

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX072518\
 Data File : VX003664.D
 Acq On : 25 Jul 2018 18:10
 Operator : JC/SP
 Sample : J4101-03
 Misc : 3.91g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP-2

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\82X072418W.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	1.069	75	79	93	rBV	413584	448313	80.92%	12.998%
2	1.636	156	172	173	rBV3	6635	22160	4.00%	0.642%
3	2.307	268	282	290	rBV	199373	271124	48.94%	7.861%
4	5.507	797	807	820	rVB	50241	141825	25.60%	4.112%
5	5.660	820	832	849	rBV	77277	215245	38.85%	6.241%
6	6.068	890	899	916	rVB	58557	151504	27.35%	4.393%
7	6.861	1018	1029	1042	rBV	138369	301049	54.34%	8.728%
8	8.714	1327	1333	1349	rBV	345235	554016	100.00%	16.062%
9	9.336	1428	1435	1440	rBV	23943	44836	8.09%	1.300%
10	10.116	1557	1563	1578	rBV	259555	364992	65.88%	10.582%
11	11.140	1726	1731	1747	rVB	259086	348053	62.82%	10.091%
12	11.372	1763	1769	1781	rVB	18549	41694	7.53%	1.209%
13	12.079	1880	1885	1898	rBV	357433	447454	80.77%	12.973%
14	12.274	1913	1917	1931	rBV	62142	79106	14.28%	2.293%
15	12.823	2000	2007	2012	rBV2	7751	11590	2.09%	0.336%
16	14.189	2221	2231	2236	rBV	4472	6189	1.12%	0.179%

