

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX030818\
Data File : VX000219.D
Acq On : 08 Mar 2018 12:56
Operator : JC/MD
Sample : J1833-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TB-002

Integration Parameters: RTEINT3.P

Integrator: RTE
Smoothing : OFF
Sampling : 1
Start Thrs: 0.2
Stop Thrs : 0
Filtering: 5
Min Area: 0 % of largest Peak
Max Peaks: 100
Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : W:\HPCHEM1\MSVOA_X\METHOD\624X022018W.M
Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.587	159	164	171	rVB5	3958	7743	1.78%	0.432%
2	2.617	324	333	340	rVB	11025	19786	4.55%	1.103%
3	5.032	718	729	743	rBV3	52480	159420	36.63%	8.884%
4	6.086	890	902	914	rBV	60354	156280	35.91%	8.709%
5	6.873	1022	1031	1043	rBV2	133946	310573	71.36%	17.308%
6	8.726	1328	1335	1344	rBV	269796	435191	100.00%	24.253%
7	10.122	1558	1564	1570	rBV	294749	389476	89.50%	21.705%
8	11.146	1726	1732	1742	rVB	260785	315920	72.59%	17.606%

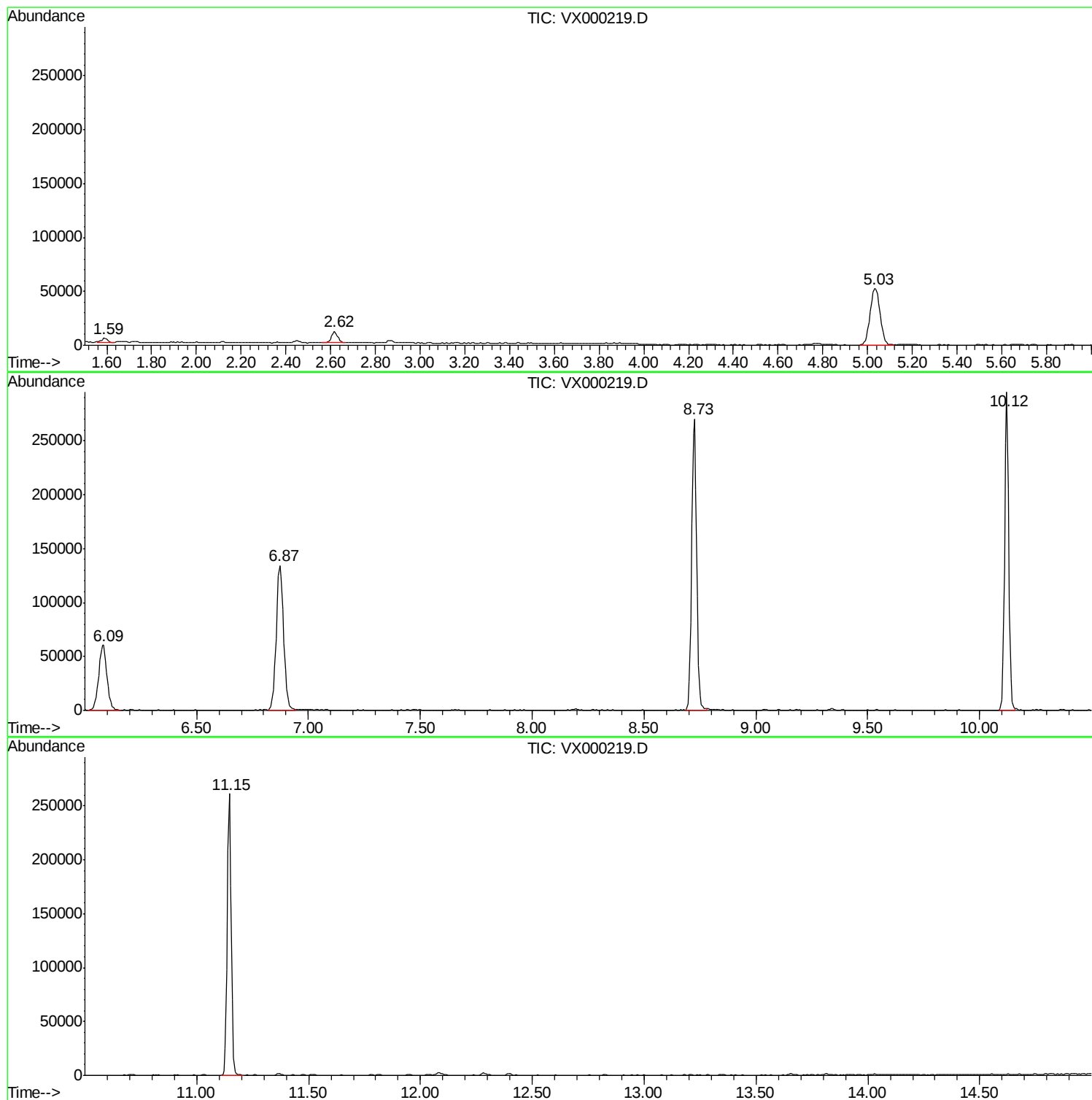
Sum of corrected areas: 1794389

