

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW081218\
 Data File : VW004583.D
 Acq On : 10 Aug 2018 05:52
 Operator : SY/AP
 Sample : J4382-03
 Misc : 17.57G/5ML/MSVOA W/SOIL
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 JTY-SB-03-5-7

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_W\METHOD\82W080618S.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	1.813	29	37	58	rVB	5407049	12233786	100.00%	58.797%
2	2.819	195	202	213	rBV	62209	158898	1.30%	0.764%
3	4.916	532	546	551	rBV2	48567	163891	1.34%	0.788%
4	5.001	551	560	584	rVB2	131790	520870	4.26%	2.503%
5	7.153	900	913	931	rBV	113855	342691	2.80%	1.647%
6	7.513	961	972	990	rBV	496078	1256135	10.27%	6.037%
7	7.952	1025	1044	1057	rBV	589504	1309448	10.70%	6.293%
8	8.848	1181	1191	1202	rBV	845056	1665929	13.62%	8.007%
9	11.634	1641	1648	1657	rBV	917582	1546921	12.64%	7.435%
10	13.567	1958	1965	1975	rBV	894596	1415014	11.57%	6.801%
11	14.707	2137	2152	2154	rBV4	60214	193358	1.58%	0.929%

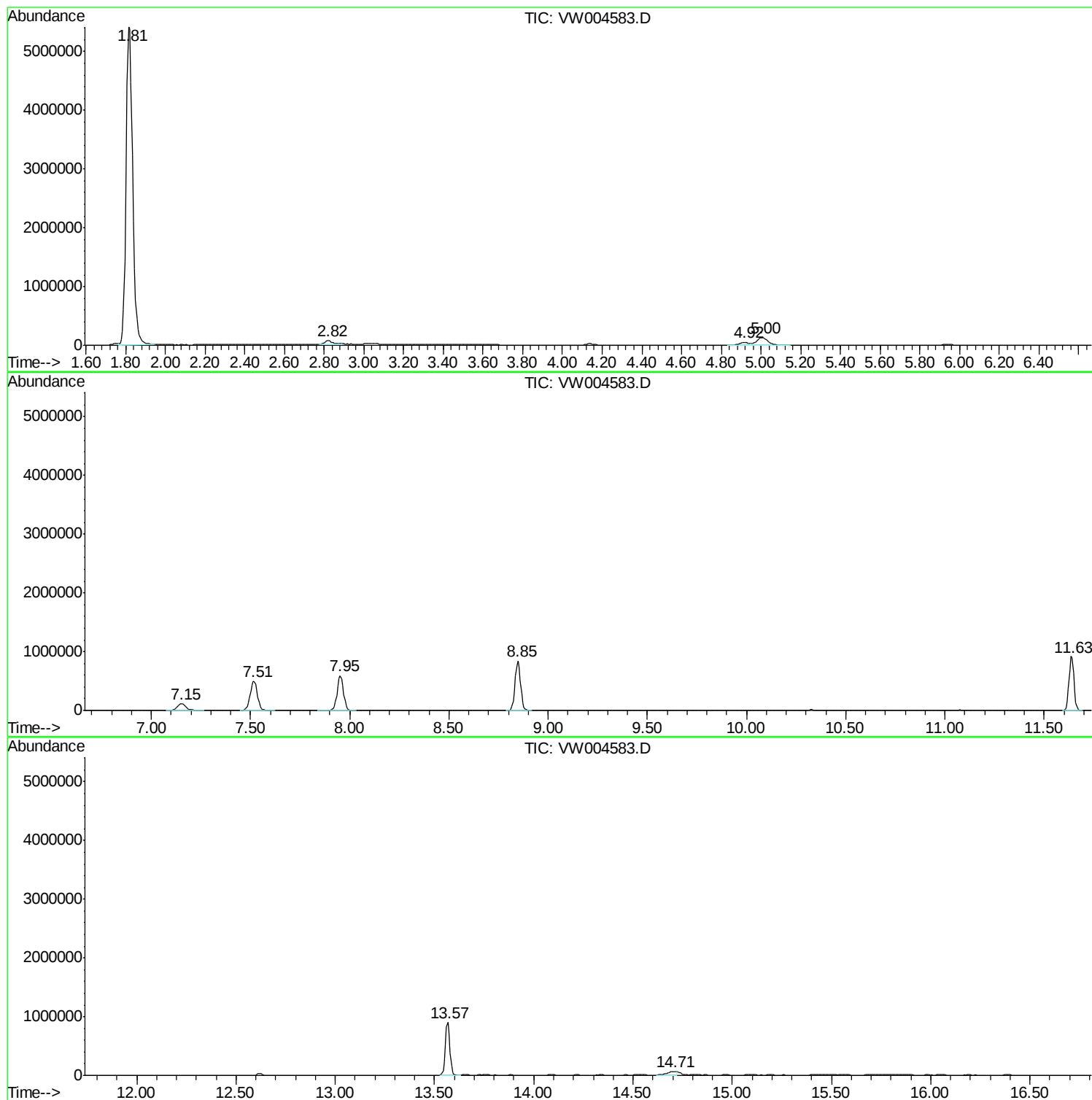
Sum of corrected areas: 20806941

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W080618S.M
Quant Title : SW846 8260

