

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121619\
Data File : VX014029.D
Acq On : 16 Dec 2019 12:54
Operator : JC/SP
Sample : K6295-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
STORAGE-BLANK-WATER-REF-4

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_X\METHOD\82X121319W.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	1.070	26	30	43	rBV	626699	776212	28.34%	5.062%
2	1.265	56	62	70	rBV	103587	143614	5.24%	0.937%
3	1.613	106	119	123	rBV8	48787	141781	5.18%	0.925%
4	2.594	272	280	286	rBV	94601	170508	6.22%	1.112%
5	5.490	744	755	771	rBV2	306127	876569	32.00%	5.717%
6	5.649	771	781	794	rVB	567053	1558606	56.90%	10.164%
7	6.051	837	847	862	rBV	352943	916355	33.45%	5.976%
8	6.850	966	978	990	rBV	874289	1995189	72.84%	13.012%
9	8.709	1276	1283	1293	rBV	1791255	2739158	100.00%	17.863%
10	10.105	1505	1512	1523	rBV	1649439	2319131	84.67%	15.124%
11	11.135	1674	1681	1687	rBV	1330765	1726368	63.03%	11.258%
12	12.074	1829	1835	1843	rBV	1619702	1970515	71.94%	12.851%

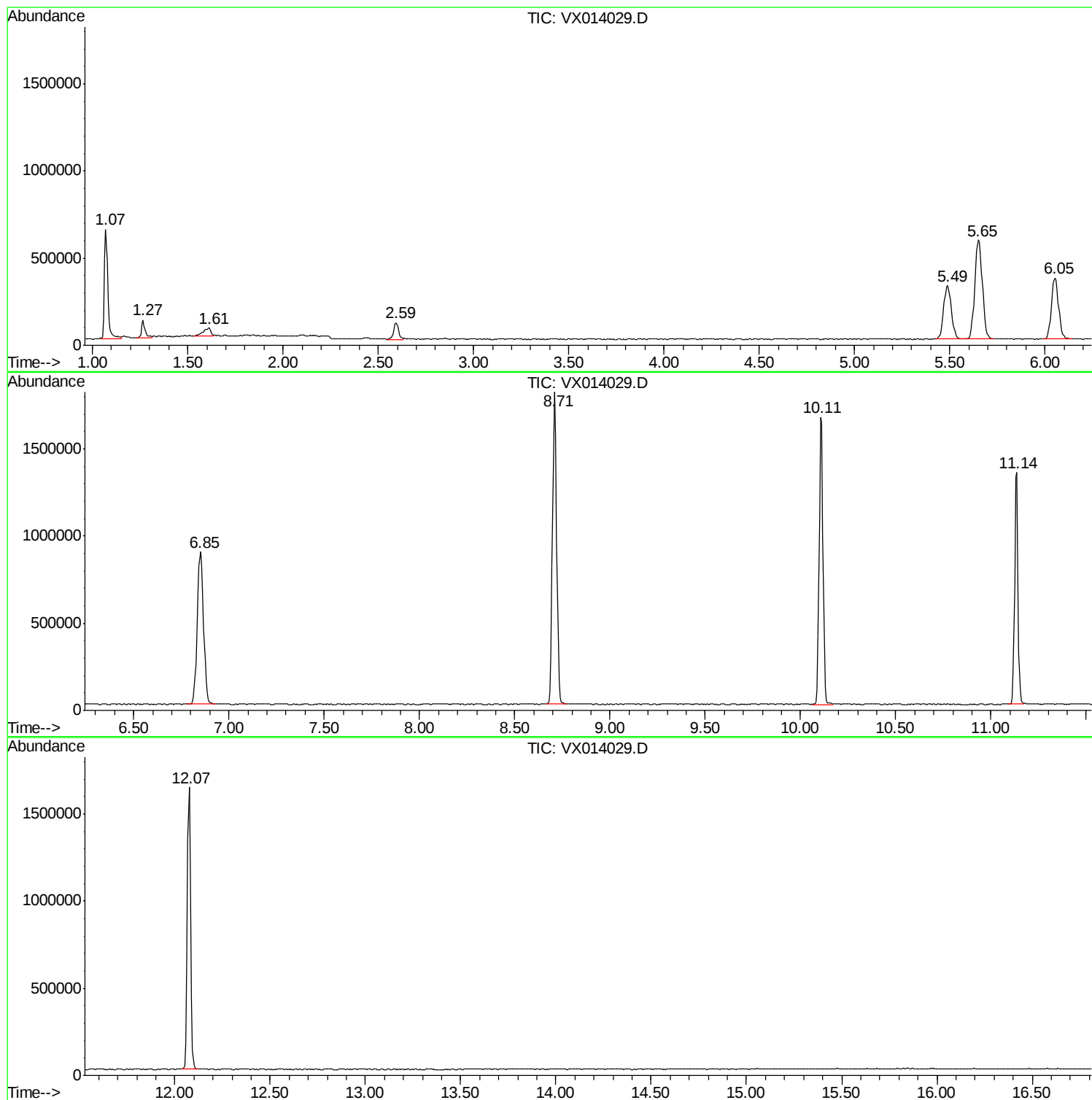
Sum of corrected areas: 15334006

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
Quant Title : SW846 8260