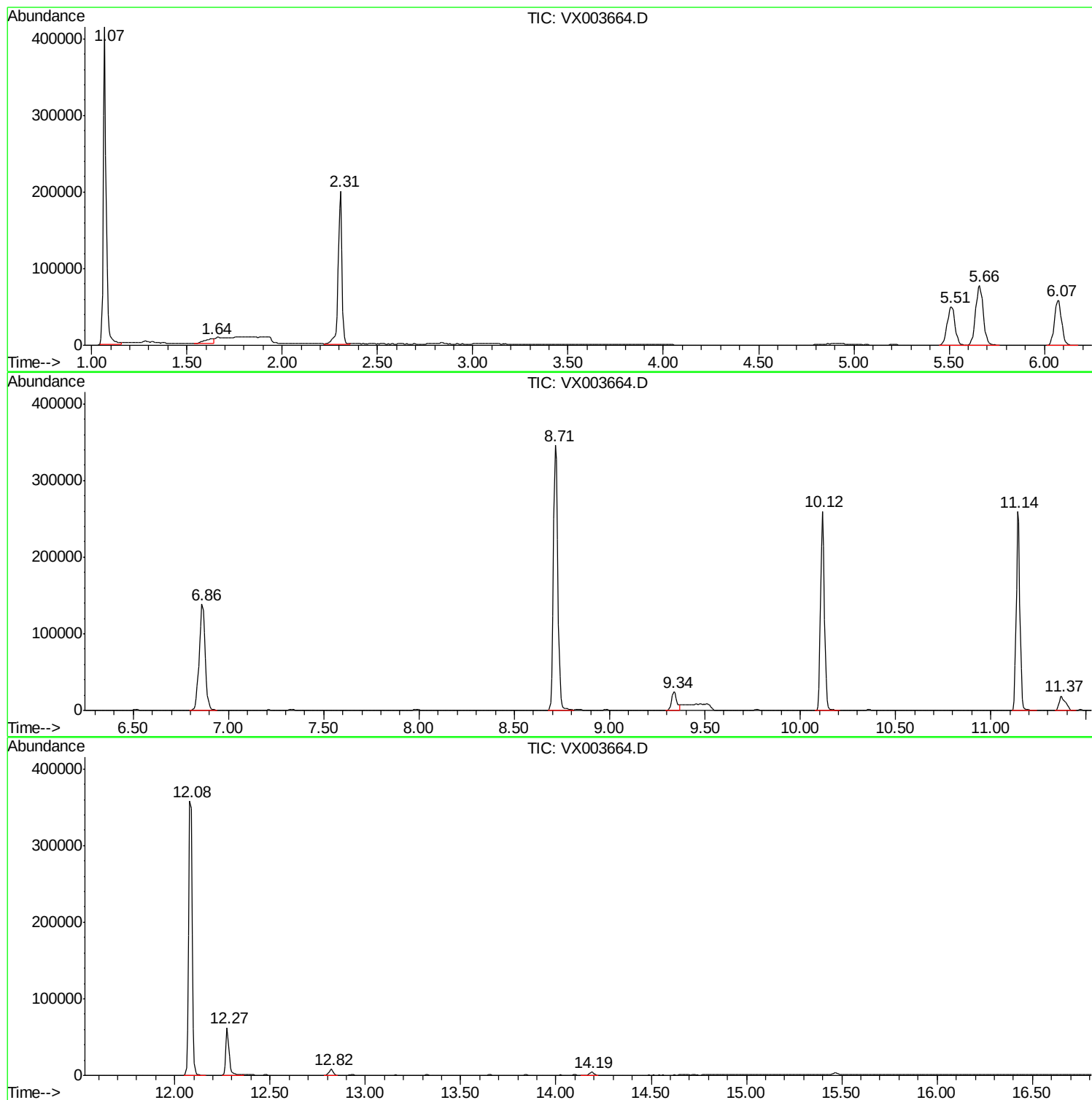
Sum of corrected areas: 3449150

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Quant Title : SW846 8260

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



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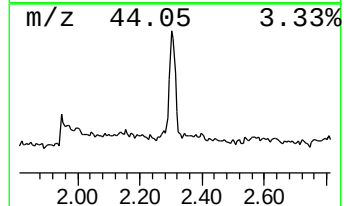
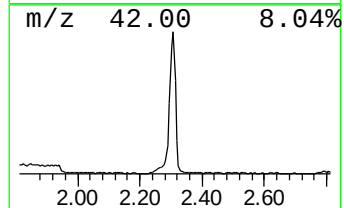
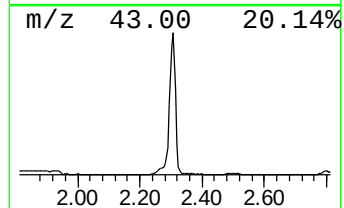
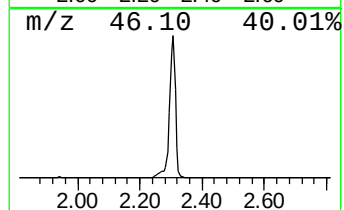
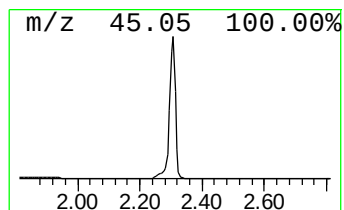
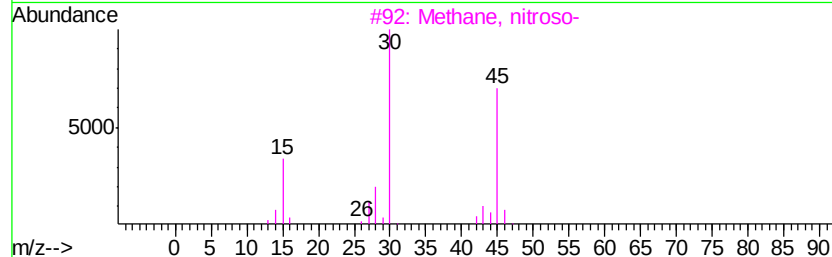