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Quant Method : W:\HPCHEM1\MSVOA X\METHOD\624X022018W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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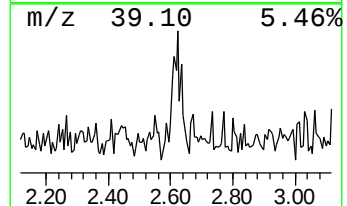
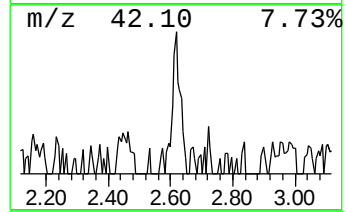
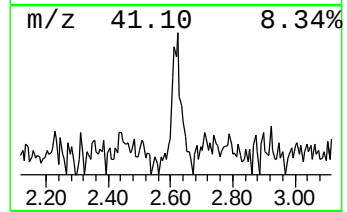
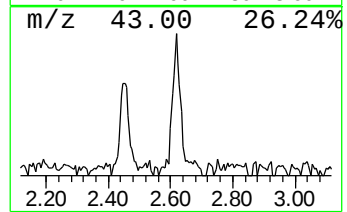
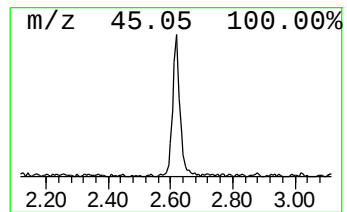
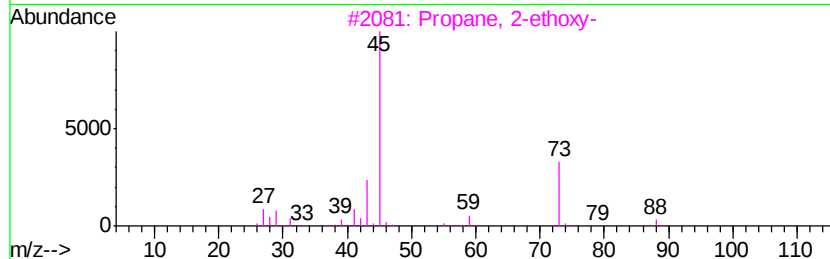
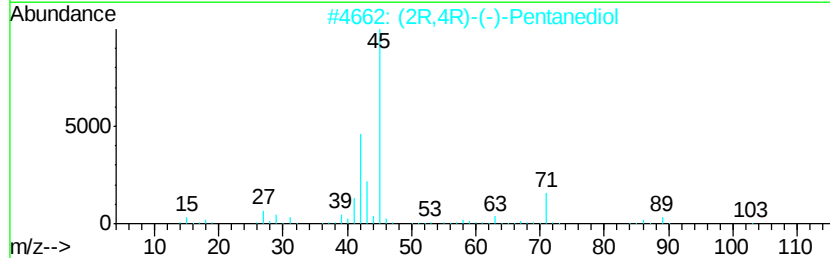
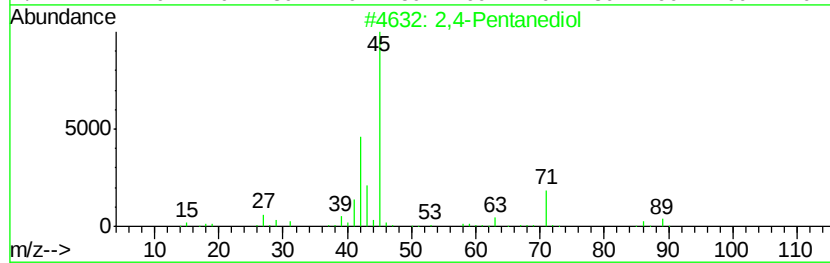
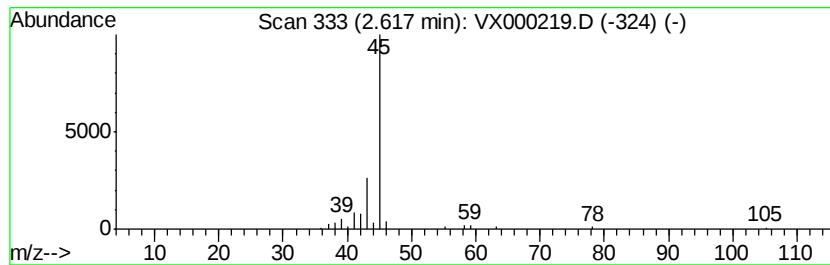
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Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown2.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.62	3.72 ug/l	19786	Bromochloromethane	5.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4-Pentanediol	104	C5H12O2	000625-69-4	40
2		(2R,4R)-(-)-Pentanediol	104	C5H12O2	042075-32-1	40
3		Propane, 2-ethoxy-	88	C5H12O	000625-54-7	40
4		Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9
5		2-Ethoxypentane	116	C7H16O	001817-89-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.62	2.62	3.7	ug/l	19786	1	5.04	159420	30.0