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



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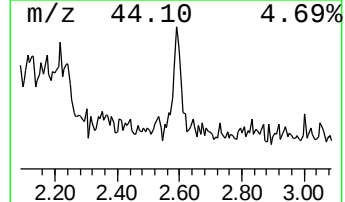
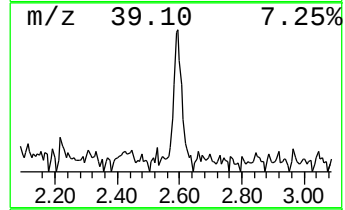
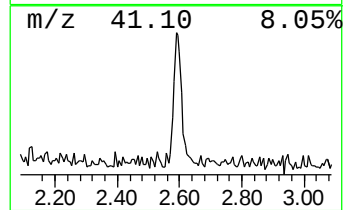
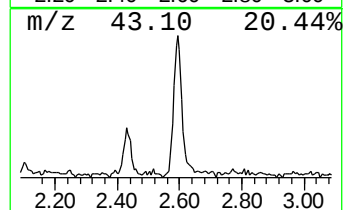
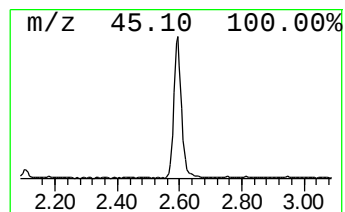
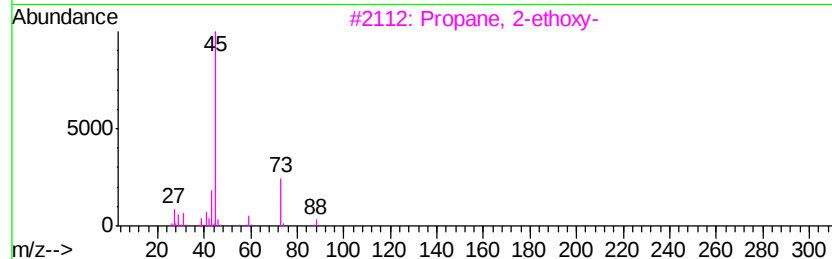
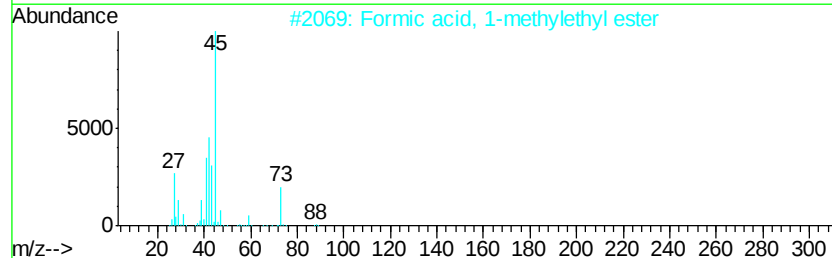
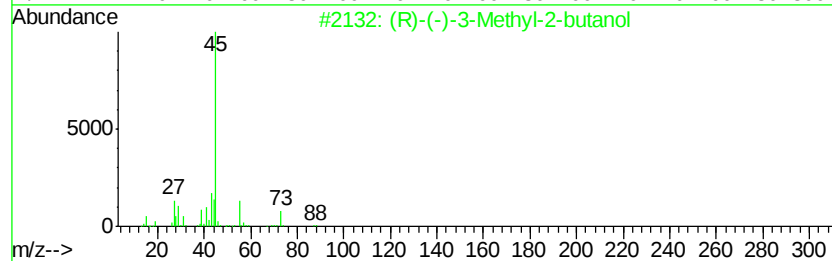
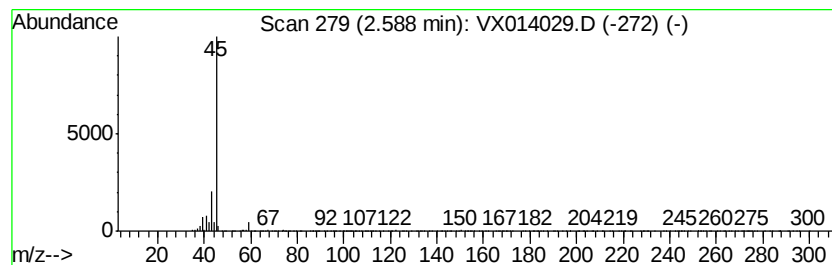
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
Quant Title : SW846 8260

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

Peak Number 2 (R)-(-)-3-Methyl-2-butanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.59	5.47 ug/l	170508	Pentafluorobenzene	5.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(R)-(-)-3-Methyl-2-butanol	88	C5H12O	001572-93-6	64
2			Formic acid, 1-methylethyl ester	88	C4H8O2	000625-55-8	64
3			Propane, 2-ethoxy-	88	C5H12O	000625-54-7	56
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
5			2-Butanol, 3-methyl-, (S)-	88	C5H12O	001517-66-4	39



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(R)-(-)-3-Methyl-...	2.59	5.5	ug/l	170508	1	5.65	1558610	50.0