

Data Path : Z:\voasrv\HPCHEM1\MSVOA D\Data\VD041819\
 Data File : VD061901.D
 Acq On : 19 Apr 2019 7:23
 Operator : VA/AP
 Sample : K2441-02
 Misc : 6.34g/5ml/MSVOA D/SOIL
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TP-18-S2

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D041519S.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	1.500	6	8	24	rVB	129413	336002	76.97%	11.533%
2	2.878	144	148	150	rBV2	2508	5058	1.16%	0.174%
3	3.222	179	183	191	rBV	22184	58867	13.48%	2.021%
4	3.833	239	245	247	rBV	18655	47583	10.90%	1.633%
5	3.882	248	250	257	rVB2	15109	33480	7.67%	1.149%
6	6.156	476	481	488	rVB2	5202	17441	4.00%	0.599%
7	6.934	554	560	567	rBV	98985	270385	61.94%	9.281%
8	7.052	567	572	581	rVB	79153	220825	50.58%	7.580%
9	7.465	608	614	624	rVB	81657	211582	48.47%	7.262%
10	8.233	685	692	703	rBV	133638	332316	76.12%	11.406%
11	10.419	909	914	922	rBV	195103	436543	100.00%	14.984%
12	10.517	922	924	929	rVB2	6615	13454	3.08%	0.462%
13	12.378	1109	1113	1118	rBV	223994	359310	82.31%	12.333%
14	13.559	1230	1233	1238	rVB	203961	283564	64.96%	9.733%
15	14.524	1328	1331	1335	rVB	242137	286994	65.74%	9.851%

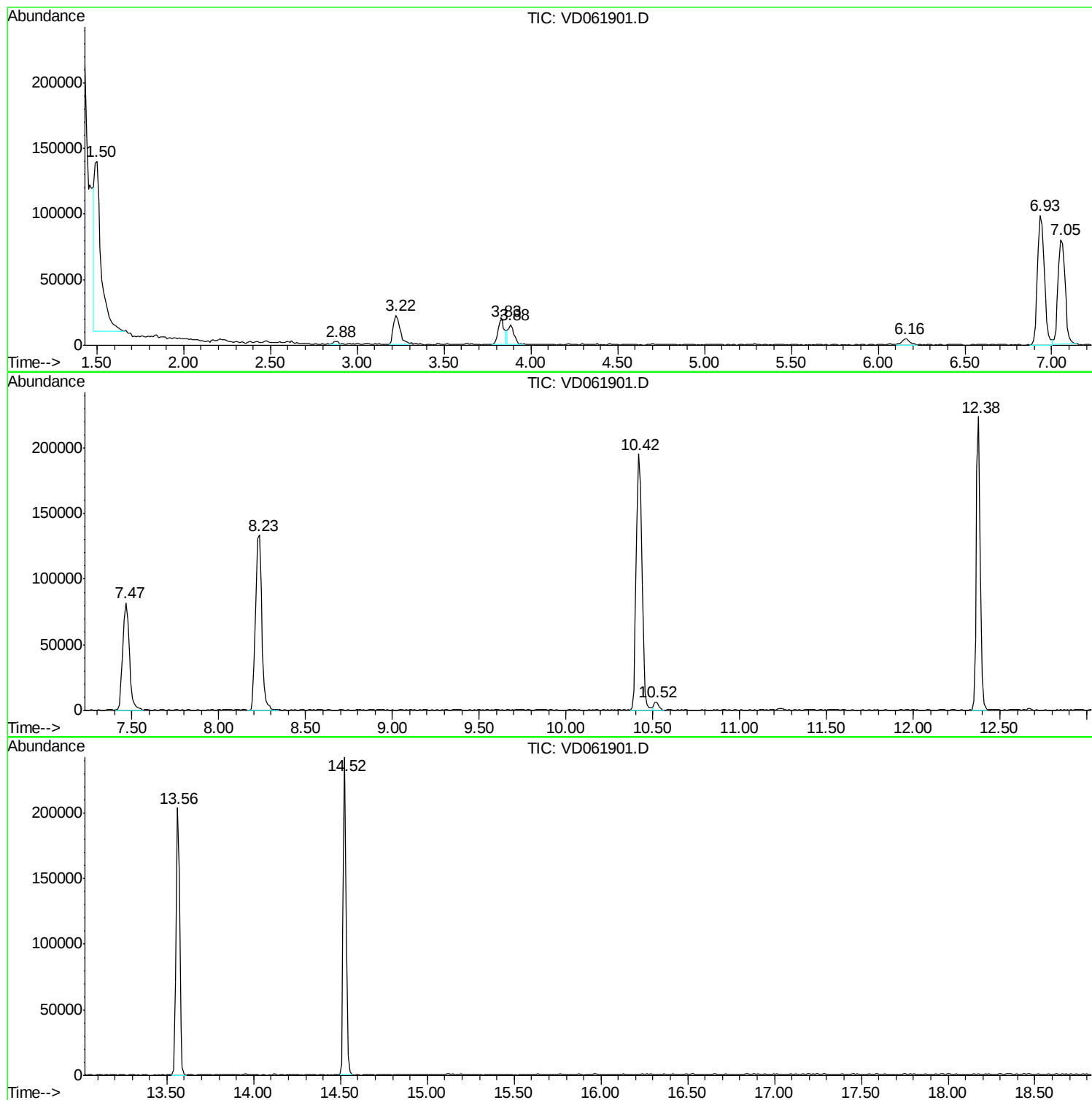
Sum of corrected areas: 2913404

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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D041519S.M
Quant Title : SW846 8260

