

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN101220\
 Data File : VN064082.D
 Acq On : 12 Oct 2020 17:01
 Operator : JC/MD
 Sample : L4315-01
 Misc : 4.60g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 346

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_N\METHODS\82N093020W.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	2.463	197	225	226	rBV4	38885	115543	8.59%	1.583%
2	7.547	1785	1806	1817	rBV	171848	430656	32.03%	5.902%
3	7.621	1817	1829	1849	rVB2	278955	668870	49.74%	9.167%
4	7.984	1923	1942	1964	rBV2	197464	469173	34.89%	6.430%
5	8.547	2102	2117	2137	rBV	455744	958631	71.29%	13.138%
6	10.058	2573	2587	2617	rBV	733589	1344734	100.00%	18.429%
7	11.379	2985	2998	3023	rBV	727047	1268367	94.32%	17.382%
8	12.376	3294	3308	3330	rBV	558868	942682	70.10%	12.919%
9	13.318	3589	3601	3617	rBV	632877	1065000	79.20%	14.595%
10	15.087	4143	4151	4155	rBV8	9569	14227	1.06%	0.195%
11	15.997	4423	4434	4441	rBV8	11804	18927	1.41%	0.259%

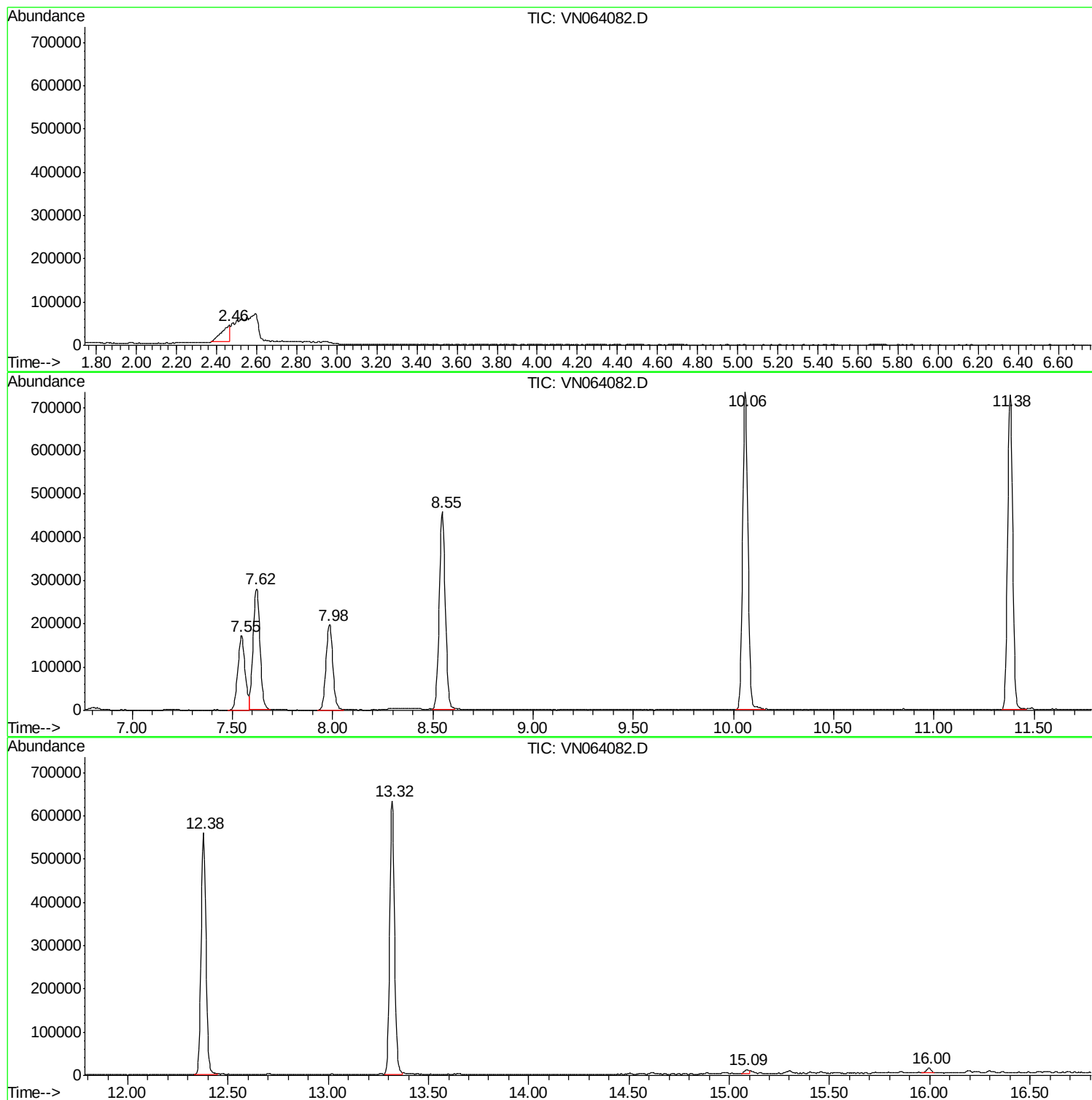
Sum of corrected areas: 7296810

