

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031819\
Data File : VX008176.D
Acq On : 18 Mar 2019 19:49
Operator : JC/SP
Sample : K1934-01 5X
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EFF-WW

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 : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\624X031819W.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	2.446	204	212	230	rBV	6545555	11872582	6.45%	5.857%
2	2.720	230	257	261	rBV2	34850531	183955243	100.00%	90.756%
3	12.268	1818	1823	1841	rBV	3697997	4521716	2.46%	2.231%
4	12.920	1925	1930	1944	rBV	2055742	2342096	1.27%	1.155%

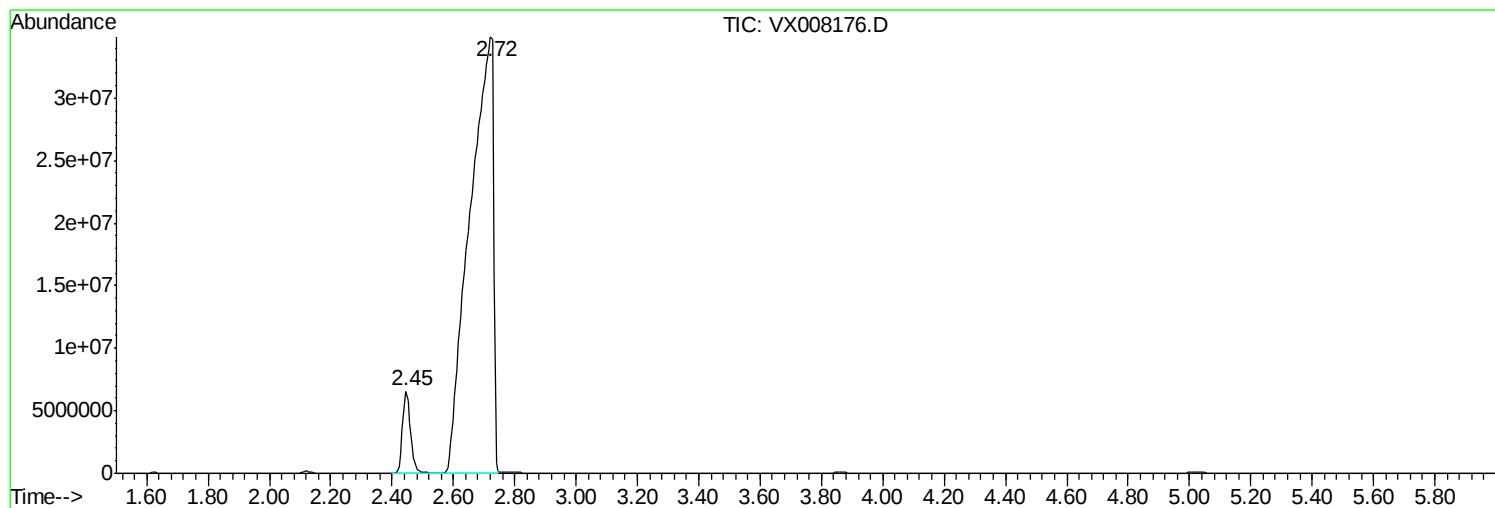
Sum of corrected areas: 202691637

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA X\METHOD\624X031819W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

