

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX050919\
 Data File : VX009429.D
 Acq On : 09 May 2019 15:48
 Operator : JC/SP
 Sample : K2748-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-11

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\82X042919W.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	1.605	65	74	78	rBV7	6008	13908	1.40%	0.273%
2	2.605	230	238	249	rBV	10964	22129	2.23%	0.435%
3	5.501	702	713	727	rBV	106495	304106	30.59%	5.974%
4	5.665	727	740	752	rVV	186758	531469	53.46%	10.441%
5	6.061	794	805	820	rBV	131527	346342	34.84%	6.804%
6	6.860	923	936	951	rBV	302333	716320	72.05%	14.072%
7	8.713	1231	1240	1259	rBV	648309	994151	100.00%	19.530%
8	10.109	1463	1469	1477	rBV	592622	800030	80.47%	15.717%
9	11.134	1630	1637	1646	rBV	471840	603088	60.66%	11.848%
10	12.079	1786	1792	1801	rVB	531545	689116	69.32%	13.538%
11	12.268	1817	1823	1834	rBV	45394	69719	7.01%	1.370%

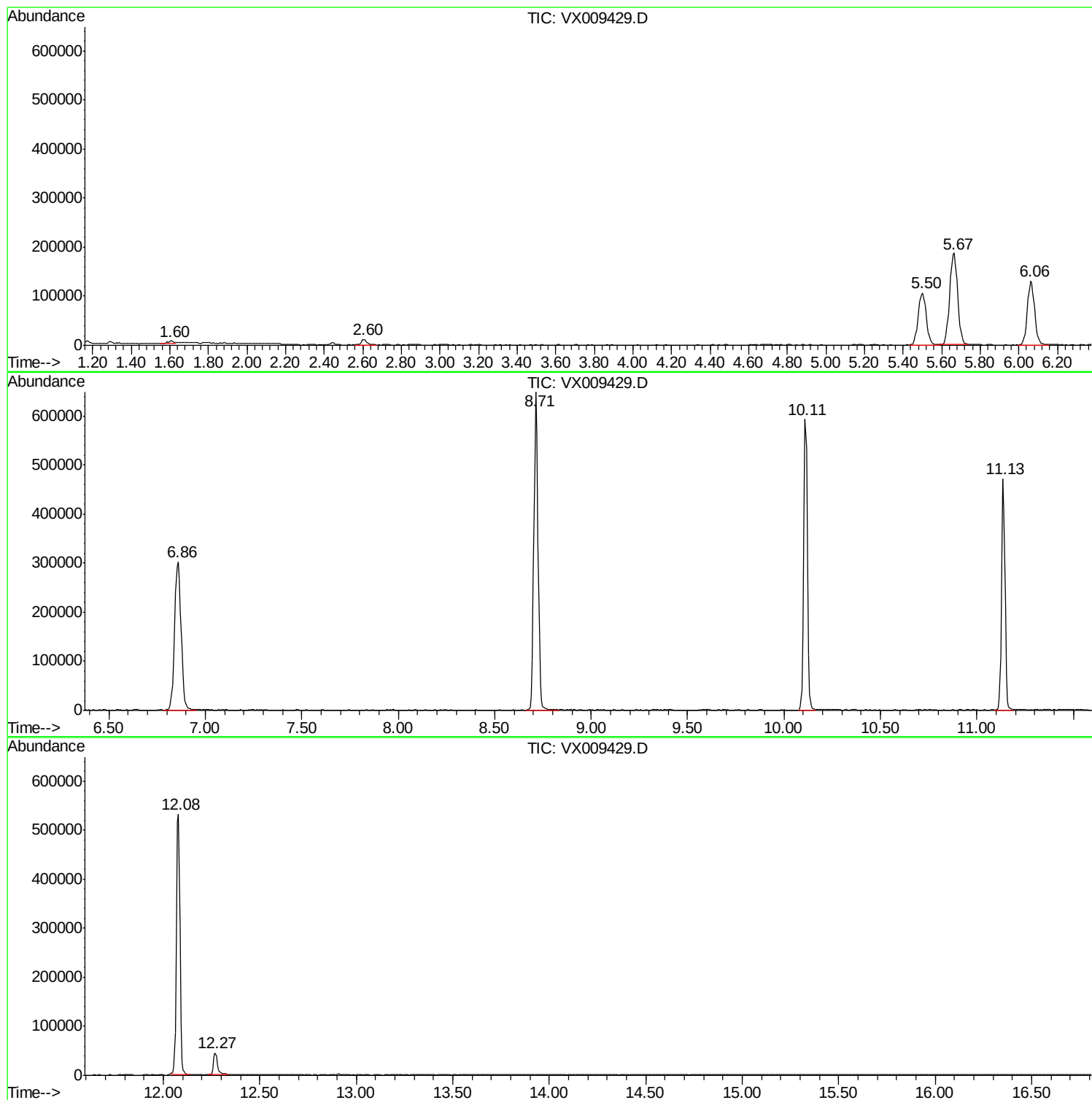
Sum of corrected areas: 5090378

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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



