

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD062218\
 Data File : VD058336.D
 Acq On : 23 Jun 2018 11:47
 Operator : VA/AP
 Sample : J3642-05
 Misc : 6.31g/5ml/MSVOA D/SOIL
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TEST-PIT-2

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\82D061918S.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	1.237	94	103	105	rBV2	1029011	2750841	56.23%	21.777%
2	1.296	105	109	125	rVB	1356234	4892310	100.00%	38.729%
3	2.625	239	244	252	rBV2	18643	59623	1.22%	0.472%
4	4.859	460	471	479	rBV2	21014	75401	1.54%	0.597%
5	5.568	535	543	548	rBV	222526	628394	12.84%	4.975%
6	5.657	548	552	569	rVB	221763	634420	12.97%	5.022%
7	6.031	584	590	604	rBV	200860	540106	11.04%	4.276%
8	6.700	652	658	673	rBV	305573	749509	15.32%	5.933%
9	8.718	857	863	870	rBV	384728	839176	17.15%	6.643%
10	10.697	1059	1064	1074	rBV	396110	679260	13.88%	5.377%
11	11.937	1186	1190	1198	rVB	314643	437869	8.95%	3.466%
12	12.912	1286	1289	1295	rBV	267425	345144	7.05%	2.732%

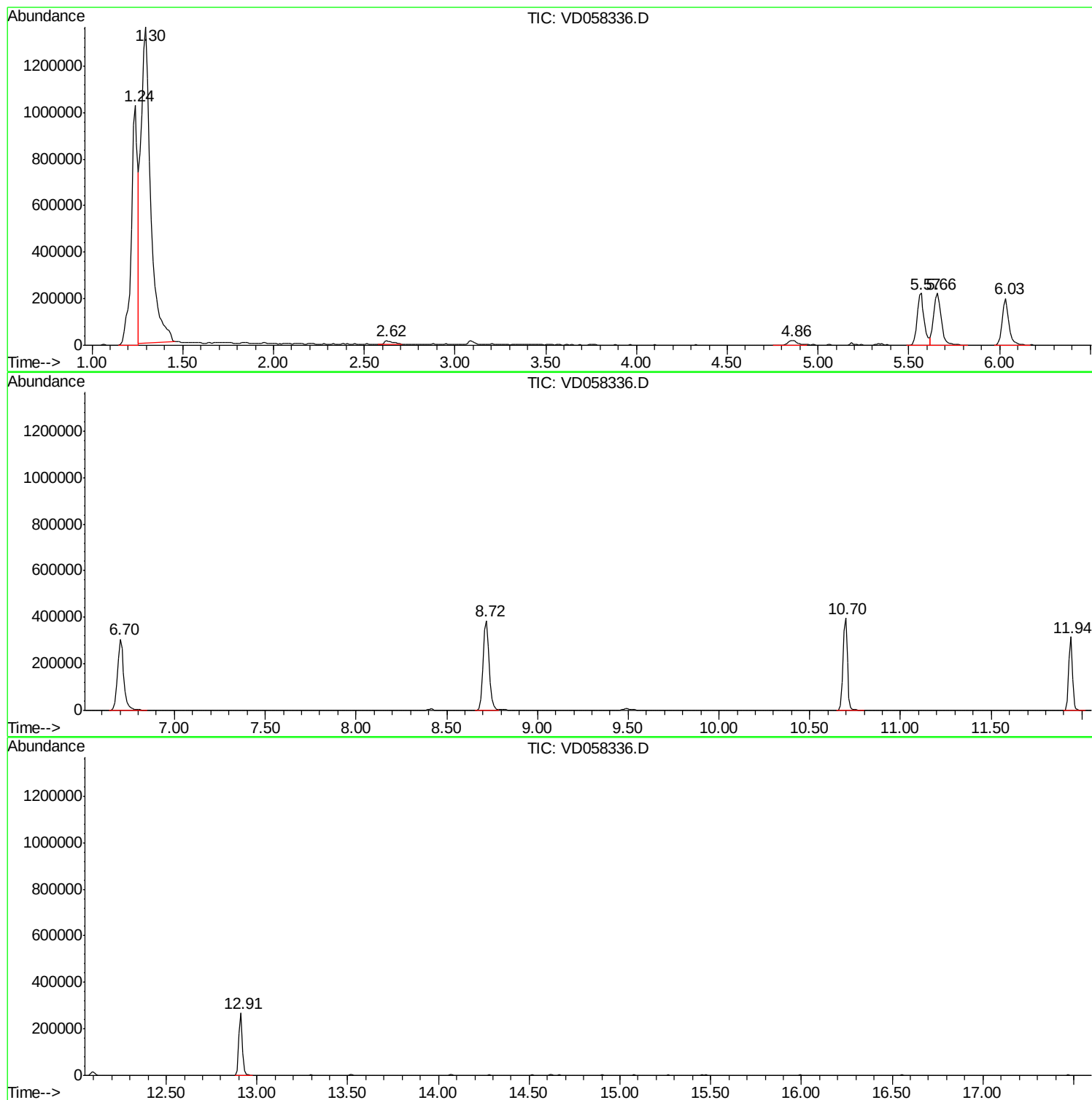
Sum of corrected areas: 12632053

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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



