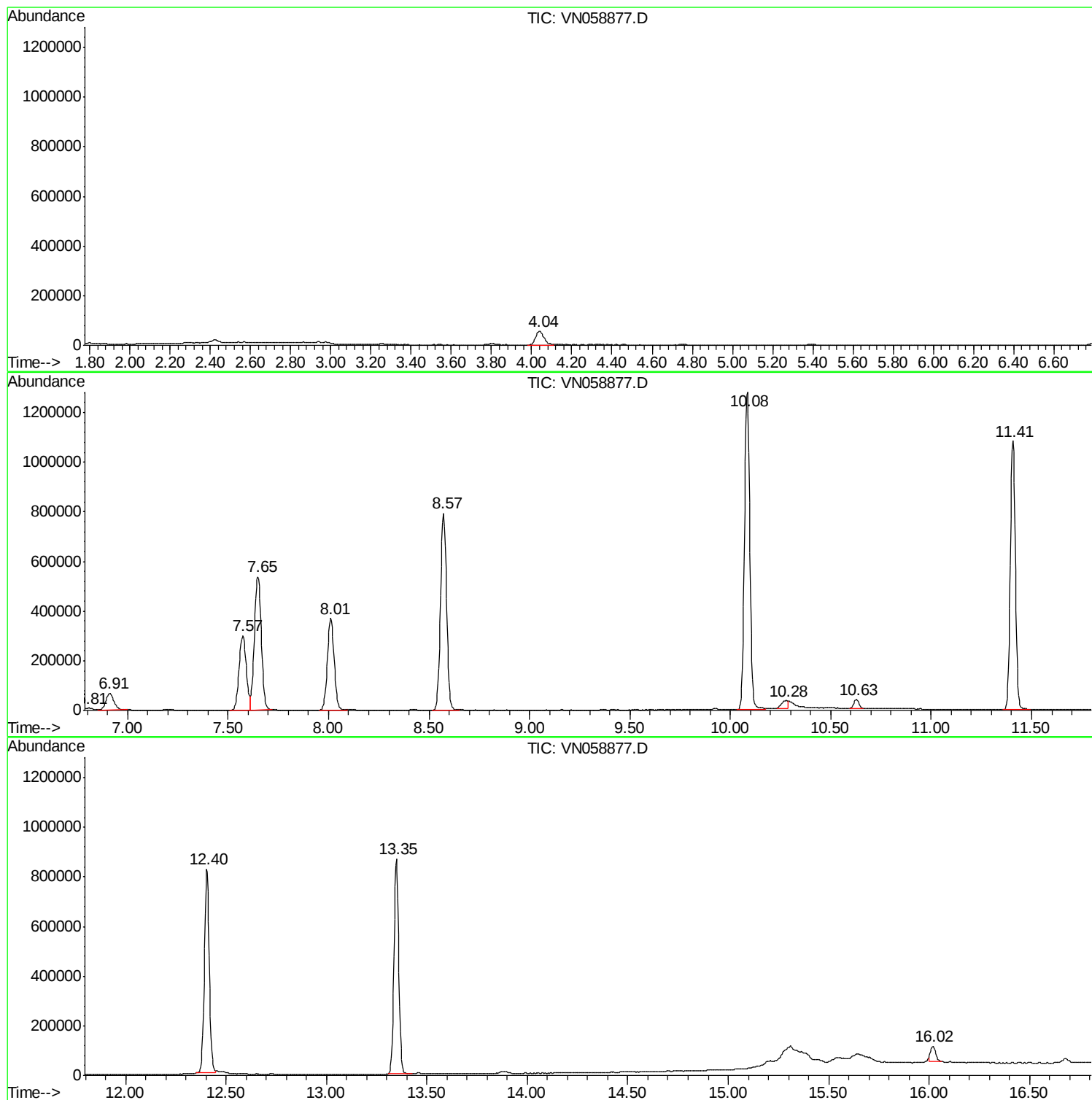
Sum of corrected areas: 12260777

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N101819W.M
Quant Title : SW846 8260

