

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU062019\
 Data File : VU032914.D
 Acq On : 20 Jun 2019 20:51
 Operator : JC/SP
 Sample : K3413-13DL 4X
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GALG1DL

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

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 : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR060419WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.180	22	28	35	rBV	22540	20995	5.71%	0.711%
2	1.283	56	60	71	rVB3	5483	8557	2.33%	0.290%
3	1.395	87	95	104	rBV2	41953	45900	12.49%	1.554%
4	1.550	136	143	153	rVB5	6161	8496	2.31%	0.288%
5	1.678	174	183	192	rVB	31422	36989	10.06%	1.252%
6	1.756	202	207	216	rVB5	2798	3810	1.04%	0.129%
7	1.900	241	252	257	rBV6	2628	4540	1.24%	0.154%
8	1.945	259	266	276	rVB4	2942	4405	1.20%	0.149%
9	2.267	356	366	377	rBV	60610	93806	25.52%	3.176%
10	2.315	377	381	392	rVB7	2215	3934	1.07%	0.133%
11	4.186	949	963	991	rBV	56780	151256	41.15%	5.122%
12	4.386	1009	1025	1044	rBV	28638	67730	18.43%	2.293%
13	4.640	1087	1104	1123	rBV	51780	121869	33.16%	4.127%
14	5.328	1296	1318	1344	rBV3	97361	278309	75.72%	9.424%
15	5.878	1476	1489	1508	rBV	103564	204378	55.61%	6.920%
16	6.324	1615	1628	1648	rBV2	69389	137293	37.35%	4.649%
17	7.222	1895	1907	1920	rBV2	34512	61737	16.80%	2.090%
18	7.559	1997	2012	2029	rBV	122353	214367	58.32%	7.259%
19	7.845	2091	2101	2116	rBV	21718	37618	10.24%	1.274%
20	8.170	2191	2202	2207	rBV2	3601	6164	1.68%	0.209%
21	8.222	2207	2218	2234	rVB	111488	190655	51.87%	6.456%
22	8.308	2234	2245	2276	rBV	194571	367539	100.00%	12.445%
23	9.083	2474	2486	2506	rBV	158954	265451	72.22%	8.988%
24	10.427	2893	2904	2918	rBV	61625	97866	26.63%	3.314%
25	10.604	2946	2959	2973	rBV2	9356	17126	4.66%	0.580%
26	11.479	3219	3231	3252	rBV	160197	264712	72.02%	8.963%
27	11.855	3337	3348	3371	rBV	138534	227791	61.98%	7.713%
28	12.479	3532	3542	3552	rBV	2764	4899	1.33%	0.166%
29	14.279	4091	4102	4109	rBV2	3042	5068	1.38%	0.172%