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Instrument :
 MSVOA_X
 ClientSampled :
 MW-11

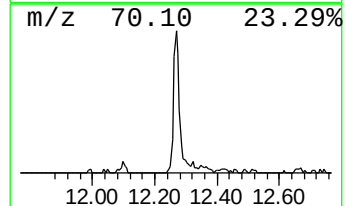
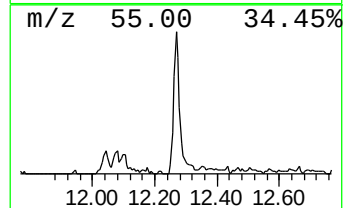
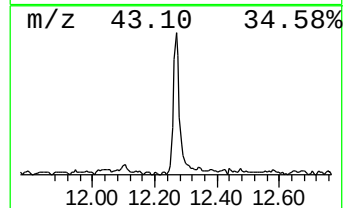
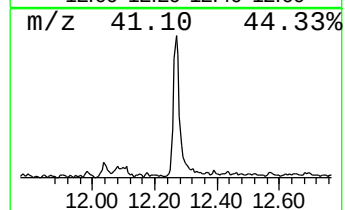
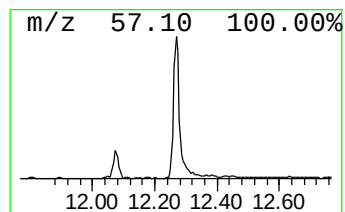
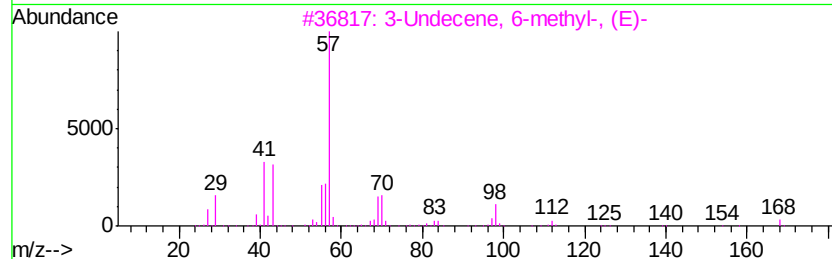
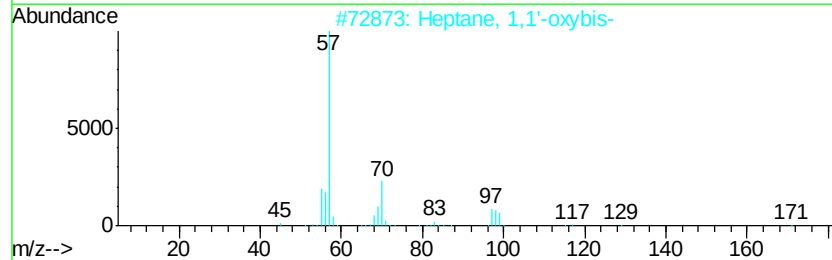
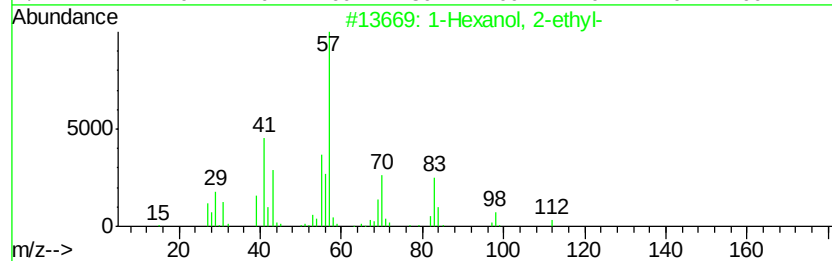
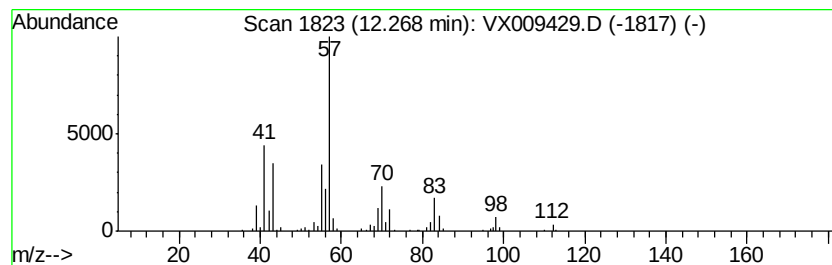
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
 Quant Title : SW846 8260

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

 Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	5.06 ug/l	69719	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2		Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	59
3		3-Undecene, 6-methyl-, (E)-	168	C12H24	074630-52-7	50
4		1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	47
5		Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	45



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Hexanol, 2-ethyl-	12.27	5.1	ug/l	69719	4	12.08	689116	50.0