

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\WX040720\
 Data File : VX015539.D
 Acq On : 07 Apr 2020 15:33
 Operator : JC/SP
 Sample : L2140-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 12 Sample Multiplier: 1

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

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\624X032520W.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	2.088	193	197	202	rVB	62835	84966	1.16%	0.345%
2	2.417	244	251	265	rBV	4522248	7328889	100.00%	29.743%
3	2.576	269	277	290	rBV	923221	1695501	23.13%	6.881%
4	4.649	608	617	626	rBV2	65933	199444	2.72%	0.809%
5	4.990	661	673	684	rBV	432127	1286404	17.55%	5.221%
6	6.039	835	845	857	rBV	455427	1189968	16.24%	4.829%
7	6.837	966	976	988	rBV	1016453	2438163	33.27%	9.895%
8	8.624	1263	1269	1275	rBV	126750	195906	2.67%	0.795%
9	8.697	1275	1281	1288	rVV	2052032	3261964	44.51%	13.238%
10	8.770	1288	1293	1302	rVB	250726	429577	5.86%	1.743%
11	9.623	1429	1433	1439	rBV	77758	101937	1.39%	0.414%
12	10.099	1505	1511	1522	rVB	2126182	2914966	39.77%	11.830%
13	10.343	1546	1551	1559	rVB	203410	277278	3.78%	1.125%
14	10.684	1602	1607	1613	rBV	82020	107628	1.47%	0.437%
15	11.123	1674	1679	1686	rVB	1891110	2338266	31.90%	9.489%
16	11.800	1786	1790	1794	rVB2	91352	118289	1.61%	0.480%
17	12.257	1861	1865	1874	rBV	97257	177341	2.42%	0.720%
18	12.861	1958	1964	1968	rBV	333375	404205	5.52%	1.640%
19	13.629	2086	2090	2099	rVB10	49574	89988	1.23%	0.365%