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



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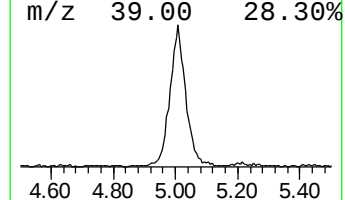
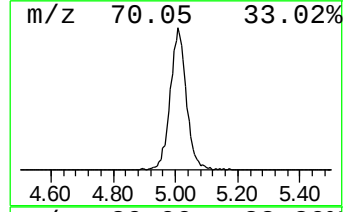
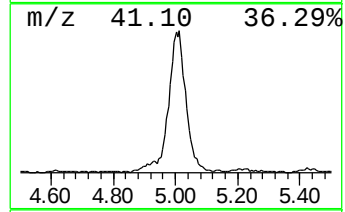
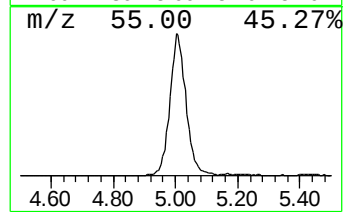
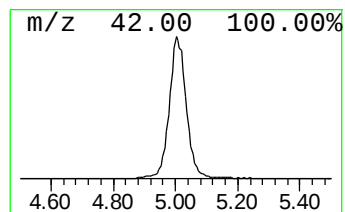
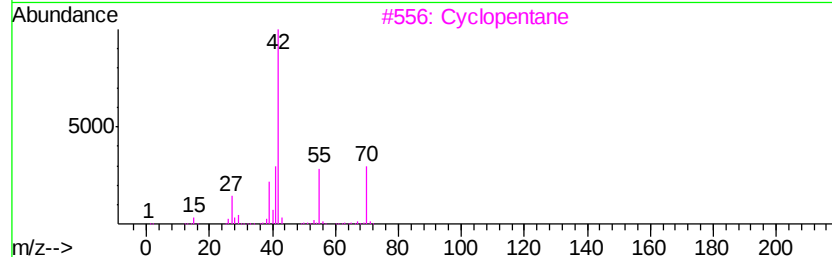
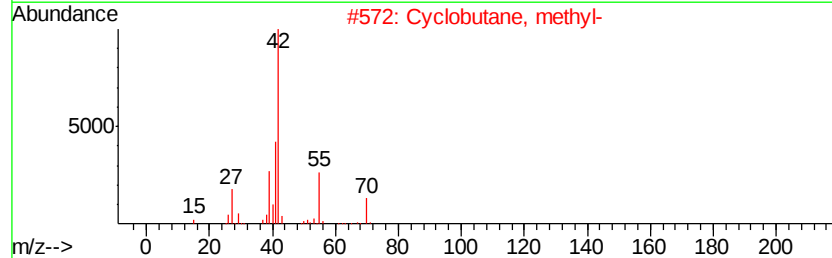
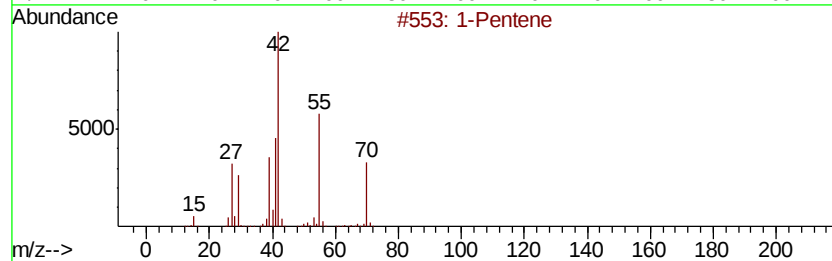
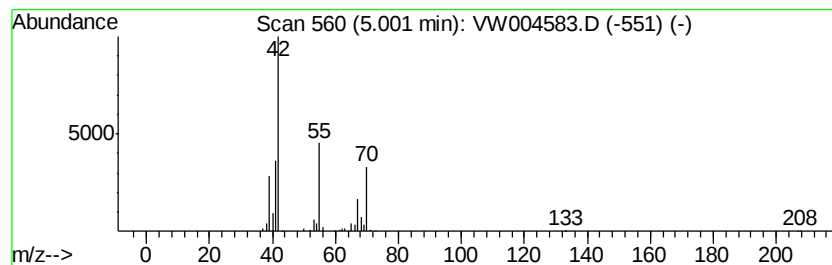
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W080618S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Pentene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	19.89 ug/l	520870	Pentafluorobenzene	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentene	70	C5H10	000109-67-1	64
2		Cyclobutane, methyl-	70	C5H10	000598-61-8	64
3		Cyclopentane	70	C5H10	000287-92-3	64
4		Cyclopropane, ethyl-	70	C5H10	001191-96-4	64
5		1-Aziridinepropanenitrile, .beta...	110	C6H10N2	004078-19-7	39