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



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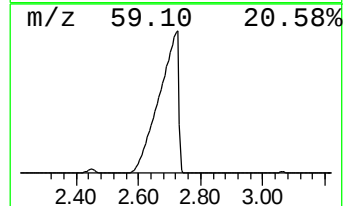
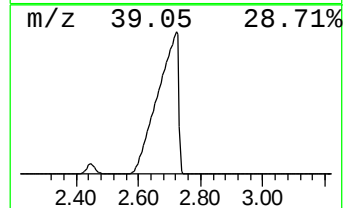
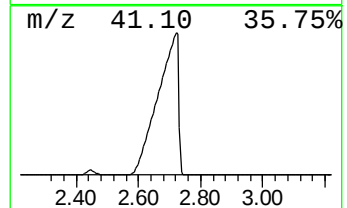
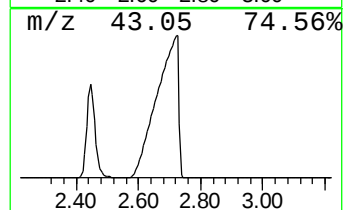
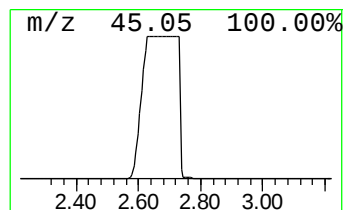
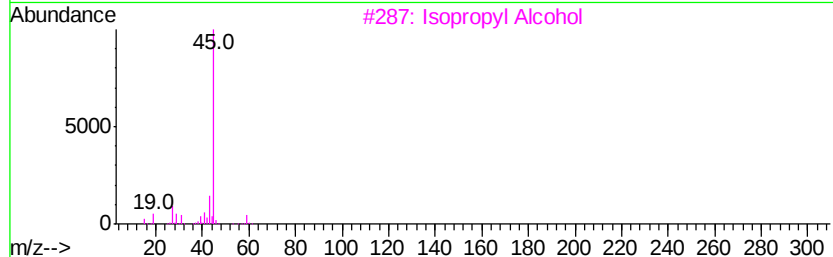
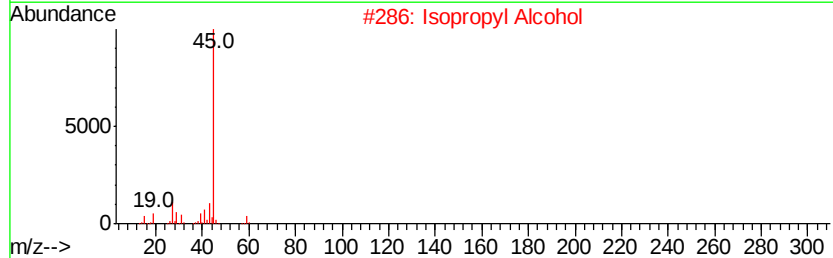
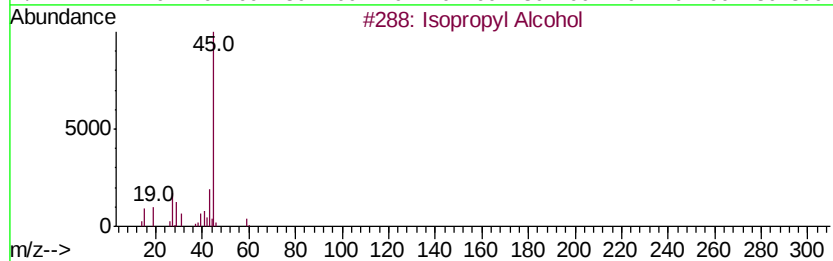
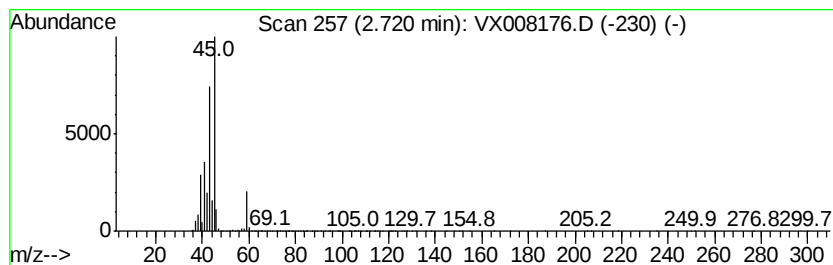
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA X\METHOD\624X031819W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

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

Peak Number 1 unknown2.72 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.72	30.00 ug/l	183955000	Bromochloromethane	5.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropyl Alcohol	60	C3H8O	000067-63-0	25
2		Isopropyl Alcohol	60	C3H8O	000067-63-0	25
3		Isopropyl Alcohol	60	C3H8O	000067-63-0	10
4		Methane, nitroso-	45	CH3NO	000865-40-7	9
5		Isobutane	58	C4H10	000075-28-5	9