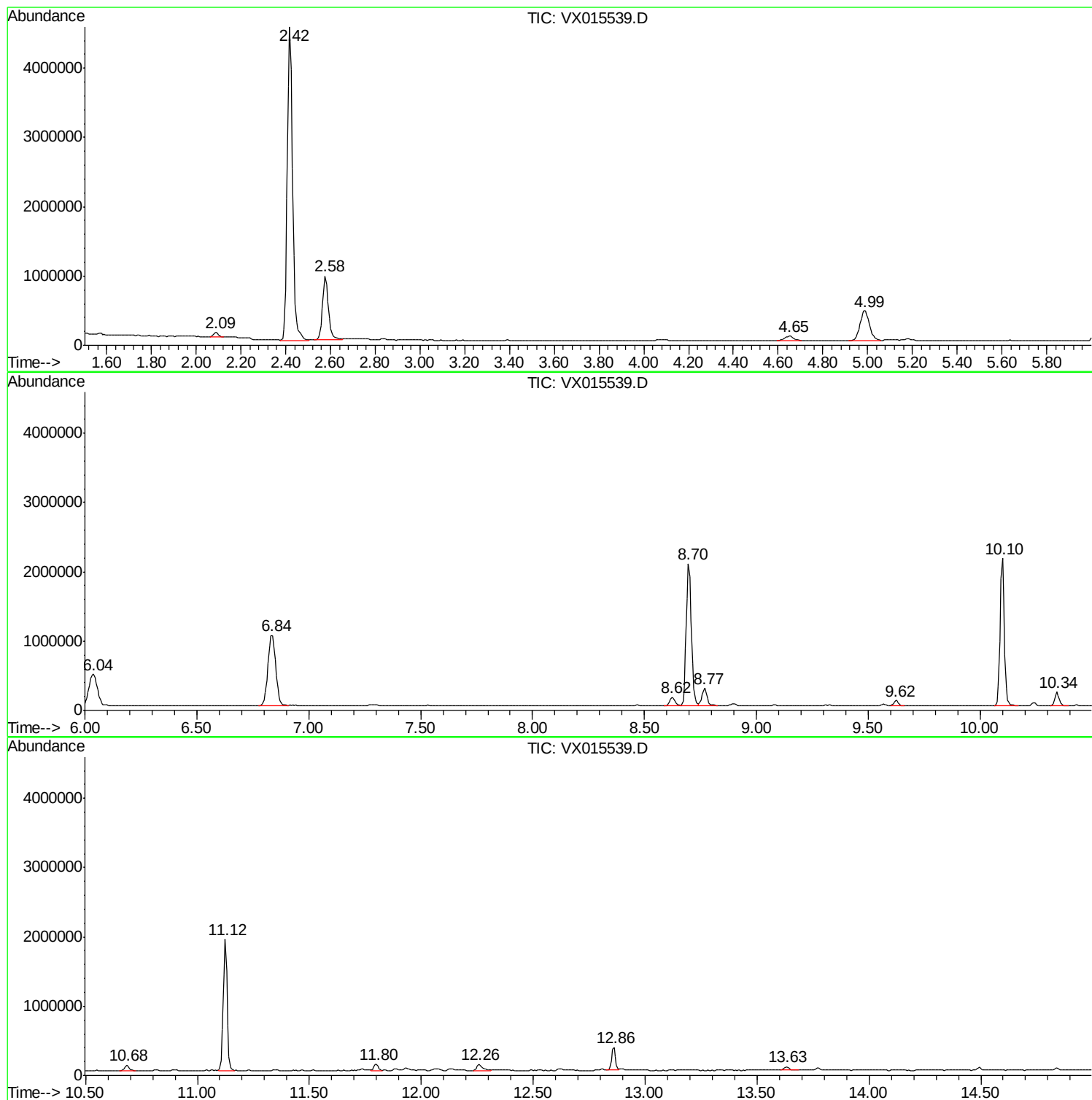
Sum of corrected areas: 24640680

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ClientSampleId :
001-WILLETS-PT-BLVD-(APR)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA X\METHOD\624X032520W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

