

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052618\
 Data File : VX002033.D
 Acq On : 26 May 2018 21:39
 Operator : JC/MD
 Sample : J3118-16
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SS052MW732R-180521

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 : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.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	1.075	76	80	93	rBV	2258551	2665137	100.00%	21.841%
2	1.172	93	96	106	rVB	121045	126530	4.75%	1.037%
3	1.270	108	112	115	rBV	21056	27429	1.03%	0.225%
4	1.404	128	134	140	rVB2	72314	107618	4.04%	0.882%
5	1.794	193	198	203	rVB3	18646	27734	1.04%	0.227%
6	2.623	330	334	344	rVB	15335	29535	1.11%	0.242%
7	5.513	796	808	820	rBV2	247580	704597	26.44%	5.774%
8	5.672	822	834	850	rVV	308221	860880	32.30%	7.055%
9	6.074	889	900	911	rBV	273963	719791	27.01%	5.899%
10	6.866	1020	1030	1051	rBV	434773	1027284	38.55%	8.419%
11	8.720	1326	1334	1343	rBV	1343032	2064469	77.46%	16.919%
12	10.116	1557	1563	1575	rBV	868210	1183526	44.41%	9.699%
13	11.140	1726	1731	1745	rVB	1092534	1415774	53.12%	11.603%
14	12.085	1880	1886	1894	rBV	981519	1241879	46.60%	10.178%

