

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU080618\
 Data File : VU025883.D
 Acq On : 06 Aug 2018 14:10
 Operator : MD/SY
 Sample : J4257-04
 Misc : 9.93g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 FK-BPM-03-001-BRE

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_U\METHOD\82U072718W.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	1.082	135	153	171	rBV	751802	828968	73.84%	11.957%
2	1.609	282	317	318	rBV6	29744	94706	8.44%	1.366%
3	2.831	685	697	716	rBV2	13978	32978	2.94%	0.476%
4	4.883	1318	1335	1351	rBV	172775	403846	35.97%	5.825%
5	4.979	1351	1365	1390	rVB	233080	516519	46.01%	7.450%
6	5.307	1453	1467	1493	rBV	175227	385278	34.32%	5.557%
7	5.886	1632	1647	1667	rBV	334304	673754	60.01%	9.718%
8	7.567	2156	2170	2191	rBV	635656	1113730	99.20%	16.064%
9	9.095	2630	2645	2668	rBV	513988	874401	77.89%	12.612%
10	10.313	3011	3024	3049	rBV	553602	886174	78.93%	12.782%
11	11.490	3377	3390	3411	rBV	715061	1122679	100.00%	16.193%

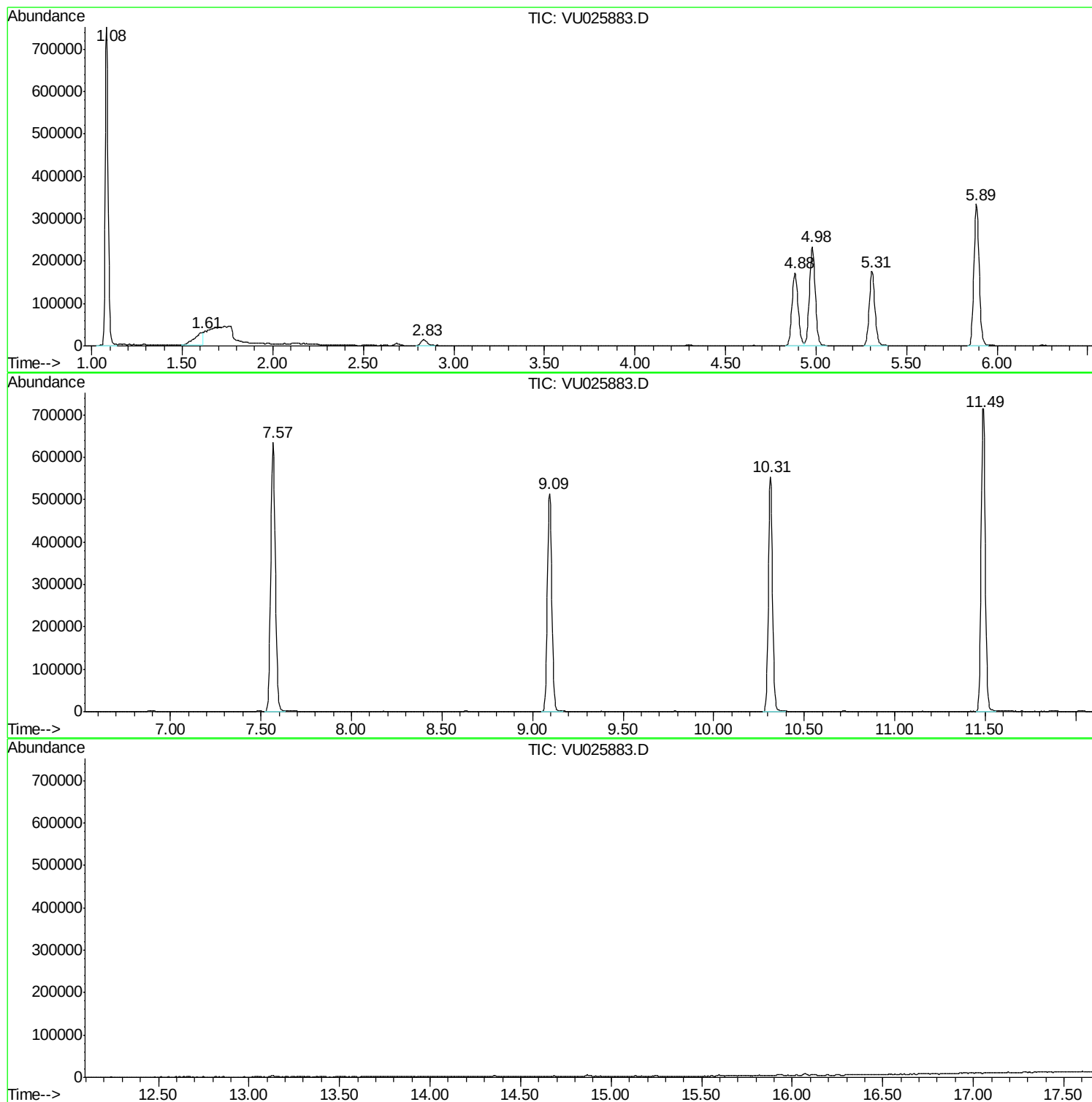
Sum of corrected areas: 6933033

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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