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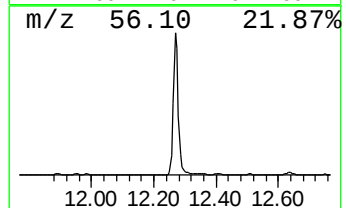
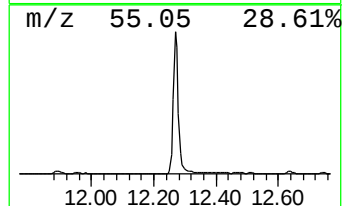
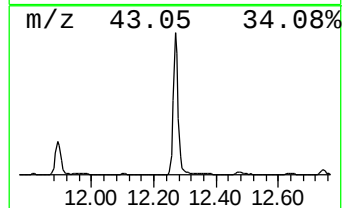
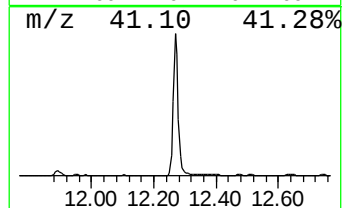
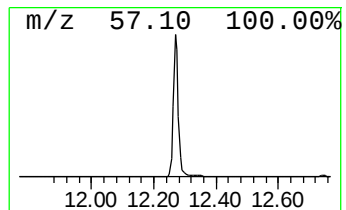
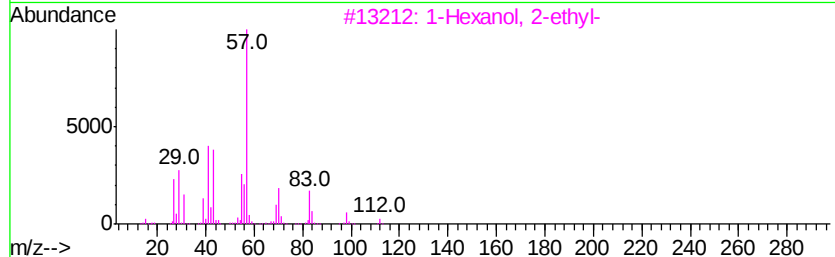
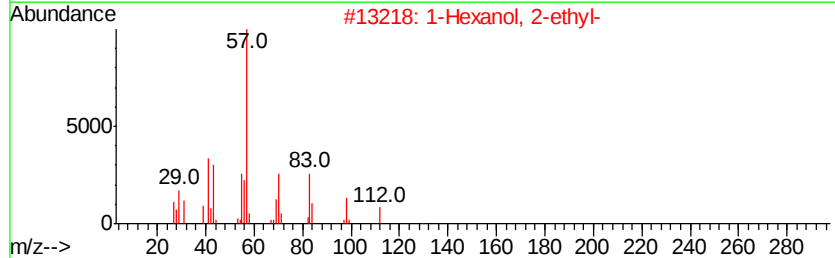
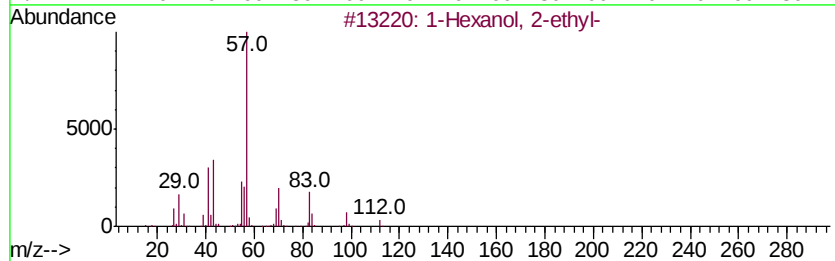
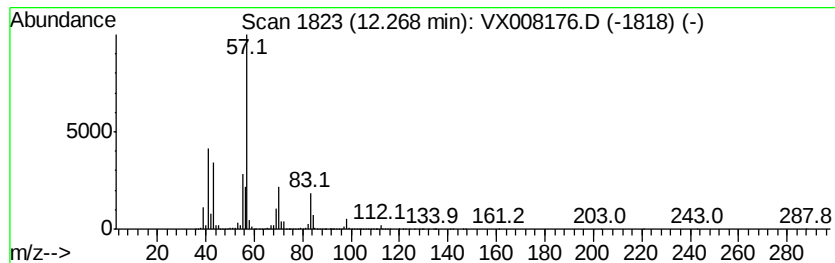
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Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	30.00 ug/l	4521720	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
3		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
4		1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	56
5		1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	50