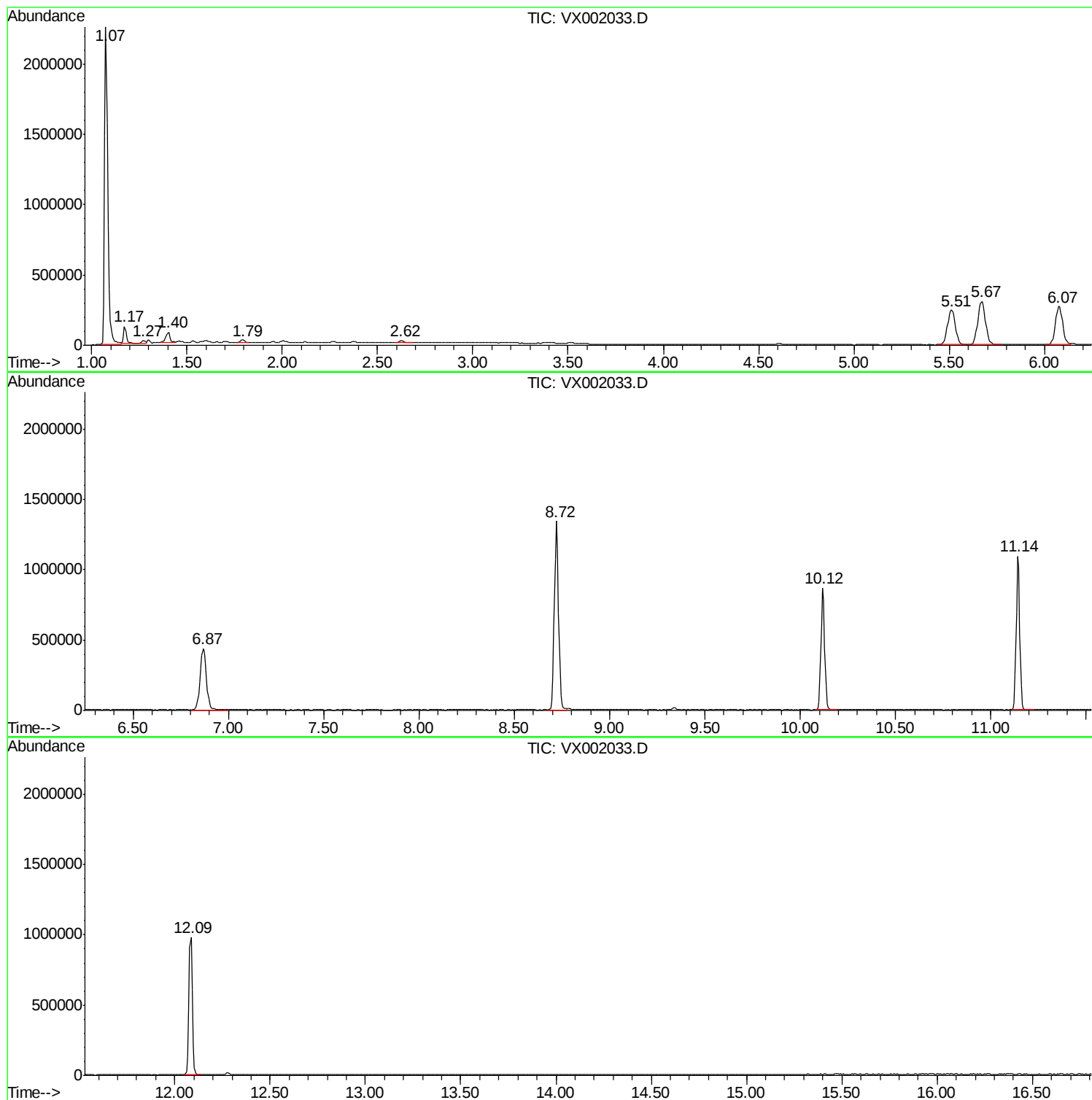
Sum of corrected areas: 12202183

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Quant Title : SW846 8260

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



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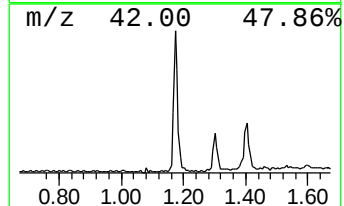
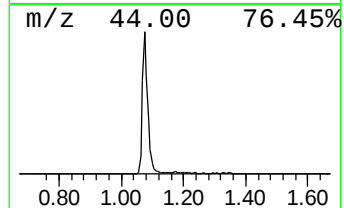
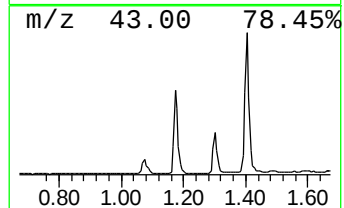
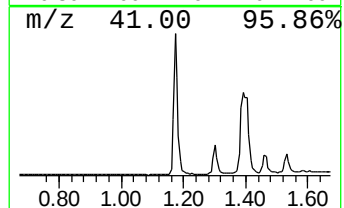
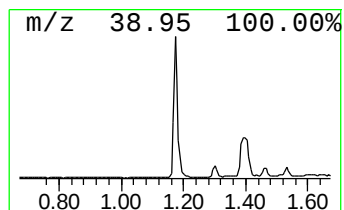
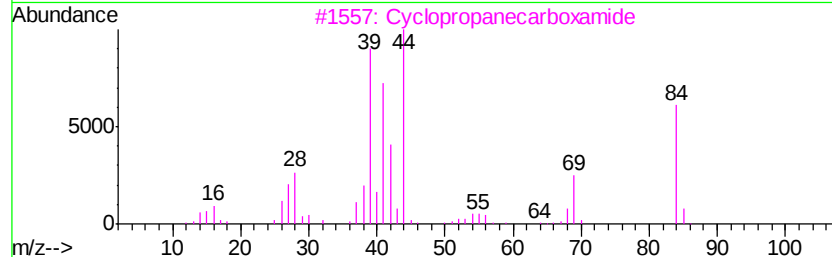
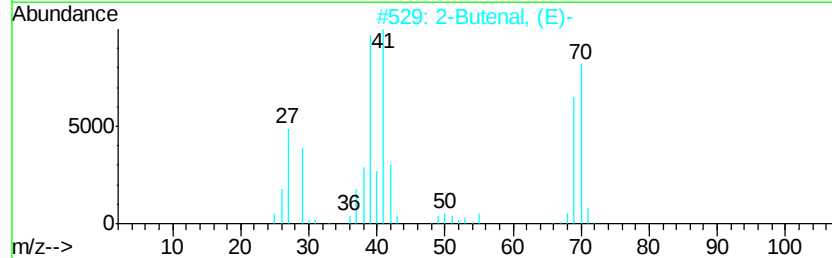
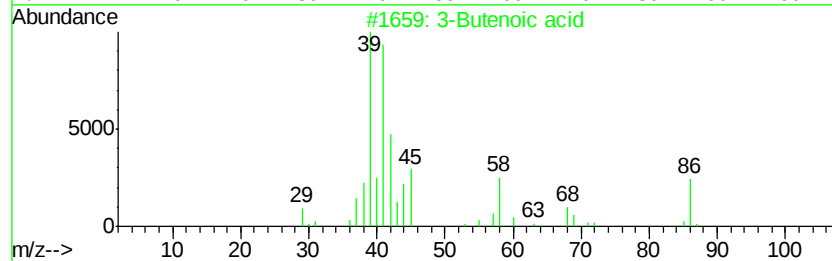
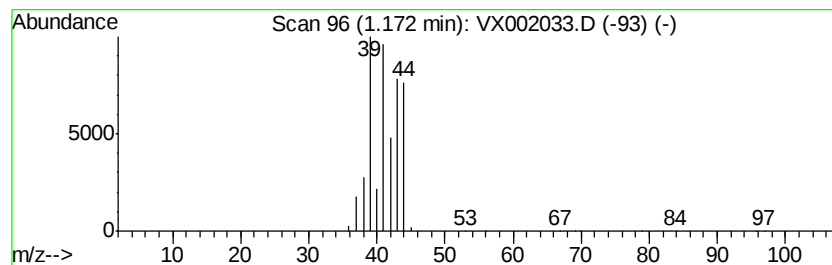
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Peak Number 2 3-Butenoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.17	7.35 ug/l	126530	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Butenoic acid	86	C4H6O2	000625-38-7	64
2		2-Butenal, (E)-	70	C4H6O	000123-73-9	64
3		Cyclopropanecarboxamide	85	C4H7NO	006228-73-5	39
4		2-Butenal	70	C4H6O	004170-30-3	38
5		2-Nonynoic acid	154	C9H14O2	001846-70-4	36