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



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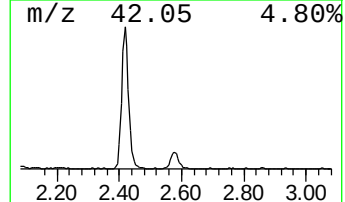
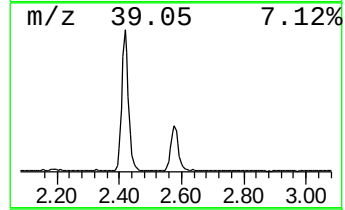
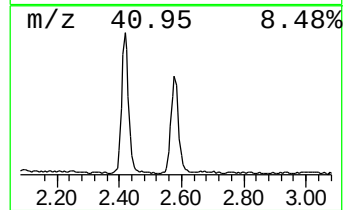
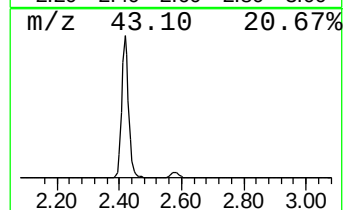
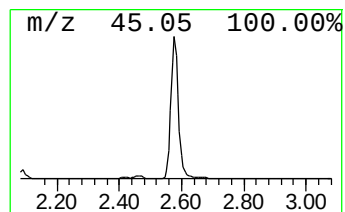
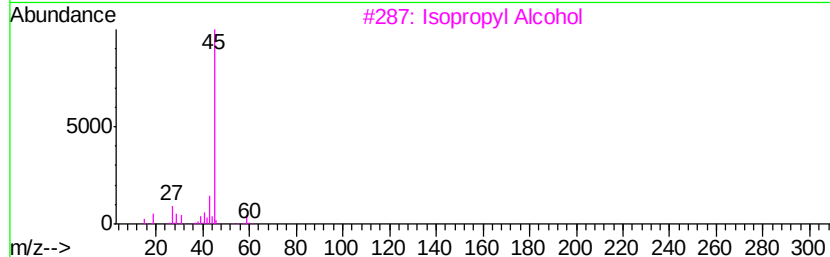
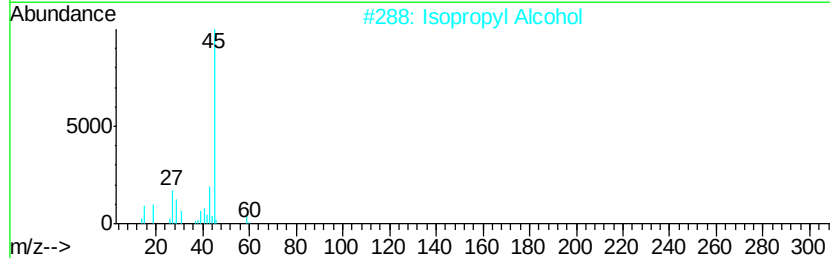
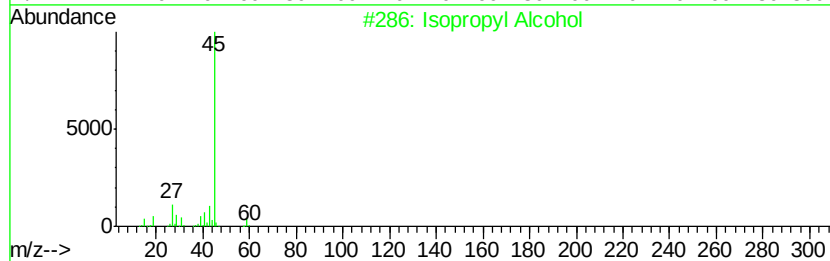
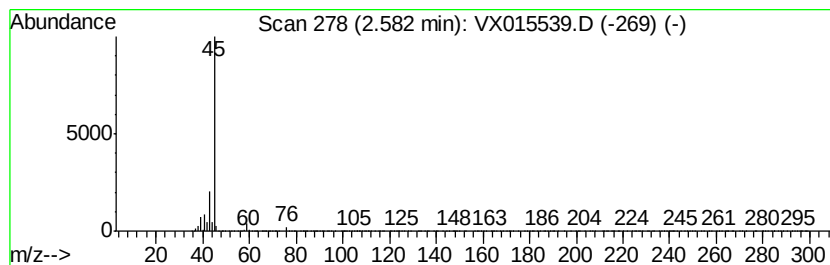
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA X\METHOD\624X032520W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

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

Peak Number 1 Isopropyl Alcohol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.58	39.54 ug/l	1695500	Bromochloromethane	4.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropyl Alcohol	60	C3H8O	000067-63-0	83
2		Isopropyl Alcohol	60	C3H8O	000067-63-0	83
3		Isopropyl Alcohol	60	C3H8O	000067-63-0	78
4		(R)-(-)-2-Pentanol	88	C5H12O	031087-44-2	9
5		2-Pentanol	88	C5H12O	006032-29-7	9