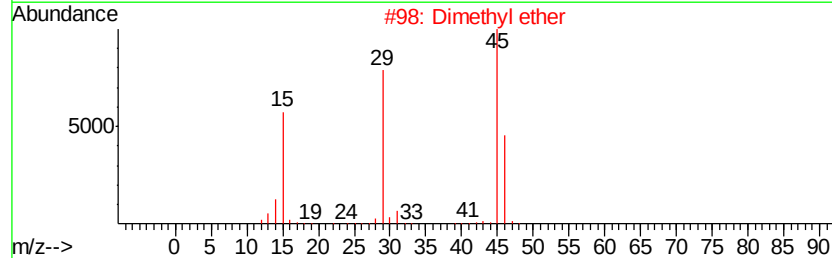
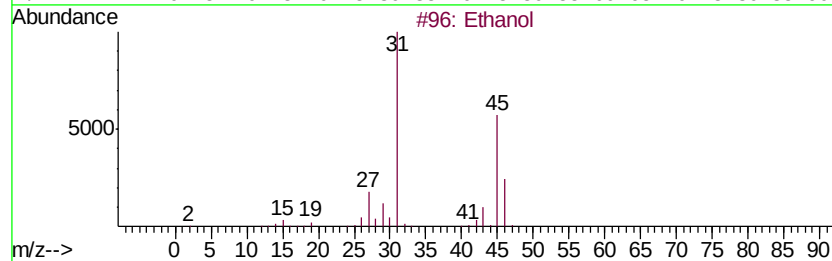
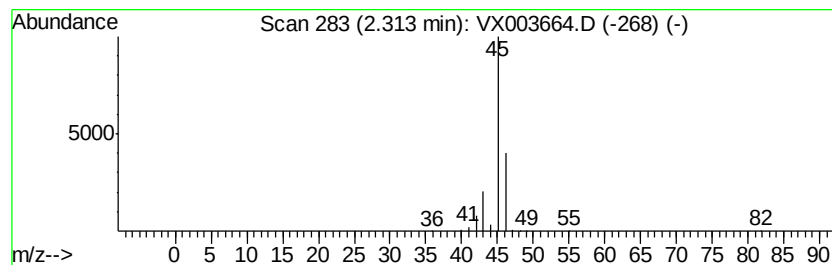
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Peak Number 3 Ethanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.31	62.98 ug/l	271124	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol	46	C2H6O	000064-17-5	74
2		Dimethyl ether	46	C2H6O	000115-10-6	9
3		Methane, nitroso-	45	CH3NO	000865-40-7	4
4		Formic acid	46	CH2O2	000064-18-6	4
5		Oxalic acid	90	C2H2O4	000144-62-7	4



