

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX020720\
 Data File : VX014871.D
 Acq On : 07 Feb 2020 15:14
 Operator : JC/SP
 Sample : L1410-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 001-WILLETS-PT-BLVD-(FEB)

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\624X011520W.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	1.576	110	113	121	rVB2	37294	53172	1.36%	0.403%
2	2.429	246	253	267	rBV	2282299	3672360	93.60%	27.840%
3	2.588	272	279	294	rBV	2253506	3923600	100.00%	29.745%
4	5.008	664	676	688	rBV2	148067	435108	11.09%	3.299%
5	6.051	839	847	859	rBV	156320	400110	10.20%	3.033%
6	6.849	970	978	993	rBV	362974	851303	21.70%	6.454%
7	8.709	1276	1283	1289	rVV	757416	1158291	29.52%	8.781%
8	8.776	1289	1294	1301	rVB	262303	412474	10.51%	3.127%
9	10.105	1507	1512	1520	rVB	729001	991657	25.27%	7.518%
10	10.587	1587	1591	1596	rVB	107588	133482	3.40%	1.012%
11	10.904	1638	1643	1651	rBV3	32331	51026	1.30%	0.387%
12	11.129	1675	1680	1687	rBV	620118	815384	20.78%	6.181%
13	11.806	1786	1791	1796	rBV4	24426	41550	1.06%	0.315%
14	12.263	1863	1866	1871	rBV	29561	43141	1.10%	0.327%
15	13.385	2046	2050	2056	rBV	74489	97411	2.48%	0.738%
16	13.641	2086	2092	2097	rBV5	30518	48759	1.24%	0.370%
17	14.854	2286	2291	2298	rBV	41723	62010	1.58%	0.470%

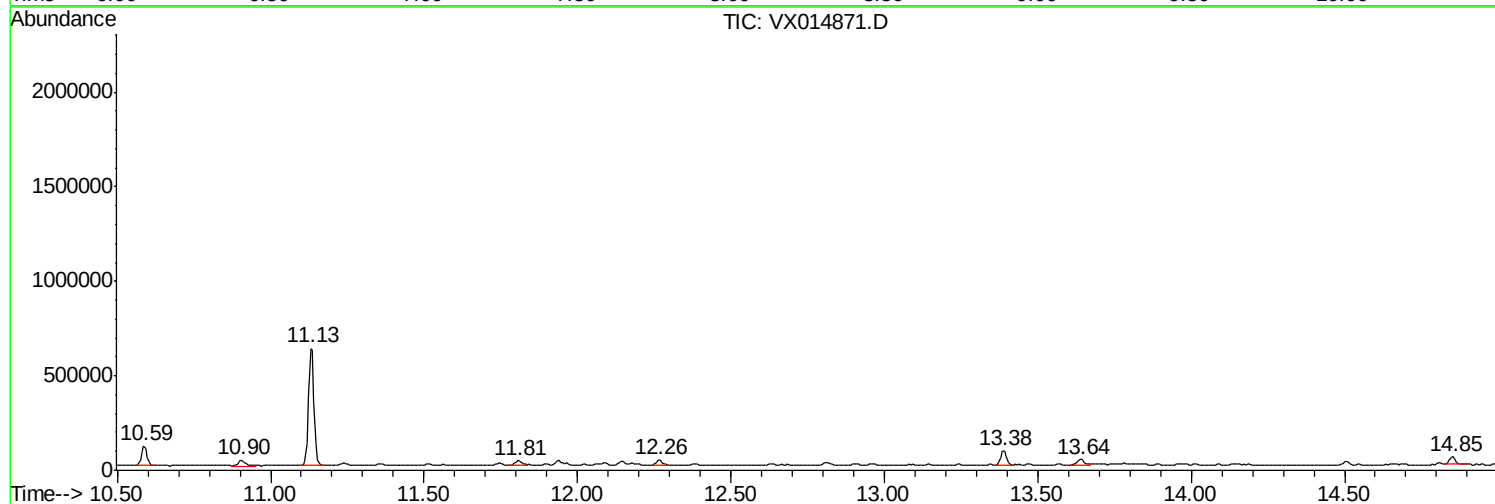
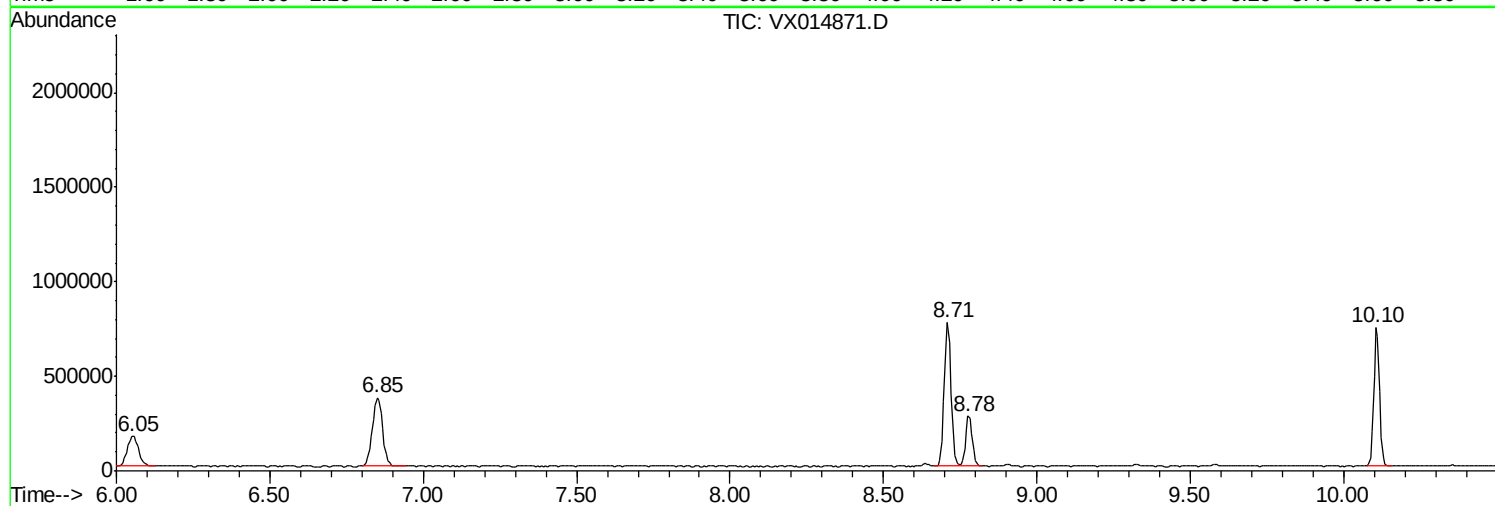
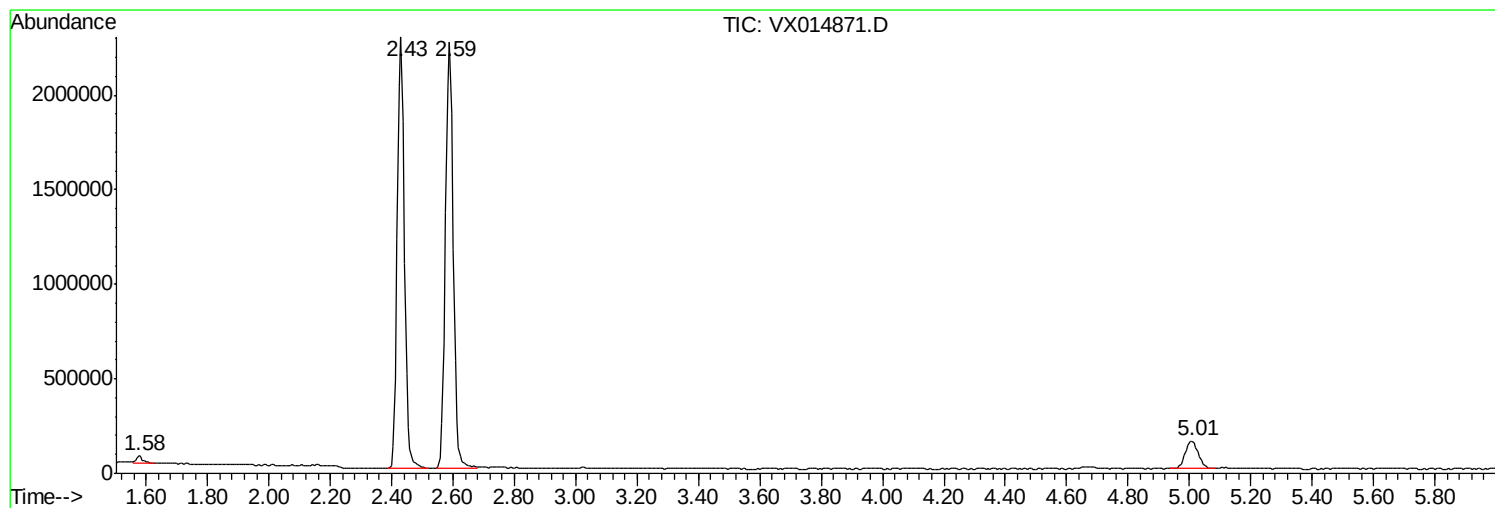
Sum of corrected areas: 13190838