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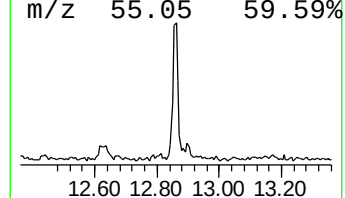
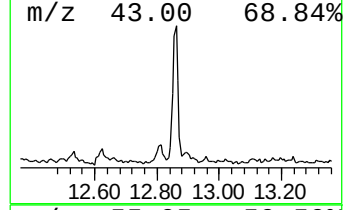
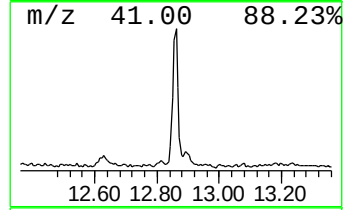
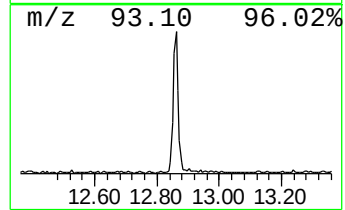
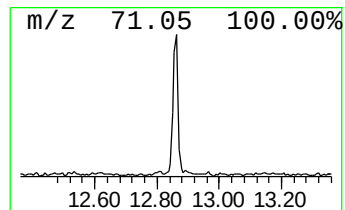
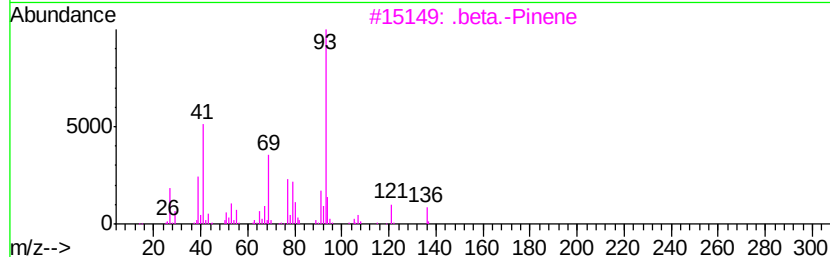
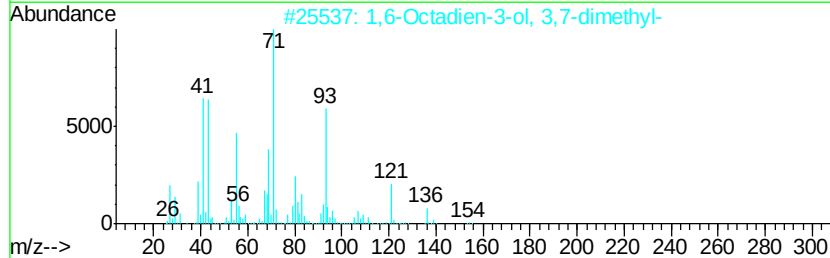
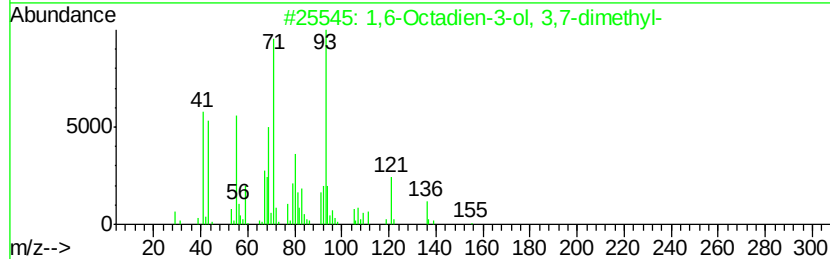
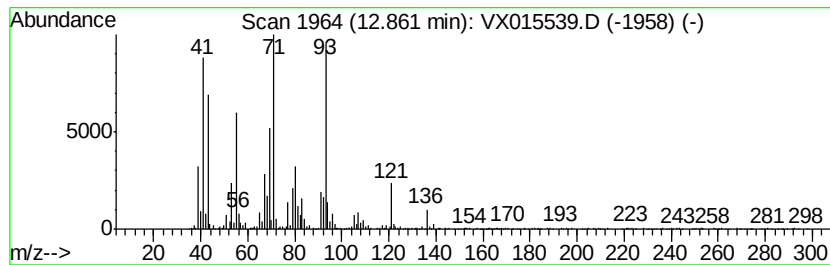
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ClientSampled :
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TIC Integration Parameters: LSCINT.P

Peak Number 2 1,6-Octadien-3-ol, 3,7-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.86	4.16 ug/l	404205	Chlorobenzene-d5	10.10	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6-Octadien-3-ol, 3,7-dimethyl-	154	C10H18O	000078-70-6	87
2	1,6-Octadien-3-ol, 3,7-dimethyl-	154	C10H18O	000078-70-6	60
3	.beta.-Pinene	136	C10H16	000127-91-3	55
4	.beta.-Myrcene	136	C10H16	000123-35-3	55
5	1,6-Octadien-3-ol, 3,7-dimethyl-...	196	C12H20O2	000115-95-7	52



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

TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Isopropyl Alcohol	2.58	39.5	ug/l	1695500	1	4.98	1286400	30.0
1,6-Octadien-3-ol...	12.86	4.2	ug/l	404205	3	10.10	2914970	30.0