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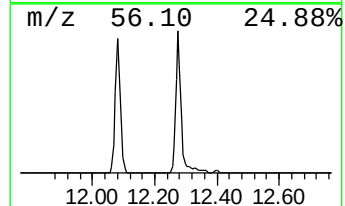
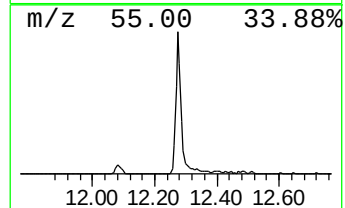
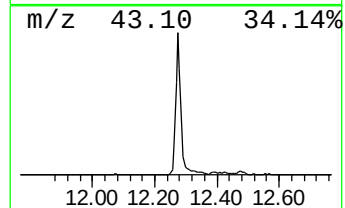
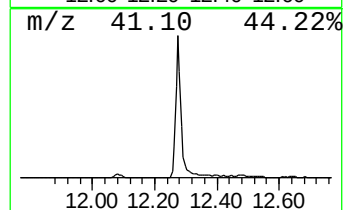
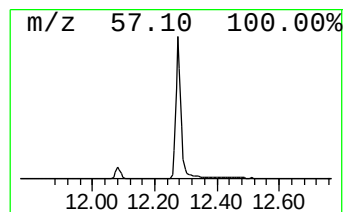
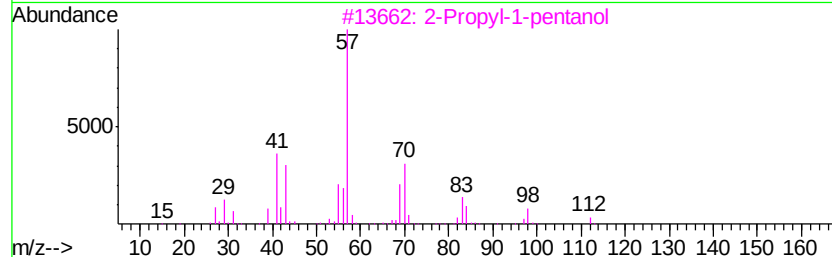
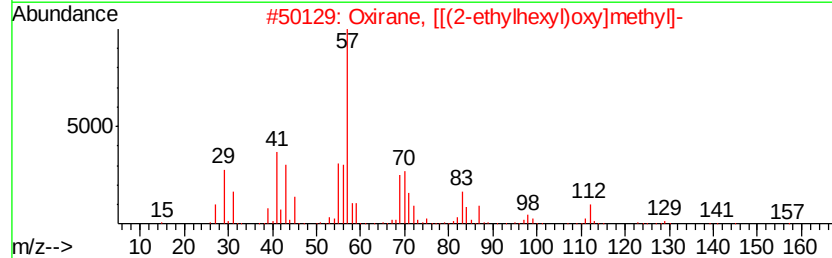
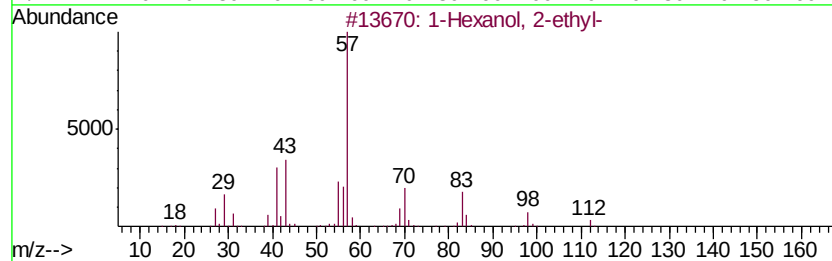
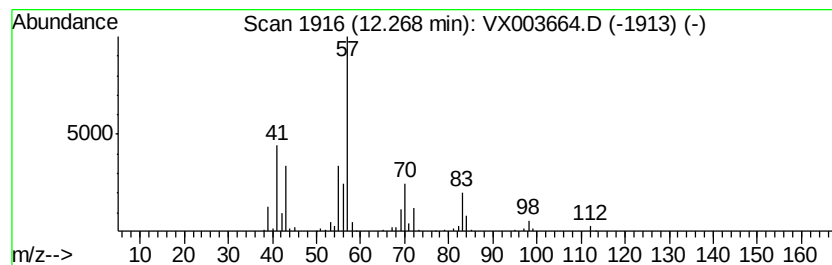
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Peak Number 5 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	8.84 ug/l	79106	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Hexanol, 2-ethyl-	130 C8H18O	000104-76-7 78
2		Oxirane, [[(2-ethylhexyl)oxy]met...	186 C11H22O2	002461-15-6 56
3		2-Propyl-1-pentanol	130 C8H18O	058175-57-8 47
4		Heptane, 1,1'-oxybis-	214 C14H30O	000629-64-1 43
5		Carbonic acid, isobutyl 2-ethylh...	230 C13H26O3	1000357-82-9 38



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol	2.31	63.0	ug/l	271124	1	5.66	215245	50.0
1-Hexanol, 2-ethyl-	12.27	8.8	ug/l	79106	4	12.09	447454	50.0