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



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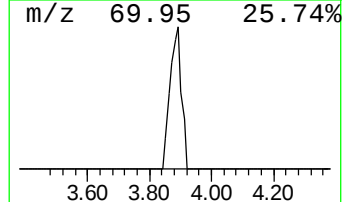
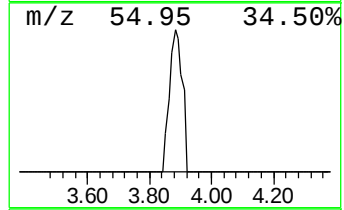
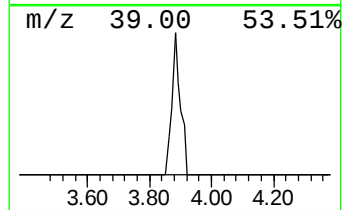
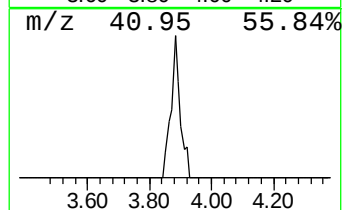
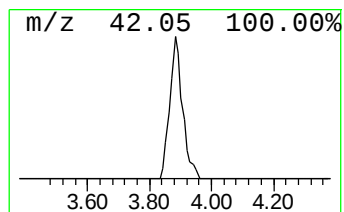
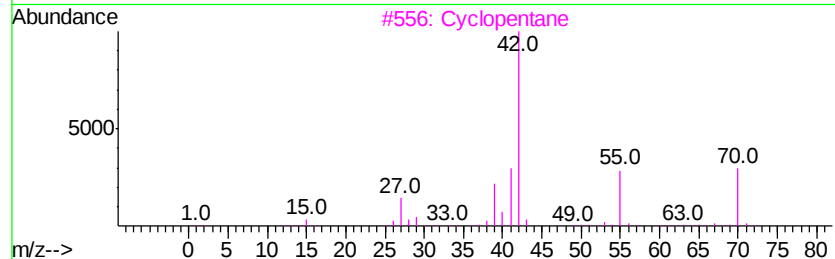
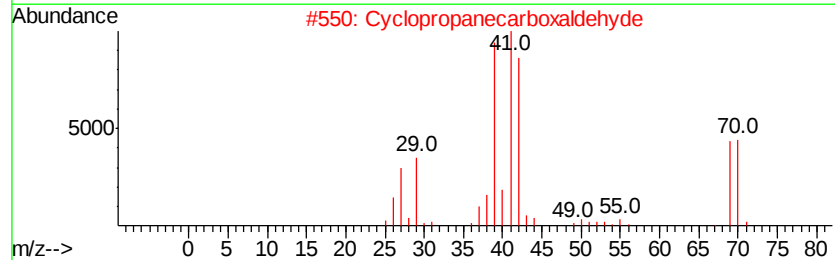
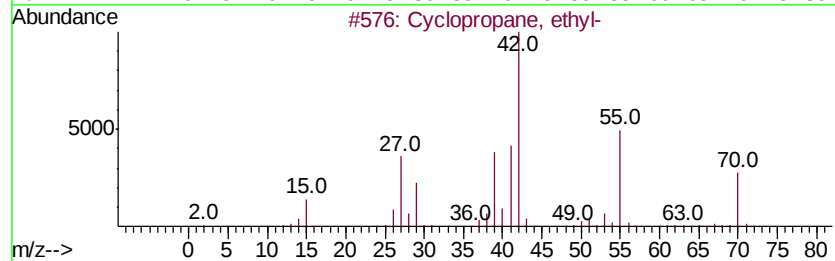
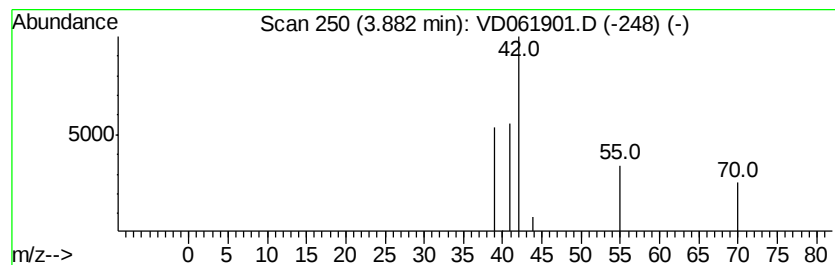
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopropane, ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.88	7.58 ug/l	33480	Pentafluorobenzene	7.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropane, ethyl-	70	C5H10	001191-96-4	64
2		Cyclopropanecarboxaldehyde	70	C4H6O	001489-69-6	9
3		Cyclopentane	70	C5H10	000287-92-3	7
4		Furan, 2,3-dihydro-	70	C4H6O	001191-99-7	7
5		Cyclobutanone	70	C4H6O	001191-95-3	4



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Cyclopropane, ethyl-	3.88	7.6	ug/l	33480	1	7.05	220825	50.0