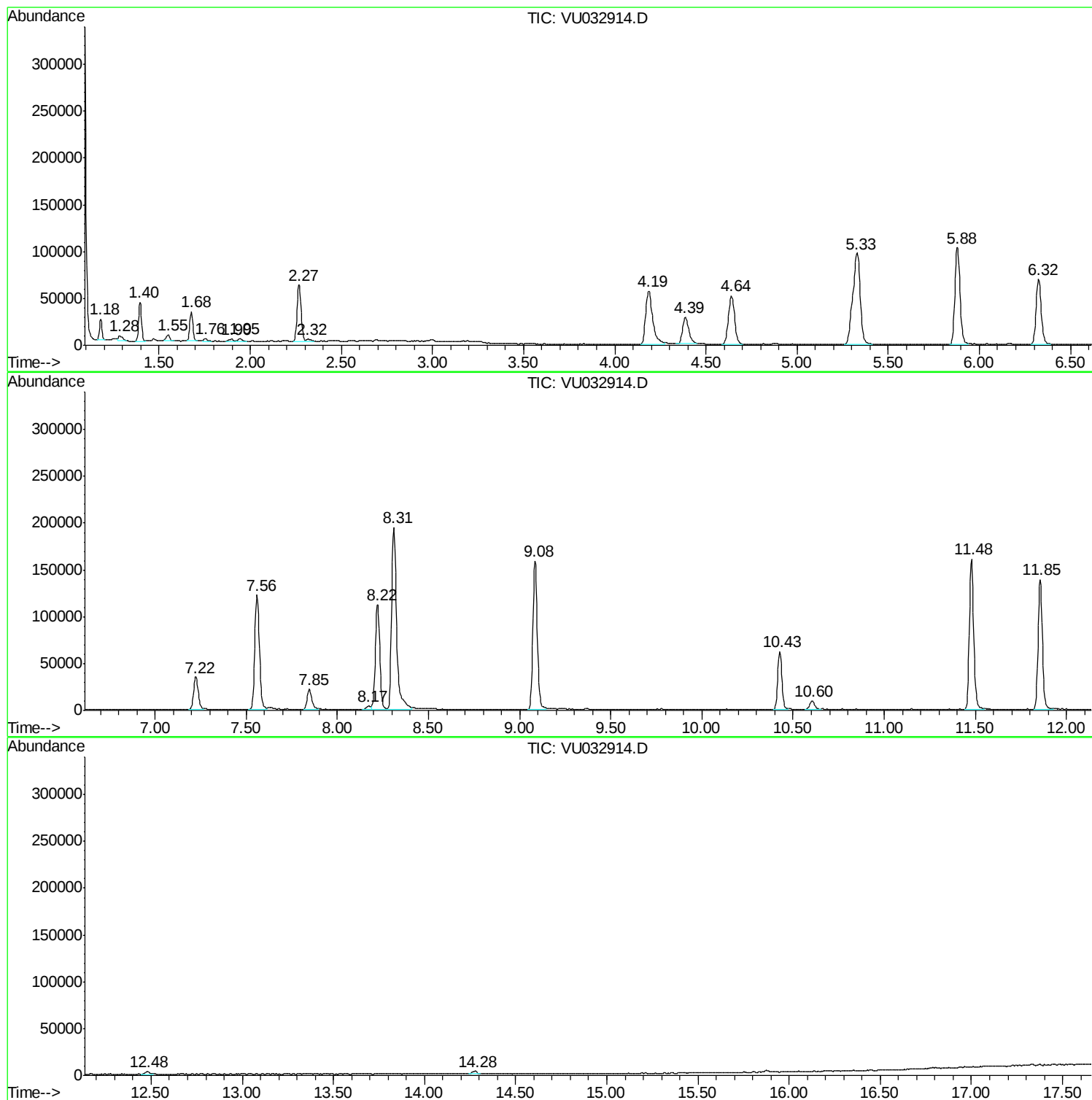
Sum of corrected areas: 2953260

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU062019\
Data File : VU032914.D
Acq On : 20 Jun 2019 20:51
Operator : JC/SP
Sample : K3413-13DL 4X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GALG1DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR060419WMA.M
Quant Title : TRACE VOA SOM01.0

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU062019\
 Data File : VU032914.D
 Acq On : 20 Jun 2019 20:51
 Operator : JC/SP
 Sample : K3413-13DL 4X
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GALG1DL