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_N
ClientSampled :
346

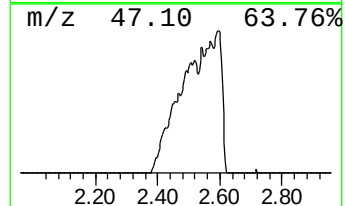
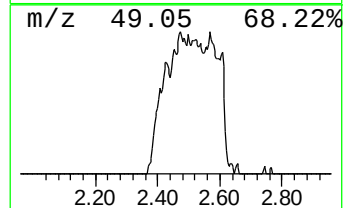
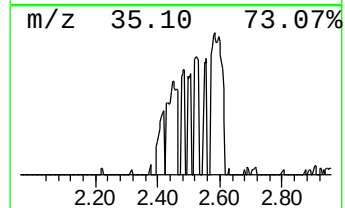
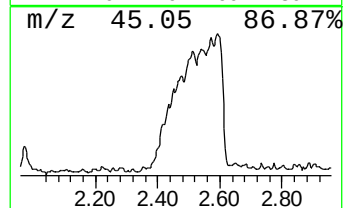
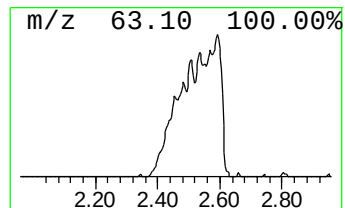
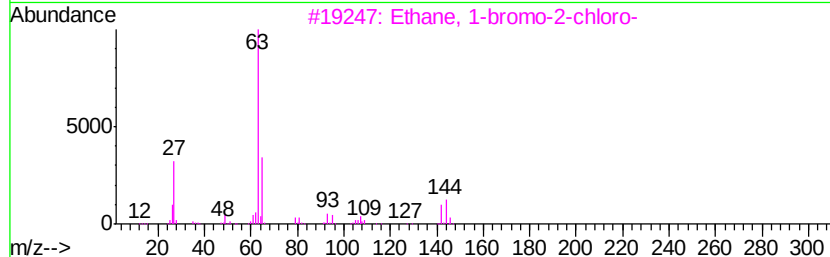
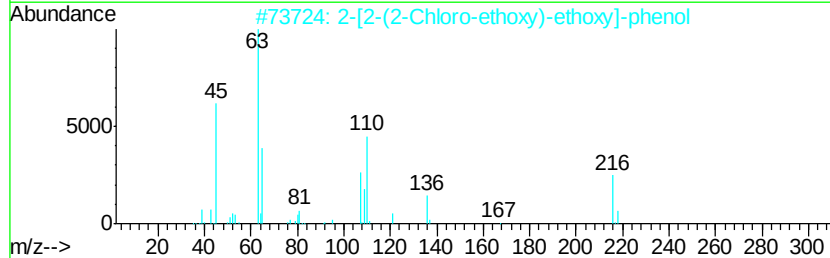
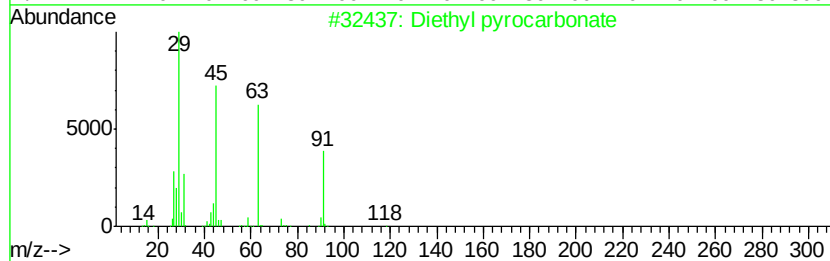
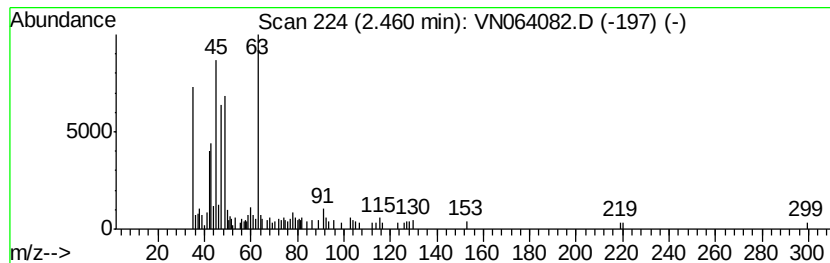
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
Quant Title : SW846 8260

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

Peak Number 1 unknown2.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.46	8.64 ug/l	115543	Pentafluorobenzene	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethyl pyrocarbonate	162	C6H10O5	001609-47-8	38
2		2-[2-(2-Chloro-ethoxy)-ethoxy]-p...	216	C10H13ClO3	1000317-46-1	23
3		Ethane, 1-bromo-2-chloro-	142	C2H4BrCl	000107-04-0	12
4		Carbonochloridic acid, 4-nitroph...	201	C7H4ClNO4	007693-46-1	10
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
unknown2.46	2.46	8.6	ug/l	115543	1	7.62	668870	50.0