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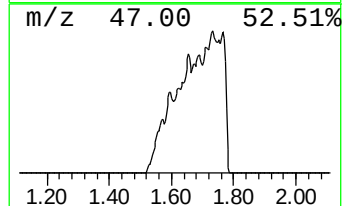
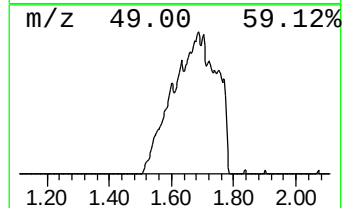
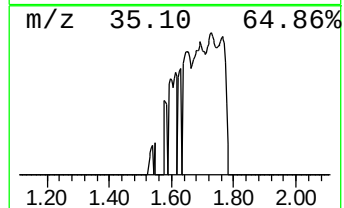
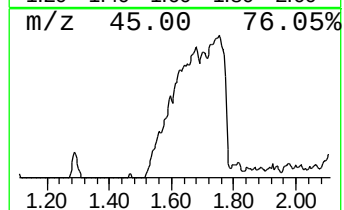
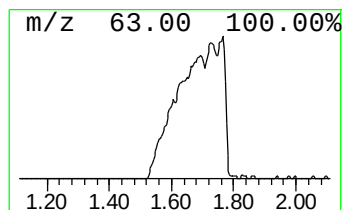
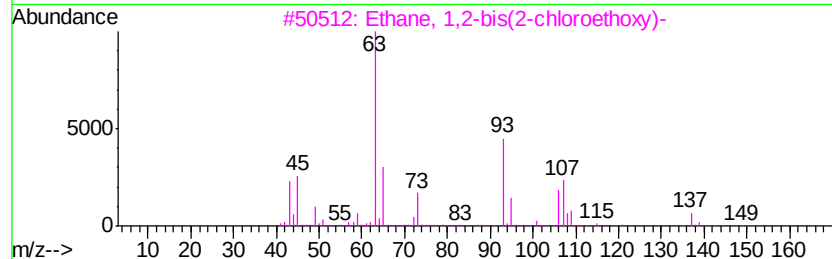
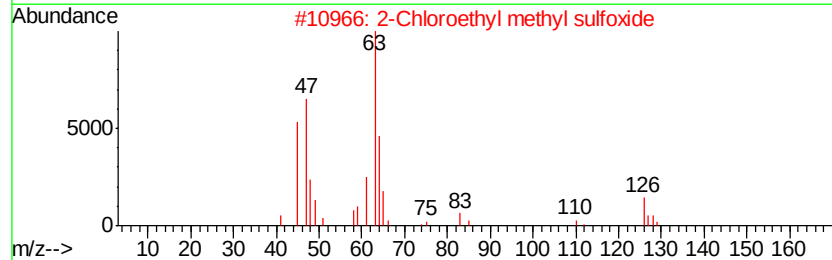
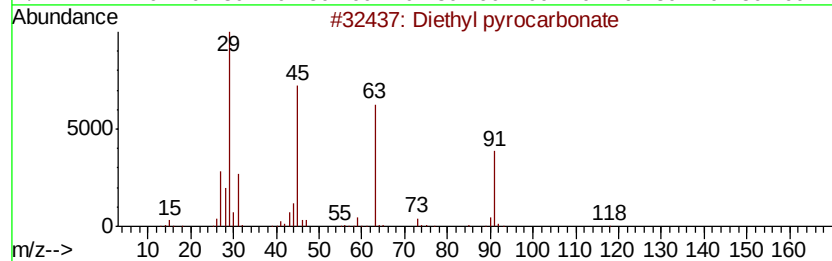
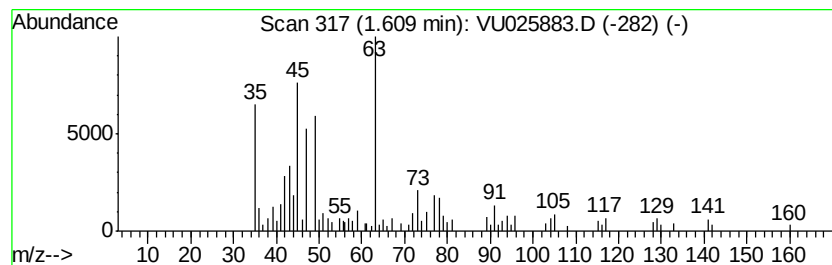
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Quant Title : SW846 8260

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

Peak Number 2 unknown1.61 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
1.61	9.17 ug/l	94706	Pentafluorobenzene	4.98		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Diethyl pyrocarbonate	162	C6H10O5	001609-47-8	38
2		2-Chloroethyl methyl sulfoxide	126	C3H7ClOS	005331-57-7	28
3		Ethane, 1,2-bis(2-chloroethoxy)-	186	C6H12Cl2O2	000112-26-5	25
4		Ethene, (methylsulfonyl)-	106	C3H6O2S	003680-02-2	16
5		Methane, chloromethoxy-	80	C2H5ClO	000107-30-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.61	1.61	9.2	ug/l	94706	1	4.98	516519	50.0