

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD081320\
 Data File : VD066217.D
 Acq On : 13 Aug 2020 18:23
 Operator : VA/SY
 Sample : L3508-07
 Misc : 4.12G/5.00ml/MSVOA D/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TR-05-081120

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\82D080520S.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	3.032	182	189	197	rVB	20018	39969	2.00%	0.361%
2	3.403	243	252	260	rBV4	9989	23799	1.19%	0.215%
3	4.973	508	519	520	rBV4	8626	24308	1.22%	0.220%
4	5.055	520	533	547	rVB2	250264	870228	43.57%	7.869%
5	5.985	682	691	701	rVB6	12232	36693	1.84%	0.332%
6	7.191	886	896	906	rBV	37490	104471	5.23%	0.945%
7	7.908	1006	1018	1023	rBV2	220310	516564	25.86%	4.671%
8	7.973	1023	1029	1040	rVB	409518	918231	45.97%	8.304%
9	8.326	1078	1089	1100	rBV	168236	382489	19.15%	3.459%
10	8.861	1171	1180	1192	rBV	568661	1162841	58.22%	10.516%
11	10.338	1422	1431	1449	rBV	908460	1655038	82.86%	14.966%
12	11.643	1643	1653	1667	rVB	1029884	1772494	88.74%	16.029%
13	12.632	1814	1821	1830	rBV	952471	1553739	77.79%	14.050%
14	13.579	1974	1982	1990	rBV	1230101	1997423	100.00%	18.063%

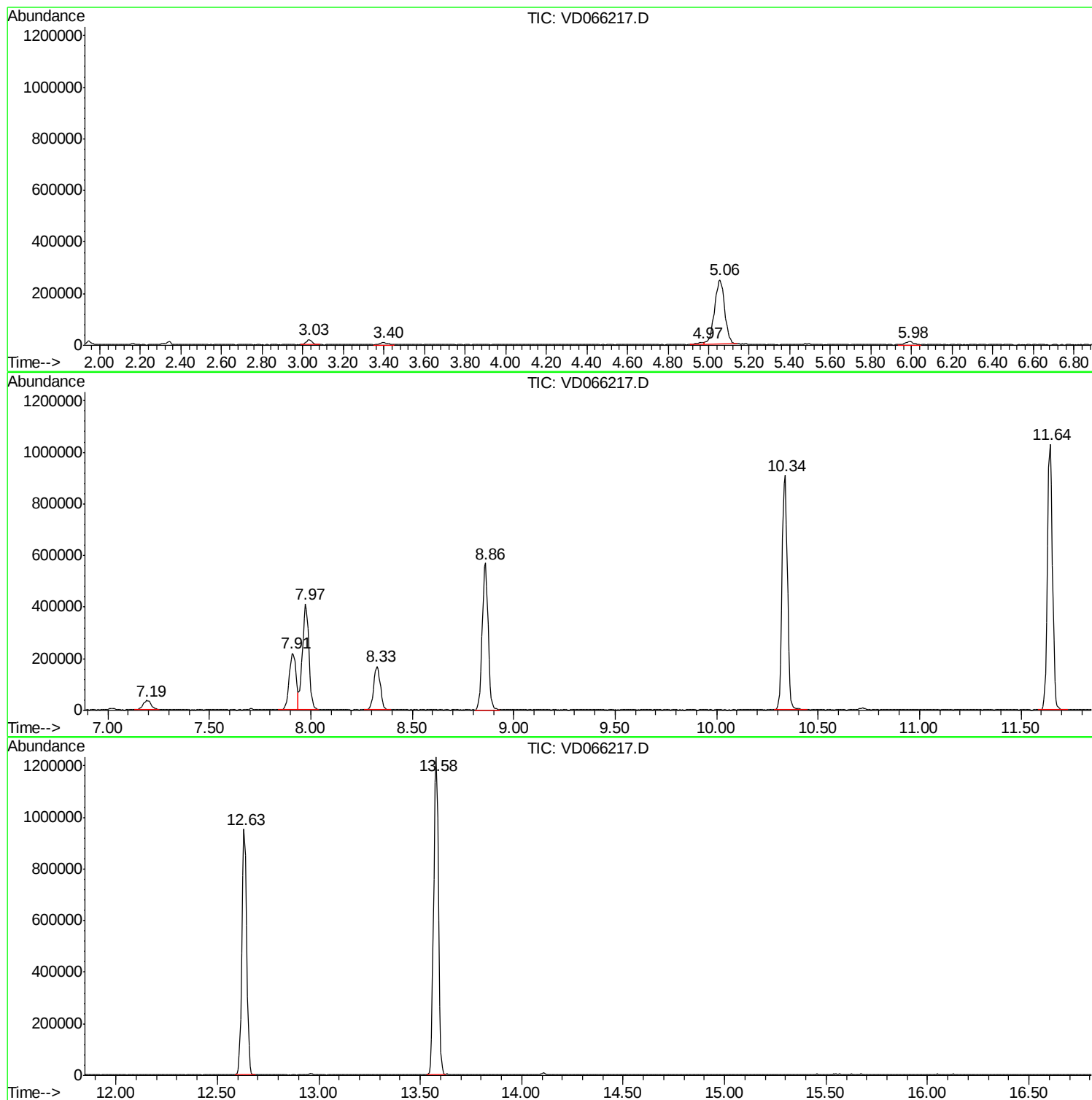
Sum of corrected areas: 11058287

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Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_D\METHOD\82D080520S.M
Quant Title : SW846 8260

