

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD091020\
 Data File : VD066635.D
 Acq On : 11 Sep 2020 02:12
 Operator : VA/SY
 Sample : L3937-04
 Misc : 6.01G/5.00ml/MSVOA D/SOIL
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 BSY-SB-03-5.0-6.0

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\82D090320S.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	3.709	294	304	312	rVB5	7387	21588	1.01%	0.183%
2	4.156	372	380	387	rBV	15187	35538	1.66%	0.301%
3	4.967	506	518	525	rBV3	59188	190938	8.92%	1.618%
4	5.056	525	533	547	rVB2	54773	185206	8.65%	1.569%
5	5.432	588	597	607	rBV	9136	24829	1.16%	0.210%
6	7.197	886	897	908	rBV	24576	76354	3.57%	0.647%
7	7.914	1009	1019	1024	rBV	308027	743407	34.72%	6.300%
8	7.979	1024	1030	1039	rVB	512877	1132979	52.91%	9.601%
9	8.326	1077	1089	1100	rBV2	286394	648148	30.27%	5.492%
10	8.861	1171	1180	1191	rBV	768041	1575123	73.56%	13.347%
11	10.338	1422	1431	1439	rBV	1205054	2141386	100.00%	18.146%
12	10.714	1488	1495	1503	rBV3	13931	27106	1.27%	0.230%
13	11.643	1644	1653	1665	rBV	1094120	1912803	89.33%	16.209%
14	12.632	1813	1821	1830	rBV	837198	1355179	63.29%	11.484%
15	13.579	1974	1982	1990	rBV	1065397	1705568	79.65%	14.453%
16	15.555	2312	2318	2324	rVB4	12180	24783	1.16%	0.210%

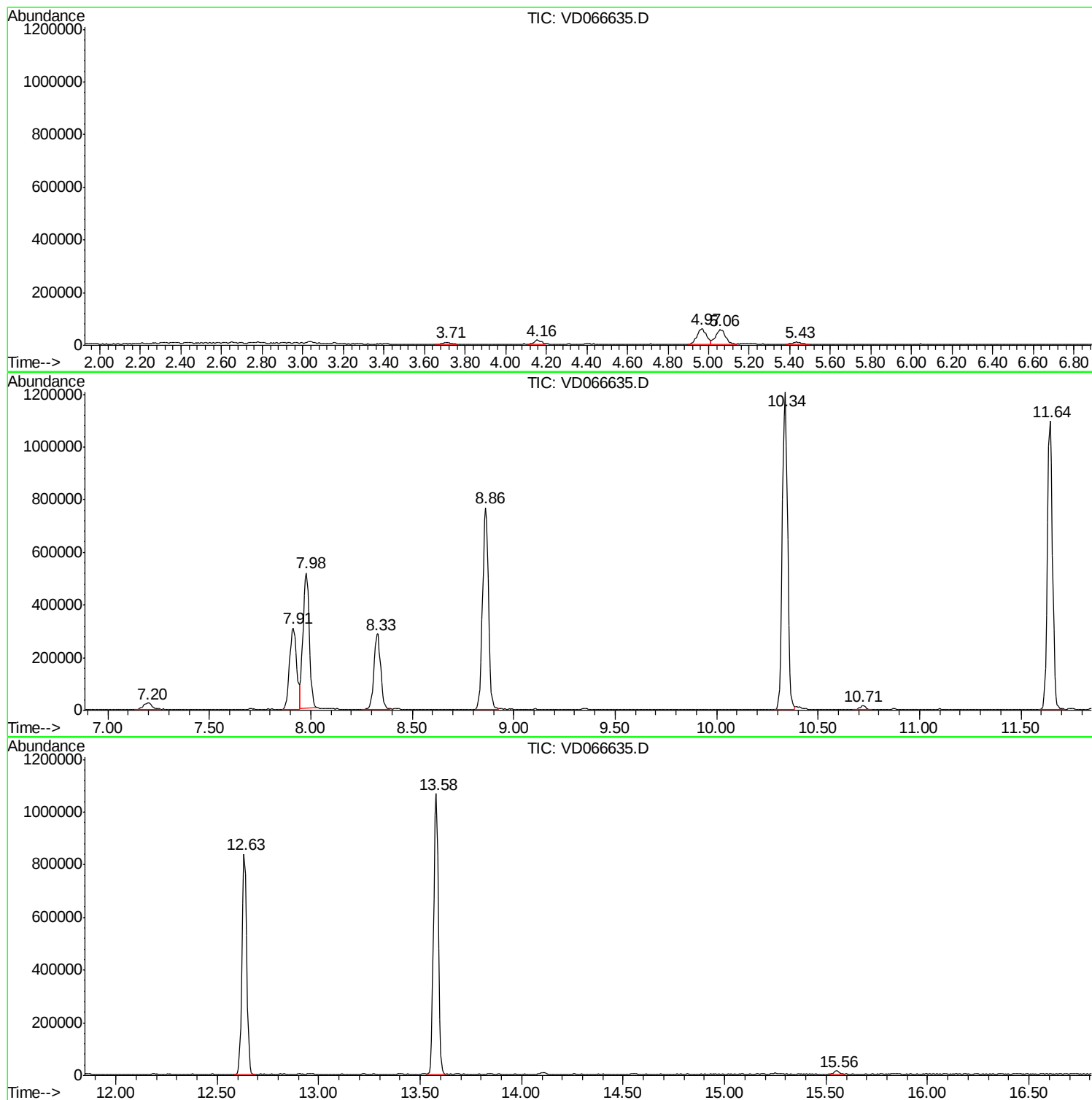
Sum of corrected areas: 11800935

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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