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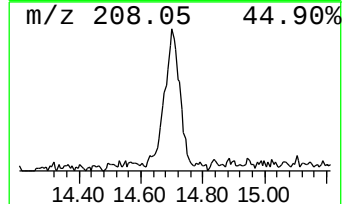
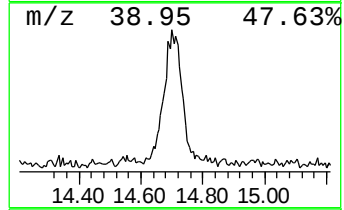
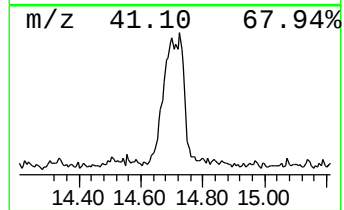
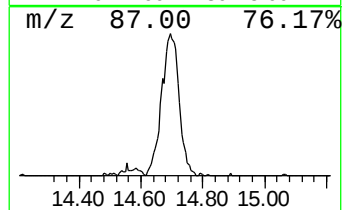
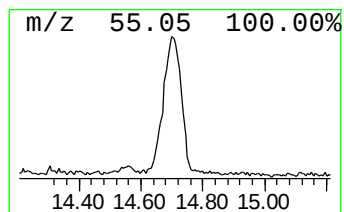
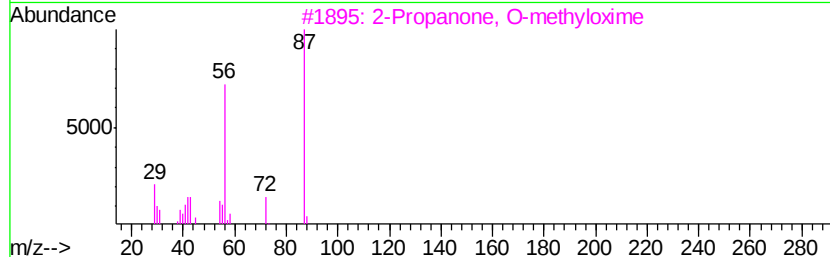
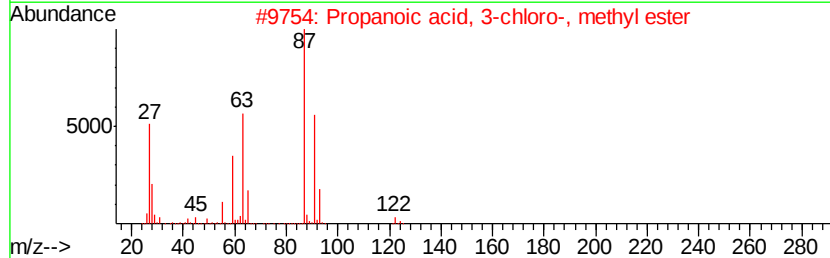
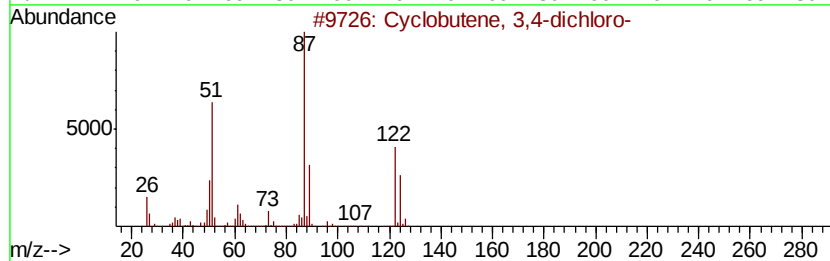
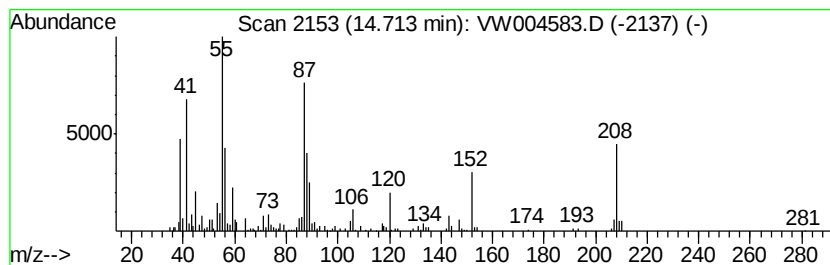
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Peak Number 6 unknown14.71 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	6.83 ug/l	193358	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclobutene, 3,4-dichloro-	122	C4H4Cl2	041326-64-1	25
2		Propanoic acid, 3-chloro-, methy...	122	C4H7ClO2	006001-87-2	22
3		2-Propanone, O-methyloxime	87	C4H9NO	003376-35-0	16
4		Hydrazine, 1,2-dipropyl-	116	C6H16N2	001615-83-4	14
5		Dimethoate	229	C5H12N03PS2	000060-51-5	14



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Pentene	5.00	19.9	ug/l	520870	1	7.95	1309450	50.0
unknown14.71	14.71	6.8	ug/l	193358	4	13.57	1415010	50.0