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA X\METHOD\624X011520W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

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



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 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

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

 Peak Number 1 Methanethiol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.58	3.67 ug/l	53172	Bromochloromethane	5.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Methanethiol	48 CH4S	000074-93-1 80
2		Methanethiol	48 CH4S	000074-93-1 72
3		Ethanol, 2-mercapto-	78 C2H6OS	000060-24-2 64
4		Propanal, 3-(methylthio)-	104 C4H8OS	003268-49-3 23
5		Ethane, 1-chloro-1-fluoro-	82 C2H4ClF	001615-75-4 23

 Peak Number 2 Isopropyl Alcohol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.59	270.53 ug/l	3923600	Bromochloromethane	5.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Isopropyl Alcohol	60 C3H8O	000067-63-0 86
2		Isopropyl Alcohol	60 C3H8O	000067-63-0 78
3		Isopropyl Alcohol	60 C3H8O	000067-63-0 78
4		Formamide	45 CH3NO	000075-12-7 5
5		Propanoic acid, 2-hydroxy-, meth...	104 C4H8O3	002155-30-8 4

 Peak Number 3 4-Octene, 2,6-dimethyl-, [S... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.59	4.04 ug/l	133482	Chlorobenzene-d5	10.10
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		4-Octene, 2,6-dimethyl-, [S-(E)]-	140 C10H20	062960-76-3 90
2		3-Octene, 4-ethyl-	140 C10H20	053966-51-1 81
3		3-Heptene, 4-propyl-	140 C10H20	004485-13-6 64
4		3-Heptene, 4-propyl-	140 C10H20	004485-13-6 64
5		2-Octene, 2,6-dimethyl-	140 C10H20	004057-42-5 64

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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Methanethiol	1.58	3.7	ug/l	53172	1	5.01	435108	30.0
Isopropyl Alcohol	2.59	270.5	ug/l	3923600	1	5.01	435108	30.0
4-Octene, 2,6-dim...	10.59	4.0	ug/l	133482	3	10.10	991657	30.0