

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD062419\
 Data File : VD062868.D
 Acq On : 25 Jun 2019 6:38
 Operator : VA/AP
 Sample : K3164-07
 Misc : 4.73g/5ml/MSVOA D/SOIL
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 BAT-B09

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\82D061919S.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	1.489	3	6	26	rVB	445264	1378688	51.35%	7.111%
2	3.231	178	183	188	rBV	43687	110935	4.13%	0.572%
3	3.822	238	243	244	rBV	38145	72347	2.69%	0.373%
4	3.881	245	249	256	rVB2	86613	248027	9.24%	1.279%
5	6.185	475	483	492	rBV	40484	142616	5.31%	0.736%
6	6.933	553	559	566	rBV	528816	1545124	57.54%	7.969%
7	7.061	566	572	591	rVB	747272	2173178	80.94%	11.208%
8	7.464	605	613	625	rBV	381750	1012032	37.69%	5.220%
9	8.232	684	691	706	rBV	947661	2429965	90.50%	12.532%
10	10.418	907	913	921	rBV	1168083	2685072	100.00%	13.848%
11	11.235	990	996	1003	rBV	21428	60253	2.24%	0.311%
12	12.377	1107	1112	1124	rBV	1623807	2614575	97.37%	13.485%
13	13.558	1229	1232	1240	rVB	1554692	2178060	81.12%	11.233%
14	13.715	1245	1248	1252	rVB	66768	100102	3.73%	0.516%
15	14.523	1327	1330	1338	rVB	2083739	2587038	96.35%	13.343%
16	15.113	1387	1390	1395	rBV2	35053	51305	1.91%	0.265%

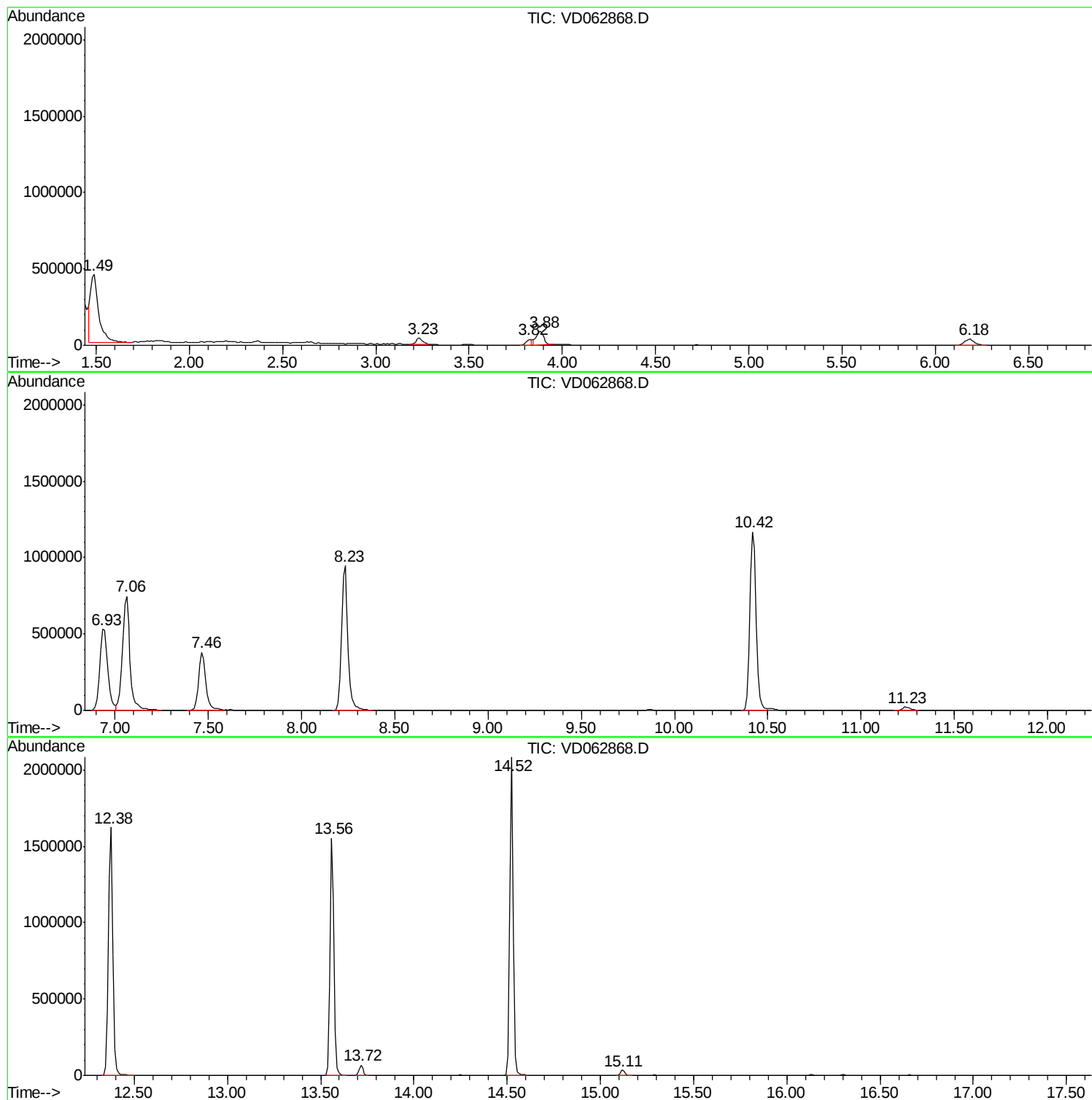
Sum of corrected areas: 19389317

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061919S.M
Quant Title : SW846 8260

