

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD101119\
 Data File : VD063947.D
 Acq On : 11 Oct 2019 19:58
 Operator : VA/SY
 Sample : K5316-03
 Misc : 7.44G/5ml/MSVOA D/SOIL
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TP-9

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 : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D100219S.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	2.315	62	67	79	rBV3	20415	53774	5.77%	1.133%
2	3.409	249	253	260	rVB6	5323	11329	1.22%	0.239%
3	4.162	371	381	390	rBV4	10138	27280	2.93%	0.575%
4	7.920	1012	1020	1025	rBV3	121492	289080	31.04%	6.091%
5	7.985	1025	1031	1040	rVB	200807	448026	48.10%	9.440%
6	8.332	1083	1090	1102	rVB	112849	248302	26.66%	5.232%
7	8.867	1172	1181	1191	rBV	314720	625351	67.14%	13.176%
8	10.338	1422	1431	1440	rBV	515429	931419	100.00%	19.625%
9	10.726	1492	1497	1502	rBV3	17328	37064	3.98%	0.781%
10	11.644	1646	1653	1662	rBV	455422	778748	83.61%	16.408%
11	12.632	1815	1821	1830	rBV	350500	580393	62.31%	12.229%
12	12.961	1872	1877	1884	rVB6	5394	10137	1.09%	0.214%
13	13.579	1975	1982	1989	rBV	411717	695573	74.68%	14.656%
14	14.108	2068	2072	2076	rVB5	6773	9643	1.04%	0.203%

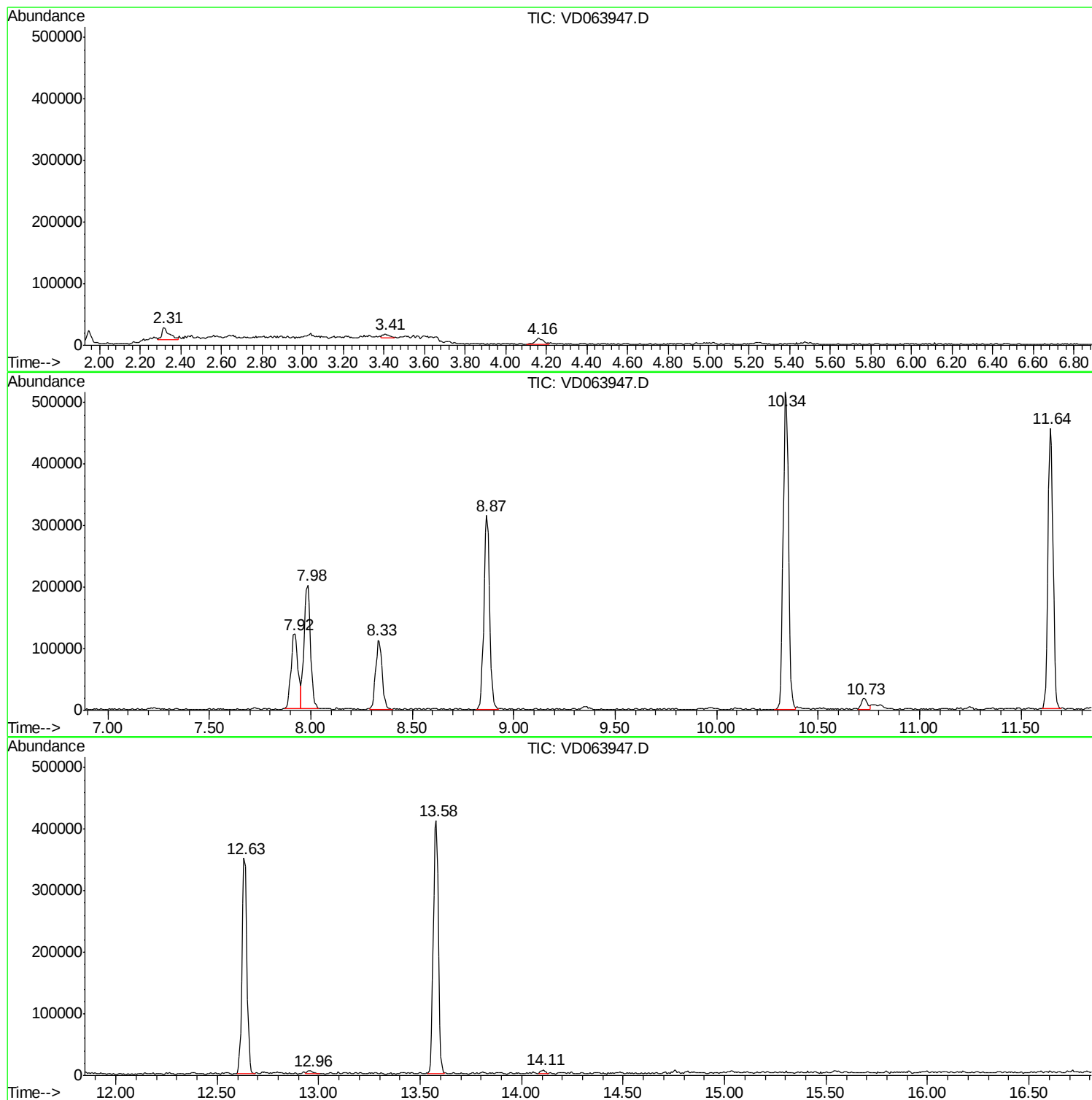
Sum of corrected areas: 4746119

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100219S.M
Quant Title : SW846 8260