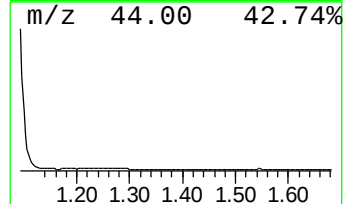
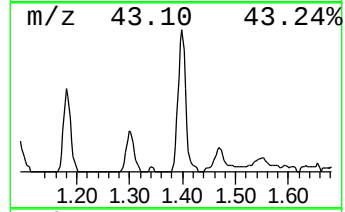
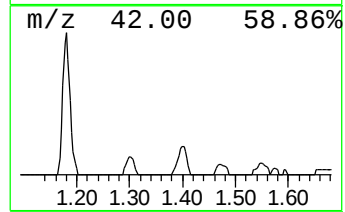
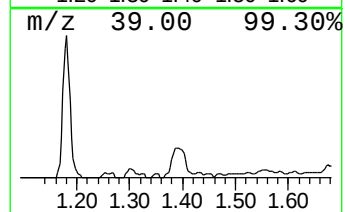
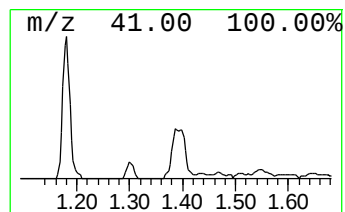
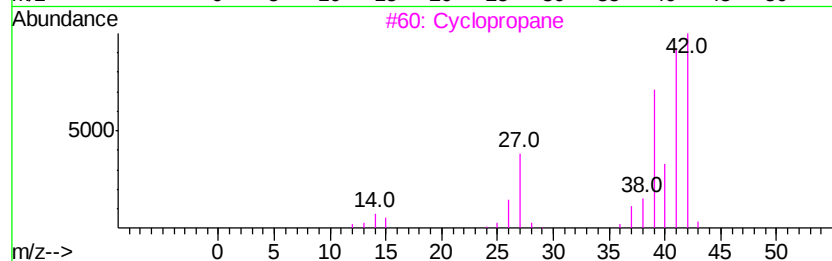
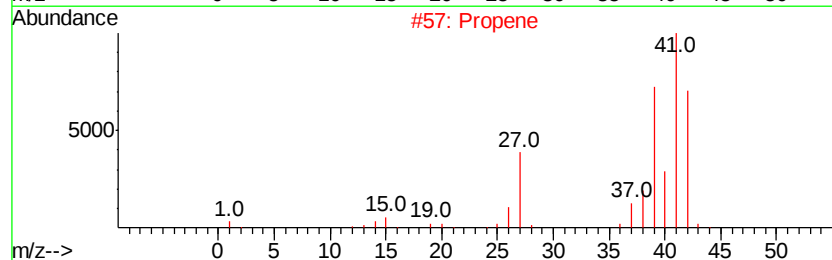
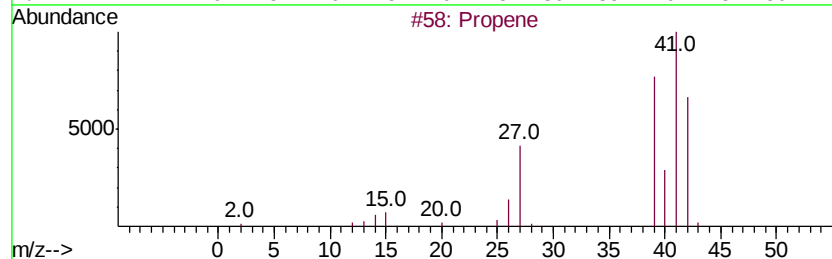
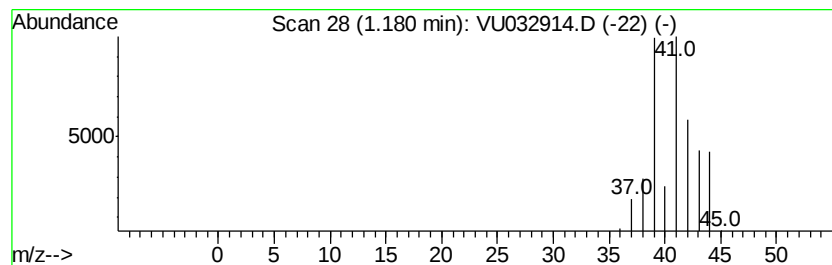
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR060419WMA.M
 Quant Title : TRACE VOA SOM01.0

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

 Peak Number 1 (DEL) Alkane: Cyclic1.18 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.18	0.51 ug/L	20995	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propene	42	C3H6	000115-07-1	40
2		Propene	42	C3H6	000115-07-1	9
3		Cyclopropane	42	C3H6	000075-19-4	4
4		Cyclopropane	42	C3H6	000075-19-4	4
5		Cyclopropene	40	C3H4	002781-85-3	4



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU062019\
Data File : VU032914.D
Acq On : 20 Jun 2019 20:51
Operator : JC/SP
Sample : K3413-13DL 4X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GALG1DL