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



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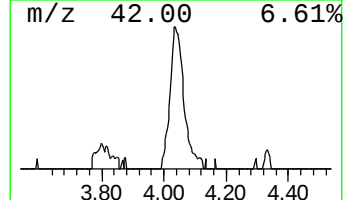
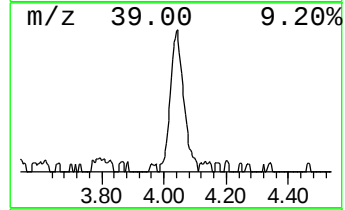
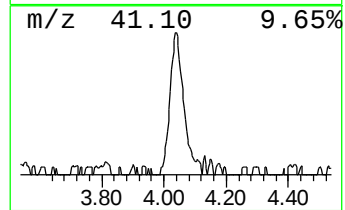
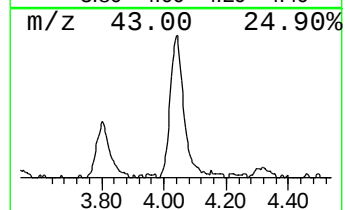
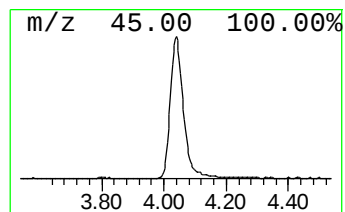
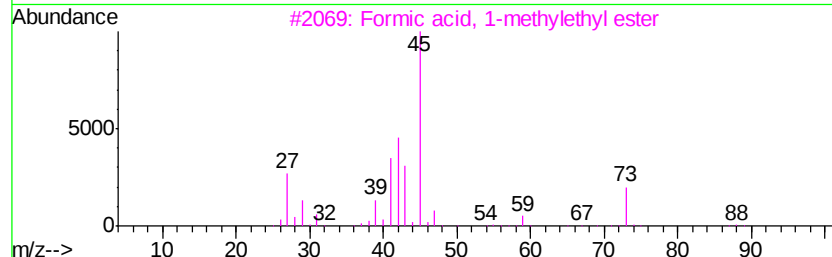
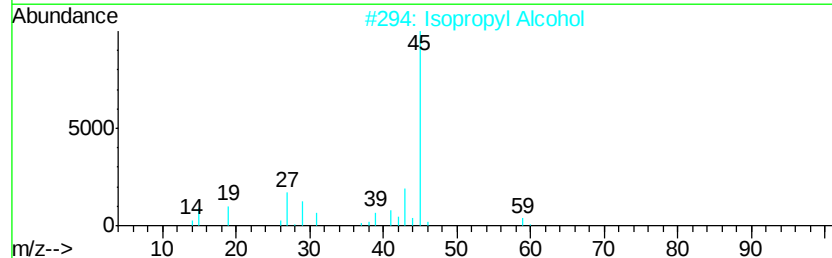
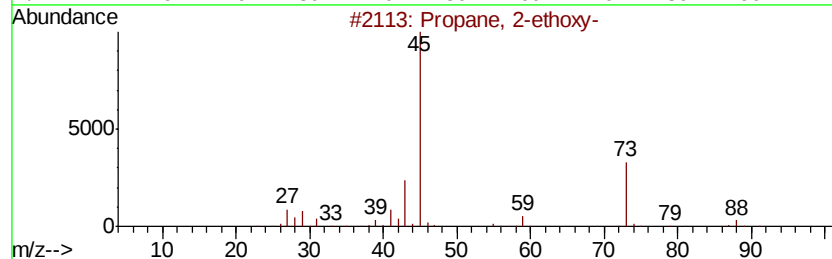
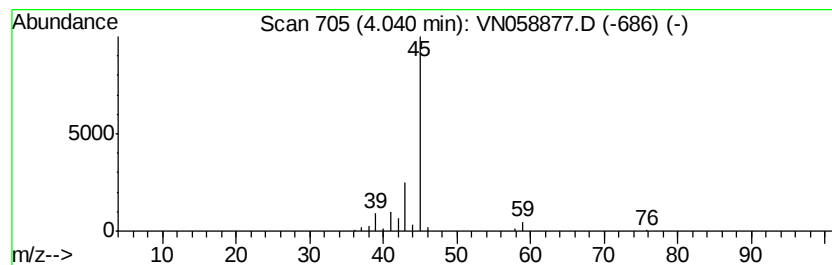
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Peak Number 1 unknown4.04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.04	6.32 ug/l	162365	Pentafluorobenzene	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2-ethoxy-	88	C5H12O	000625-54-7	43
2		Isopropyl Alcohol	60	C3H8O	000067-63-0	43
3		Formic acid, 1-methylethyl ester	88	C4H8O2	000625-55-8	43
4		R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	9
5		2-Pentanol	88	C5H12O	006032-29-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown4.04	4.04	6.3	ug/l	162365	1	7.65	1285120	50.0