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



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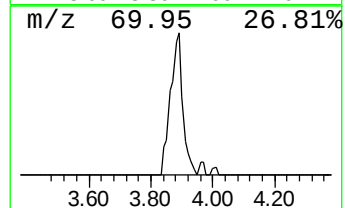
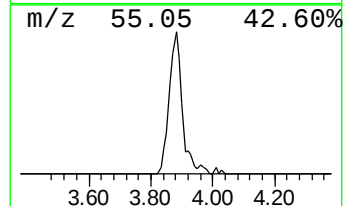
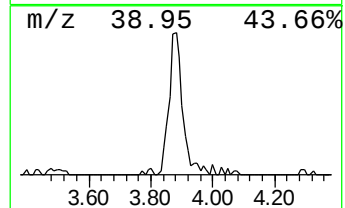
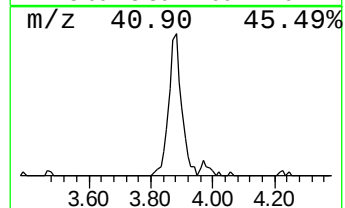
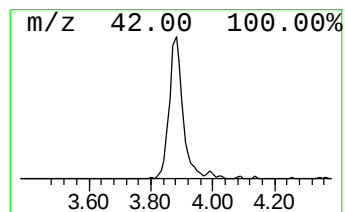
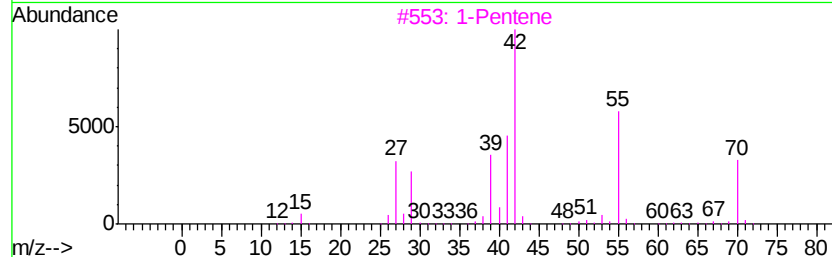
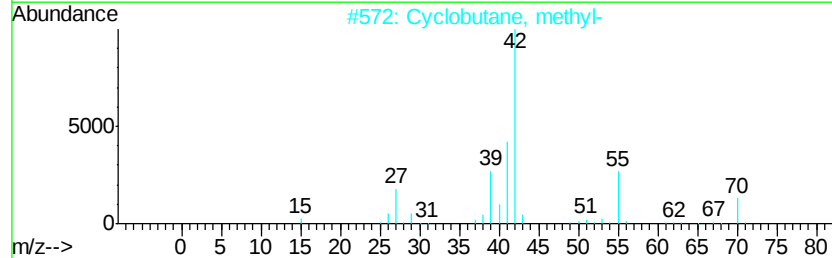
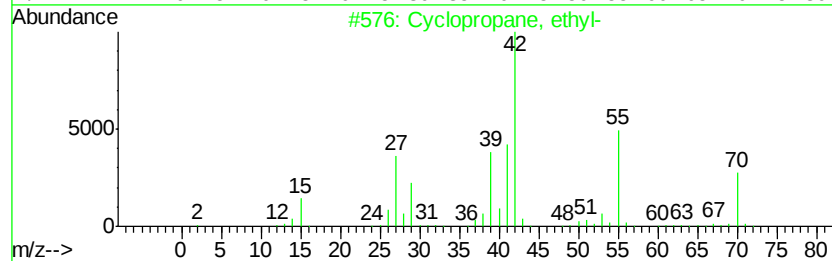
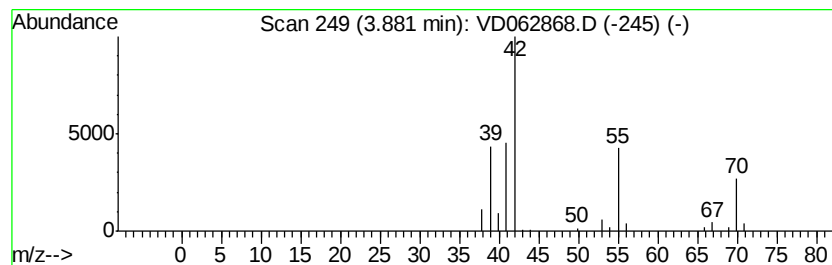
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopropane, ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.88	5.71 ug/l	248027	Pentafluorobenzene	7.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropane, ethyl-	70	C5H10	001191-96-4	86
2		Cyclobutane, methyl-	70	C5H10	000598-61-8	53
3		1-Pentene	70	C5H10	000109-67-1	52
4		Cyclopentane	70	C5H10	000287-92-3	9
5		Ethoxyacetylene	70	C4H6O	000927-80-0	4



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclopropane, ethyl-	3.88	5.7	ug/l	248027	1	7.06	2173180	50.0