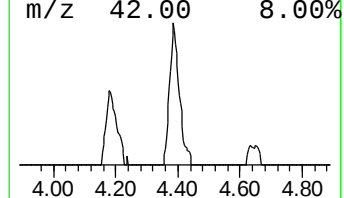
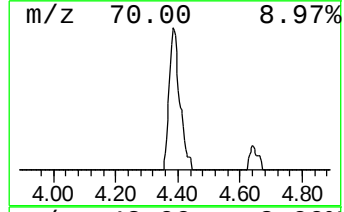
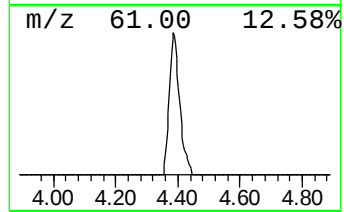
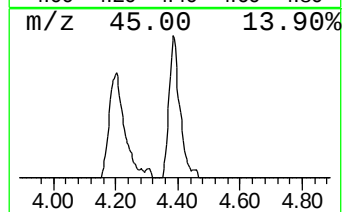
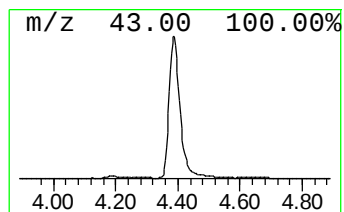
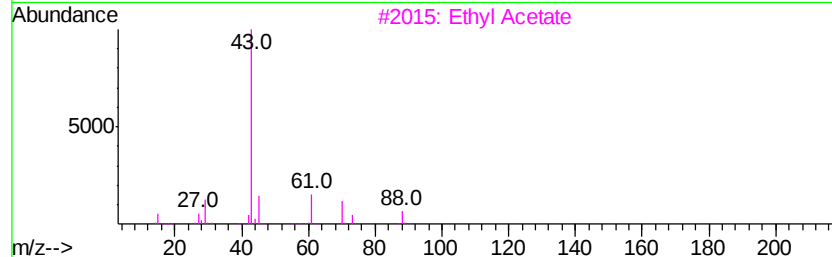
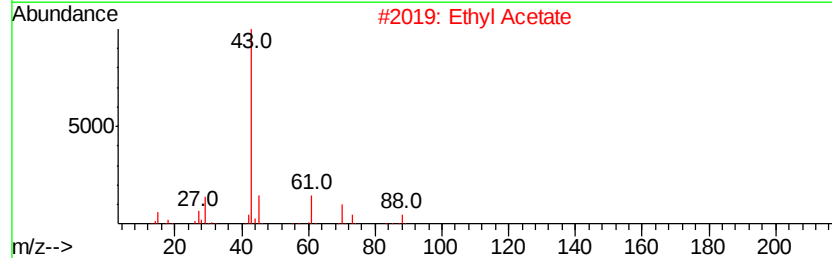
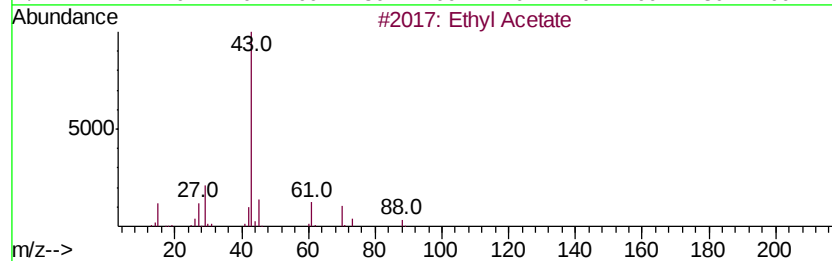
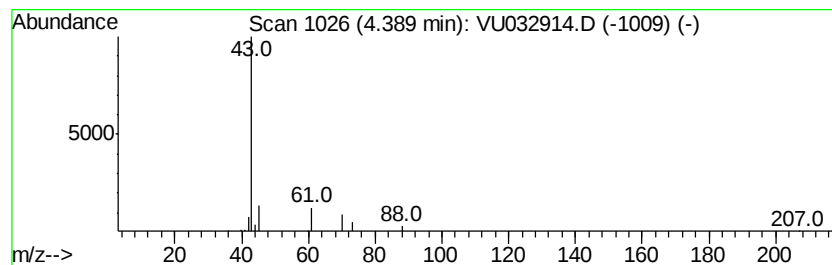
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR060419WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 Ethyl Acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	1.66 ug/L	67730	1,4-Difluorobenzene	5.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethyl Acetate	88	C4H8O2	000141-78-6	90
2			Ethyl Acetate	88	C4H8O2	000141-78-6	78
3			Ethyl Acetate	88	C4H8O2	000141-78-6	74
4			Ethyl Acetate	88	C4H8O2	000141-78-6	72
5			CH3C(O)CH2CH2OH	88	C4H8O2	000590-90-9	36



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU062019\
Data File : VU032914.D
Acq On : 20 Jun 2019 20:51
Operator : JC/SP
Sample : K3413-13DL 4X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GALG1DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR060419WMA.M
Quant Title : TRACE VOA SOM01.0

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Cyc...	1.18	0.5	ug/L	20995	1	5.88	204378	5.0
Ethyl Acetate	4.39	1.7	ug/L	67730	1	5.88	204378	5.0