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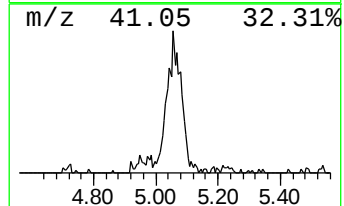
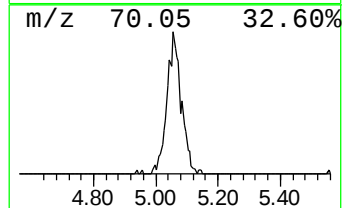
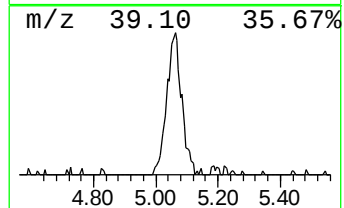
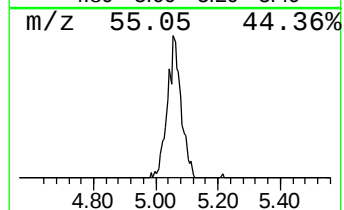
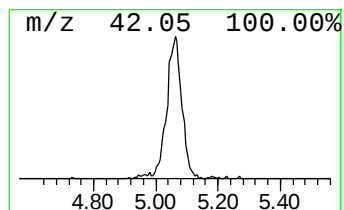
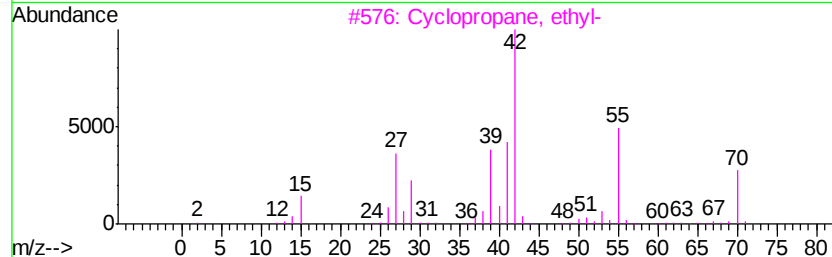
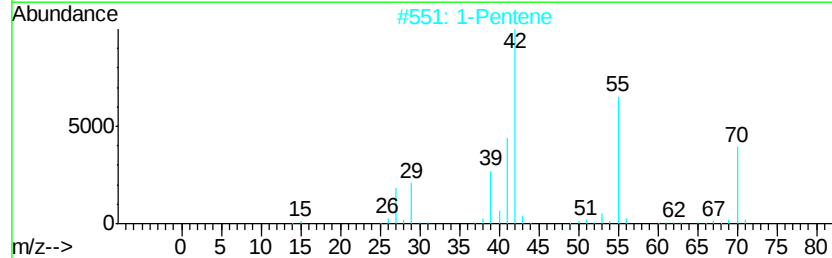
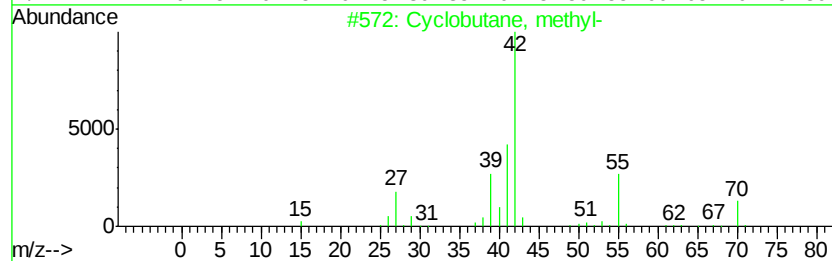
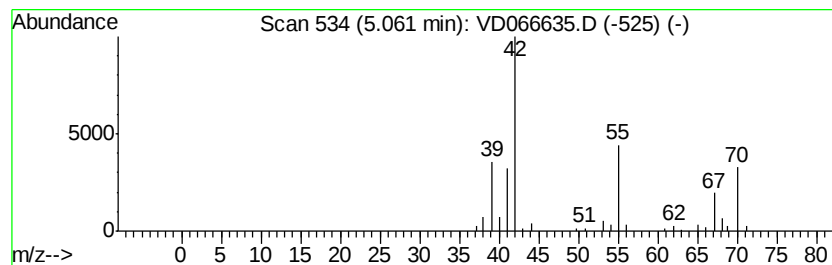
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclobutane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	8.17 ug/l	185206	Pentafluorobenzene	7.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclobutane, methyl-	70	C5H10	000598-61-8	72
2		1-Pentene	70	C5H10	000109-67-1	72
3		Cyclopropane, ethyl-	70	C5H10	001191-96-4	64
4		Cyclopentane	70	C5H10	000287-92-3	64
5		1-Aziridinepropanenitrile, .beta...	110	C6H10N2	004078-19-7	40



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Cyclobutane, methyl-	5.06	8.2	ug/l	185206	1	7.98	1132980	50.0