

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110918\
 Data File : VU028037.D
 Acq On : 09 Nov 2018 12:18
 Operator : MD/SY
 Sample : J5762-01
 Misc : 4.07g/5mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BU-01-110518

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\82U110218W.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	1.624	104	140	146	rBV2	61157	222968	18.35%	3.426%
2	2.627	441	452	464	rBV	6922	13475	1.11%	0.207%
3	2.836	504	517	530	rBV	8657	17976	1.48%	0.276%
4	4.881	1137	1153	1170	rBV	144859	354585	29.18%	5.449%
5	4.977	1170	1183	1208	rVB	232706	539569	44.40%	8.291%
6	5.305	1270	1285	1312	rBV2	184288	421305	34.67%	6.474%
7	5.884	1447	1465	1493	rBV2	403502	850395	69.97%	13.068%
8	7.565	1974	1988	2010	rBV	666380	1215307	100.00%	18.675%
9	9.093	2449	2463	2480	rBV	589399	1042171	85.75%	16.015%
10	10.311	2830	2842	2873	rBV	509087	825896	67.96%	12.691%
11	11.485	3195	3207	3231	rBV	634337	1003932	82.61%	15.427%

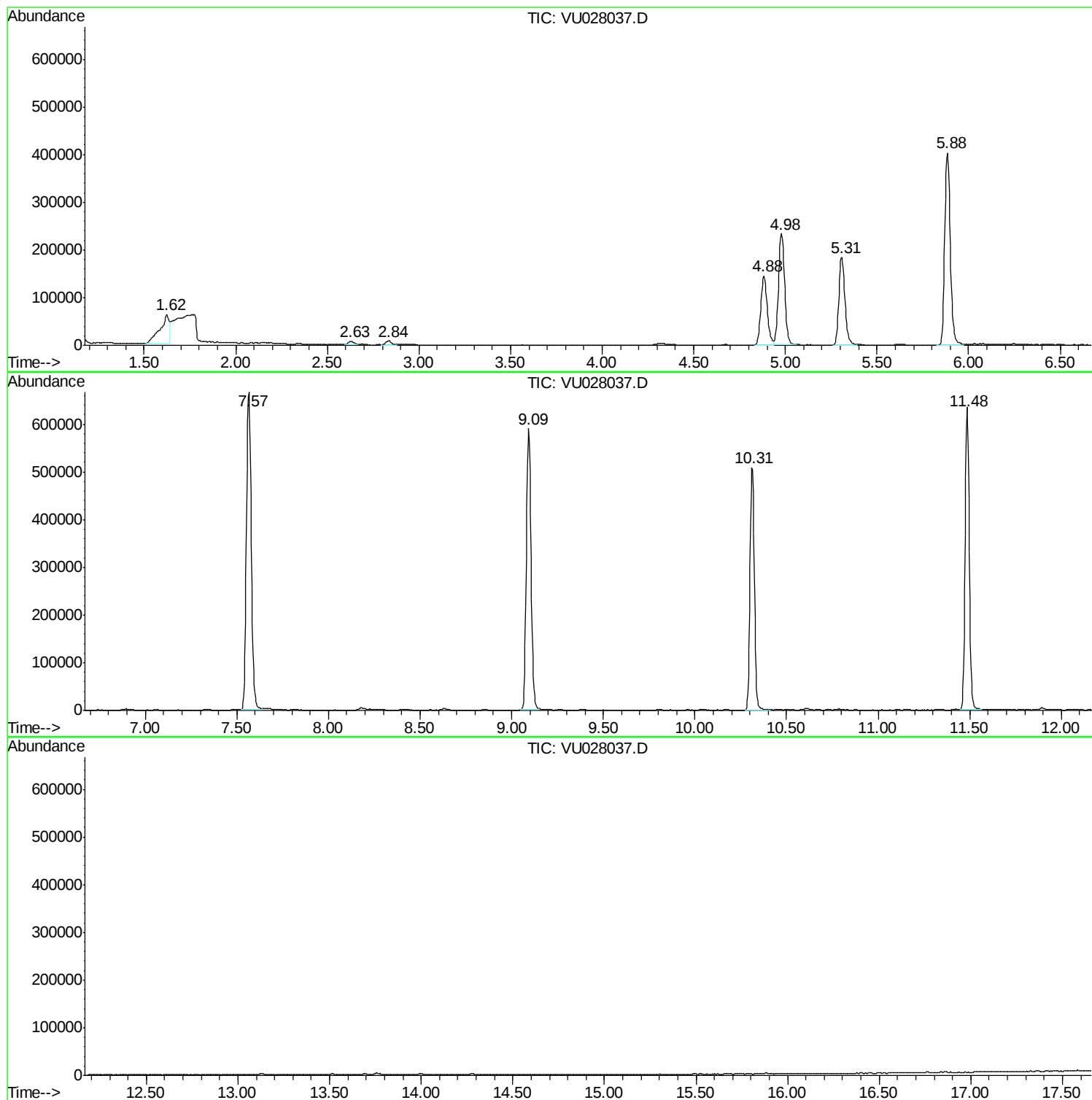
Sum of corrected areas: 6507579

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Instrument :
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ClientSampled :
BU-01-110518