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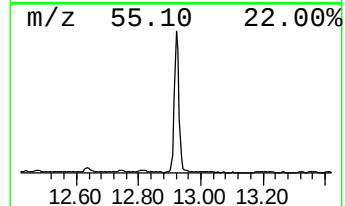
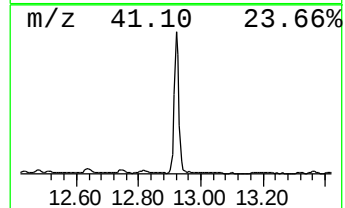
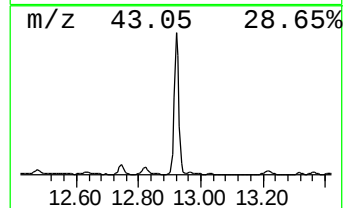
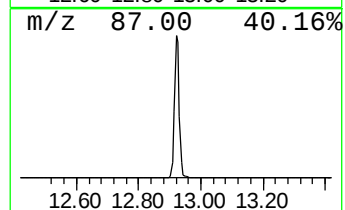
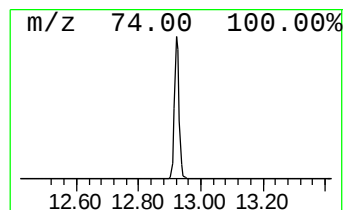
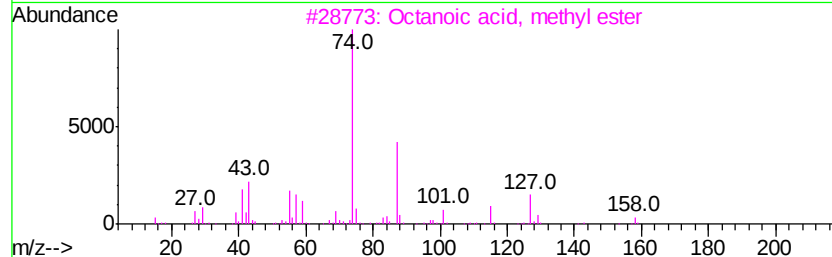
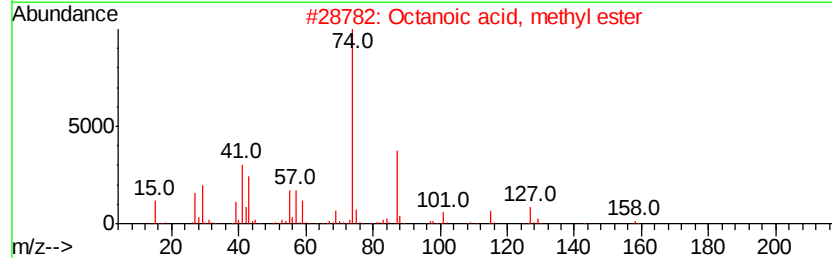
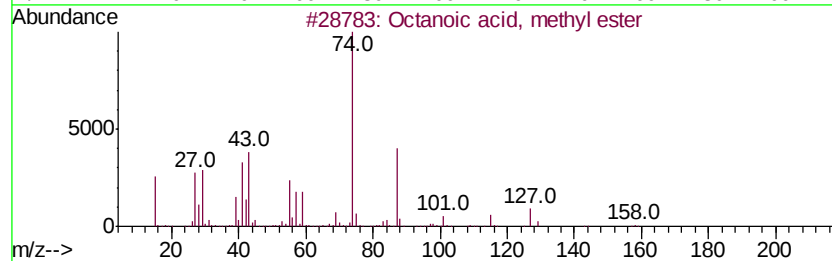
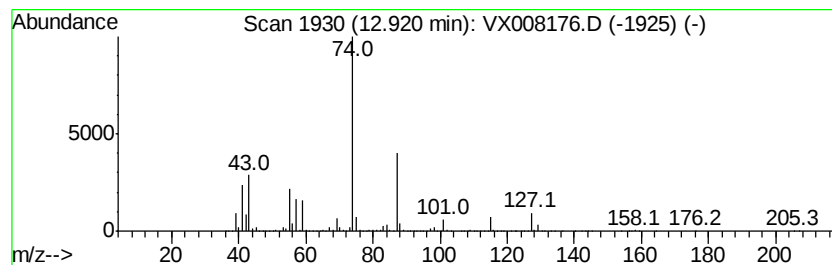
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Peak Number 3 Octanoic acid, methyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.92	15.54 ug/l	2342100	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octanoic acid, methyl ester	158	C9H18O2	000111-11-5	91
2		Octanoic acid, methyl ester	158	C9H18O2	000111-11-5	90
3		Octanoic acid, methyl ester	158	C9H18O2	000111-11-5	83
4		Nonanoic acid, methyl ester	172	C10H20O2	001731-84-6	78
5		Nonanoic acid, methyl ester	172	C10H20O2	001731-84-6	72



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.72	2.72	30.0	ug/l	183955000	1	5.02	183955000	30.0
1-Hexanol, 2-ethyl-	12.27	30.0	ug/l	4521720	3	10.11	4521720	30.0
Octanoic acid, me...	12.92	15.5	ug/l	2342100	3	10.11	4521720	30.0