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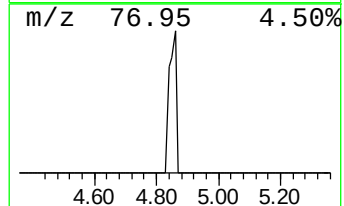
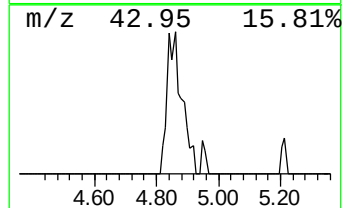
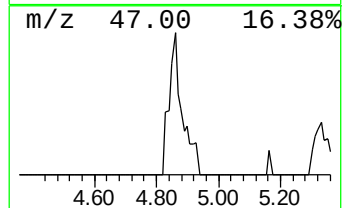
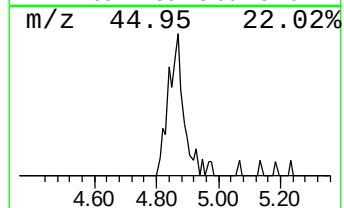
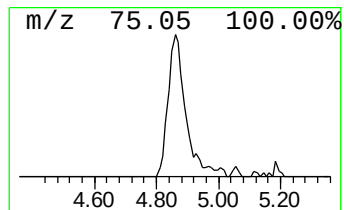
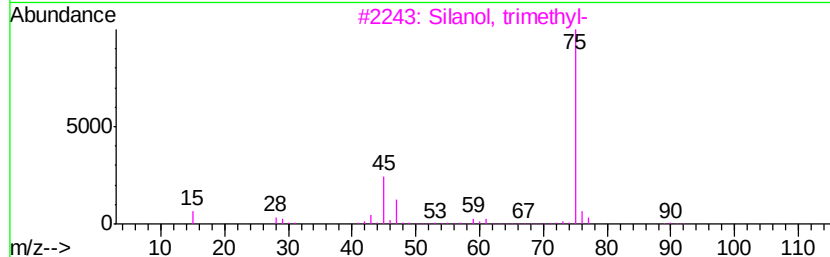
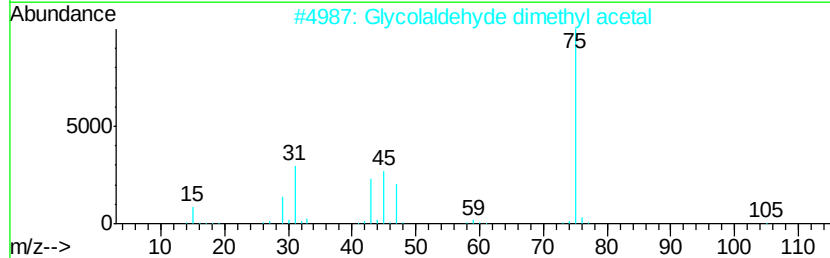
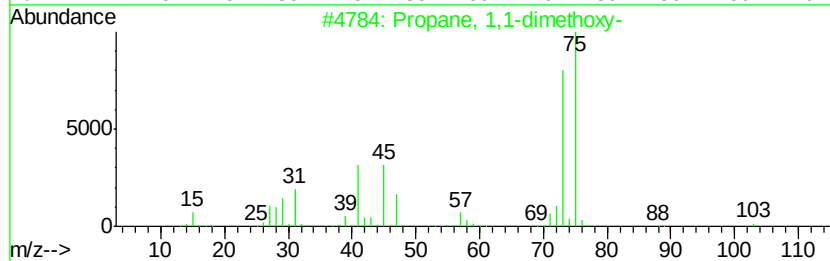
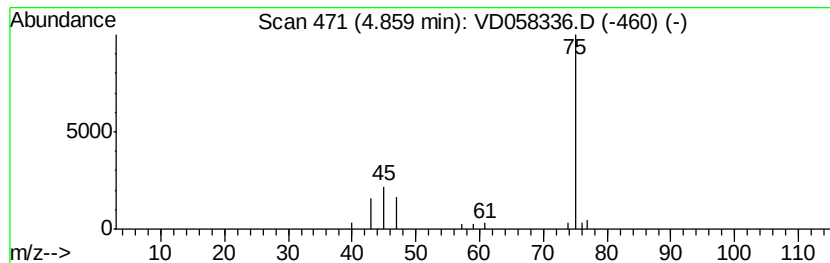
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 3 Propane, 1,1-dimethoxy- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	5.94 ug/l	75401	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	64
2		Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	43
3		Silanol, trimethyl-	90	C3H10OSi	001066-40-6	43
4		Formamide, N-methylthio	75	C2H5NS	018952-41-5	7
5		Ethanethioamide	75	C2H5NS	000062-55-5	7



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-dime...	4.86	5.9	ug/l	75401	1	5.66	634420	50.0