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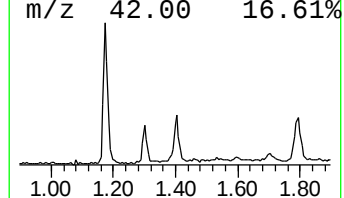
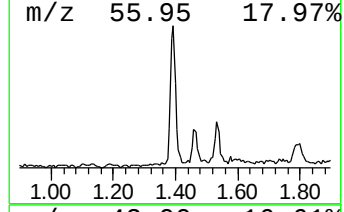
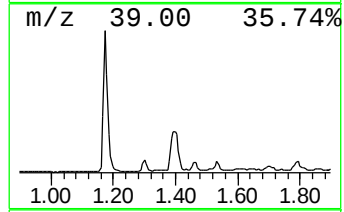
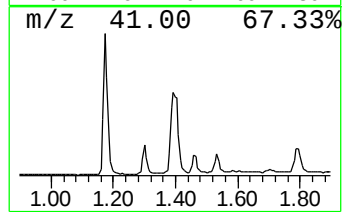
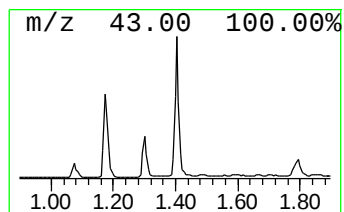
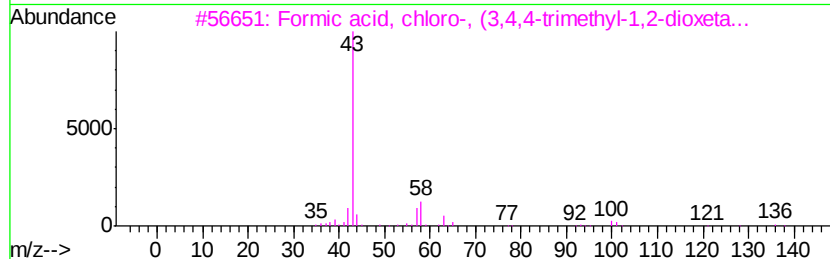
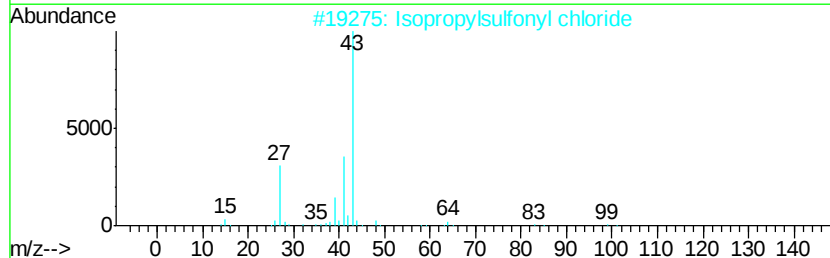
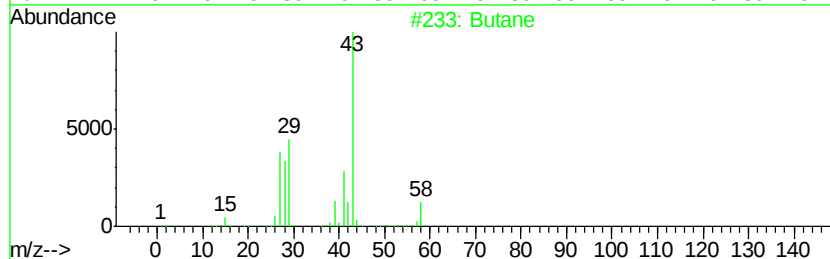
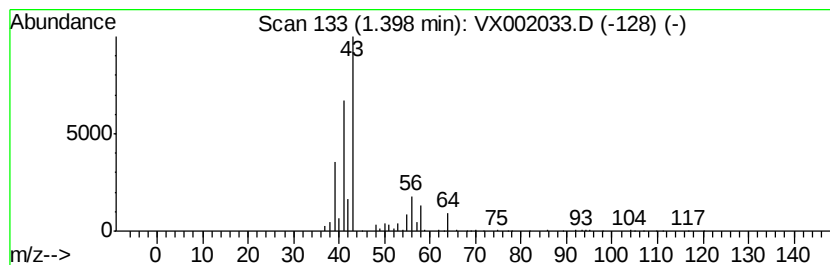
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Peak Number 3 Butane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.40	6.25 ug/l	107618	Pentafluorobenzene	5.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane		58	C4H10	000106-97-8	59
2	Isopropylsulfonyl chloride		142	C3H7ClO2S	010147-37-2	33
3	Formic acid, chloro-, (3,4,4-tri...		194	C7H11ClO4	107323-92-2	33
4	Oxirane, trimethyl-		86	C5H10O	005076-19-7	4
5	Pentane		72	C5H12	000109-66-0	4



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Butenoic acid	1.17	7.3	ug/l	126530	1	5.67	860880	50.0
Butane	1.40	6.3	ug/l	107618	1	5.67	860880	50.0