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



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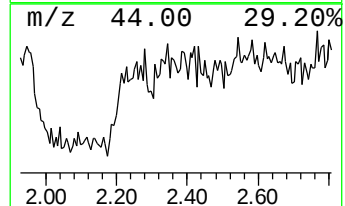
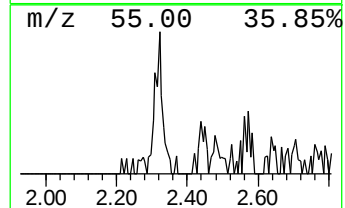
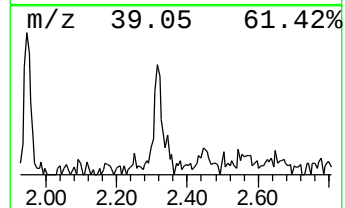
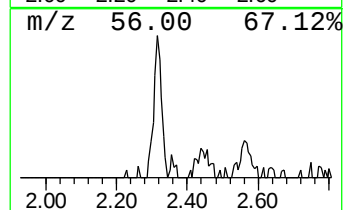
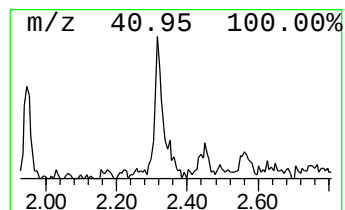
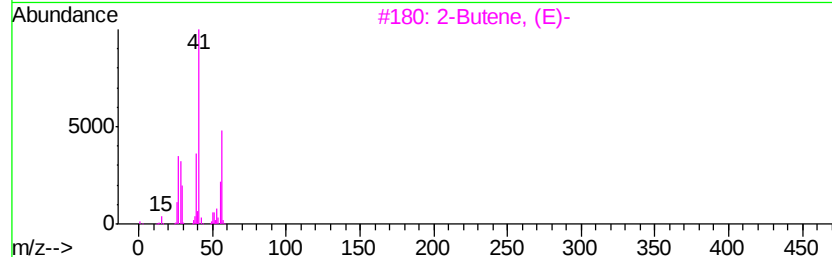
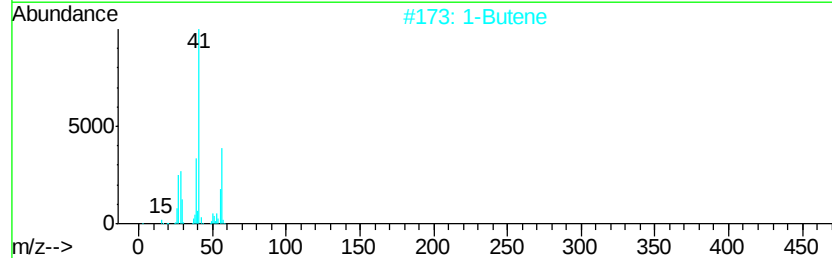
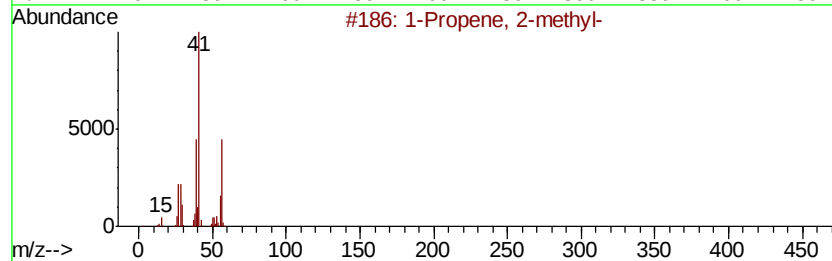
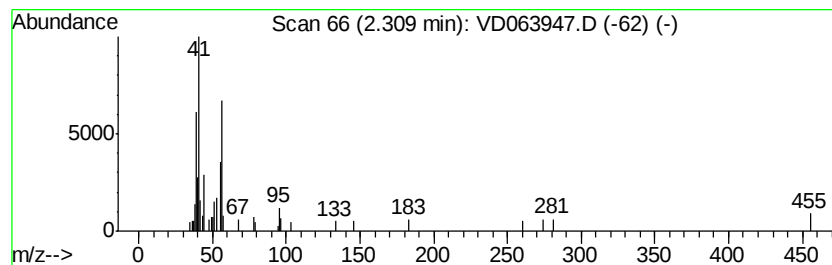
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Quant Title : SW846 8260

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

Peak Number 1 1-Propene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.31	6.00 ug/l	53774	Pentafluorobenzene	7.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Propene, 2-methyl-	56	C4H8	000115-11-7	80
2		1-Butene	56	C4H8	000106-98-9	80
3		2-Butene, (E)-	56	C4H8	000624-64-6	72
4		2-Butene	56	C4H8	000107-01-7	59
5		3-Buten-1-ol, 3-methyl-	86	C5H10O	000763-32-6	45



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Propene, 2-methyl-	2.31	6.0	ug/l	53774	1	7.98	448026	50.0