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



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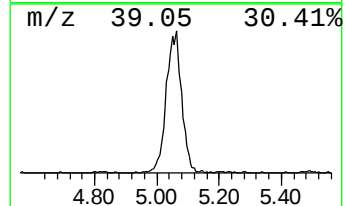
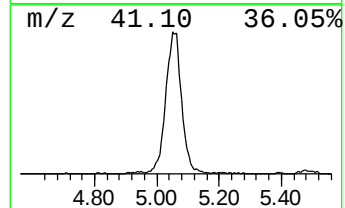
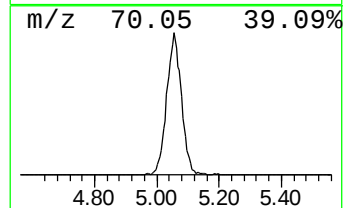
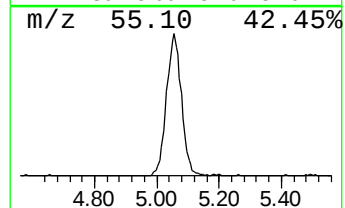
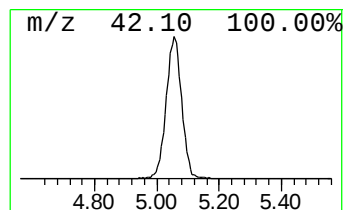
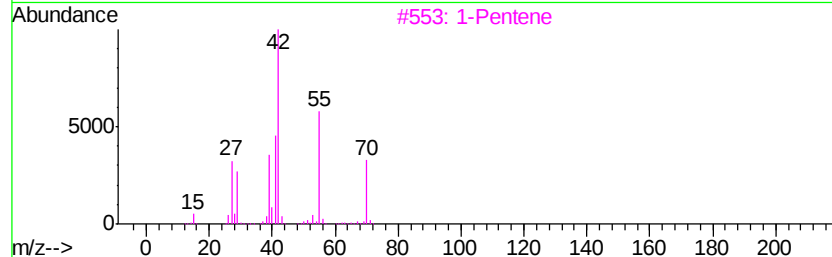
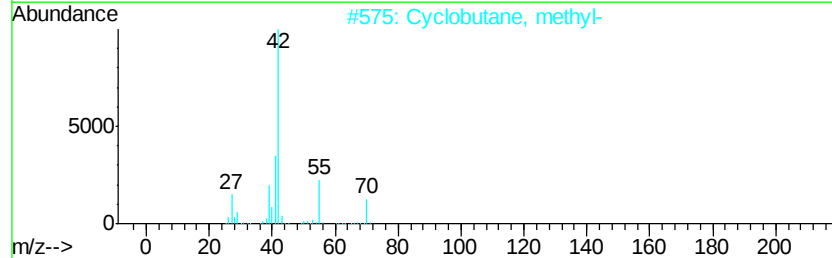
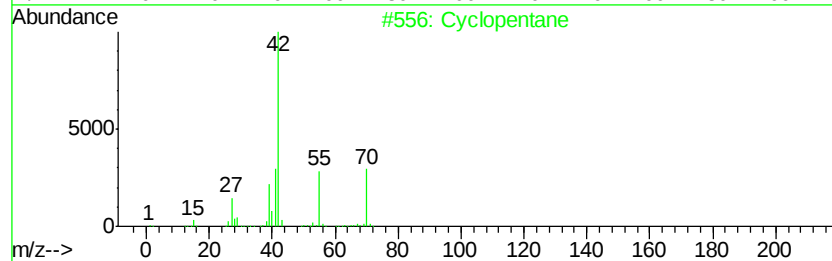
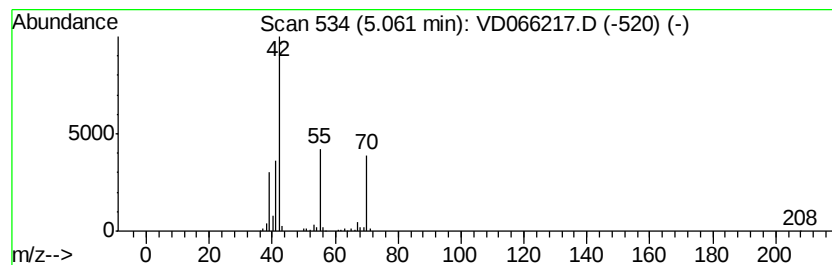
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Cyclopentane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	47.39 ug/l	870228	Pentafluorobenzene	7.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane	70	C5H10	000287-92-3	86
2		Cyclobutane, methyl-	70	C5H10	000598-61-8	78
3		1-Pentene	70	C5H10	000109-67-1	64
4		Cyclopropane, ethyl-	70	C5H10	001191-96-4	37
5		Methyl propargyl ether	70	C4H6O	000627-41-8	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclopentane	5.06	47.4	ug/l	870228	1	7.97	918231	50.0