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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



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ClientSampleId :
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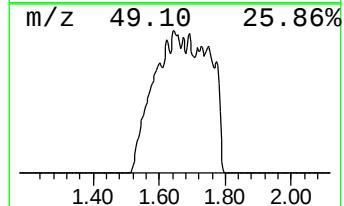
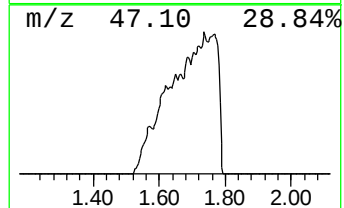
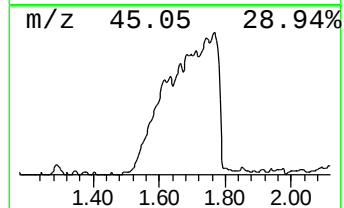
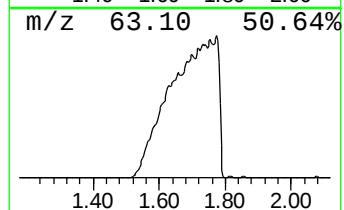
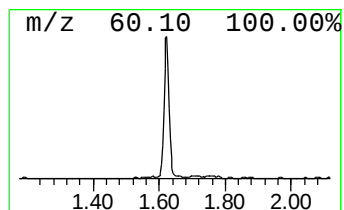
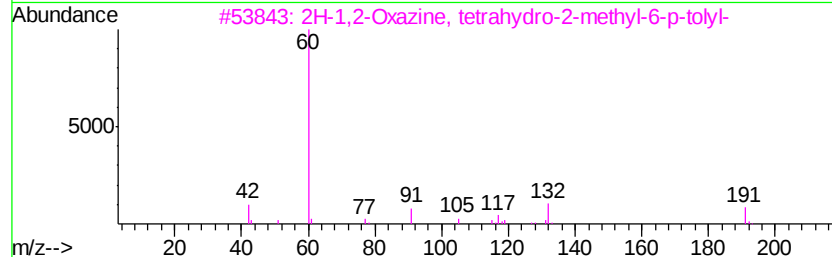
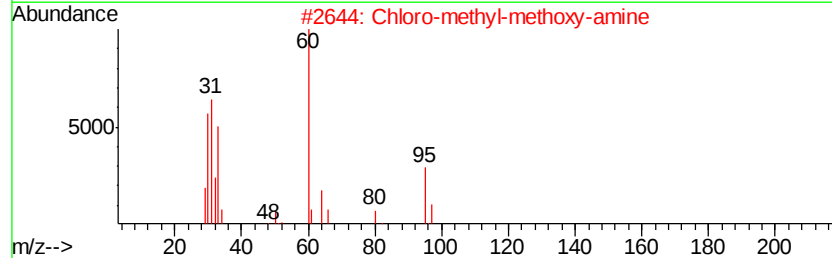
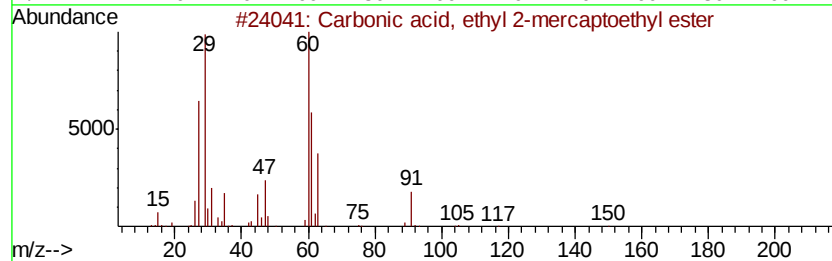
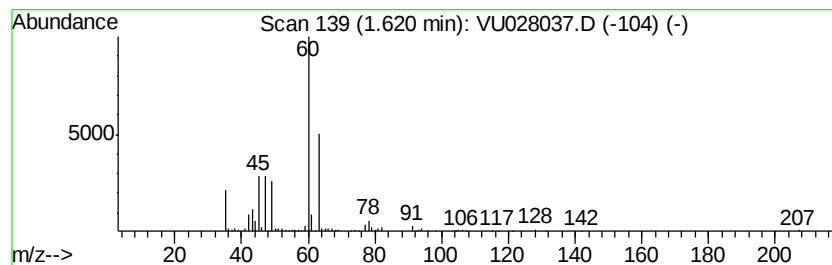
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown1.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.62	20.66 ug/l	222968	Pentafluorobenzene	4.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbonic acid, ethyl 2-mercaptoe...	150	C5H10O3S	005628-96-6	38
2		Chloro-methyl-methoxy-amine	95	C2H6ClNO	070569-68-5	9
3		2H-1,2-Oxazine, tetrahydro-2-met...	191	C12H17NO	015769-90-1	9
4		2-Amino-1,3-propanediol	91	C3H9NO2	000534-03-2	9
5		L-Galactose, 6-deoxy-	164	C6H12O5	002438-80-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.62	1.62	20.7	ug/l	222968	1	4.98	539569	50.0