

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071921\
 Data File : VN067789.D
 Acq On : 19 Jul 2021 13:26
 Operator : JC/MD
 Sample : M3032-01
 Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 NB-07-071421

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\82N070721W.M

Title : SW846 8260

Signal : TIC: VN067789.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.746	246	282	284	rBV3	39375	115700	5.52%	1.082%
2	5.066	1122	1147	1176	rBV4	18865	67550	3.22%	0.632%
3	8.016	2226	2247	2258	rBV2	238362	572953	27.32%	5.358%
4	8.080	2259	2271	2299	rVB2	440486	1021703	48.72%	9.555%
5	8.432	2381	2402	2427	rBV	287845	669858	31.94%	6.264%
6	8.660	2471	2487	2508	rVB6	11541	28182	1.34%	0.264%
7	8.965	2581	2601	2628	rBV	668826	1431205	68.25%	13.384%
8	10.440	3133	3151	3189	rBV	1092525	2097080	100.00%	19.611%
9	11.746	3618	3638	3670	rBV	1000368	1815581	86.58%	16.979%
10	12.736	3990	4007	4034	rVB	768953	1311251	62.53%	12.262%
11	13.677	4341	4358	4379	rBV	913704	1562198	74.49%	14.609%

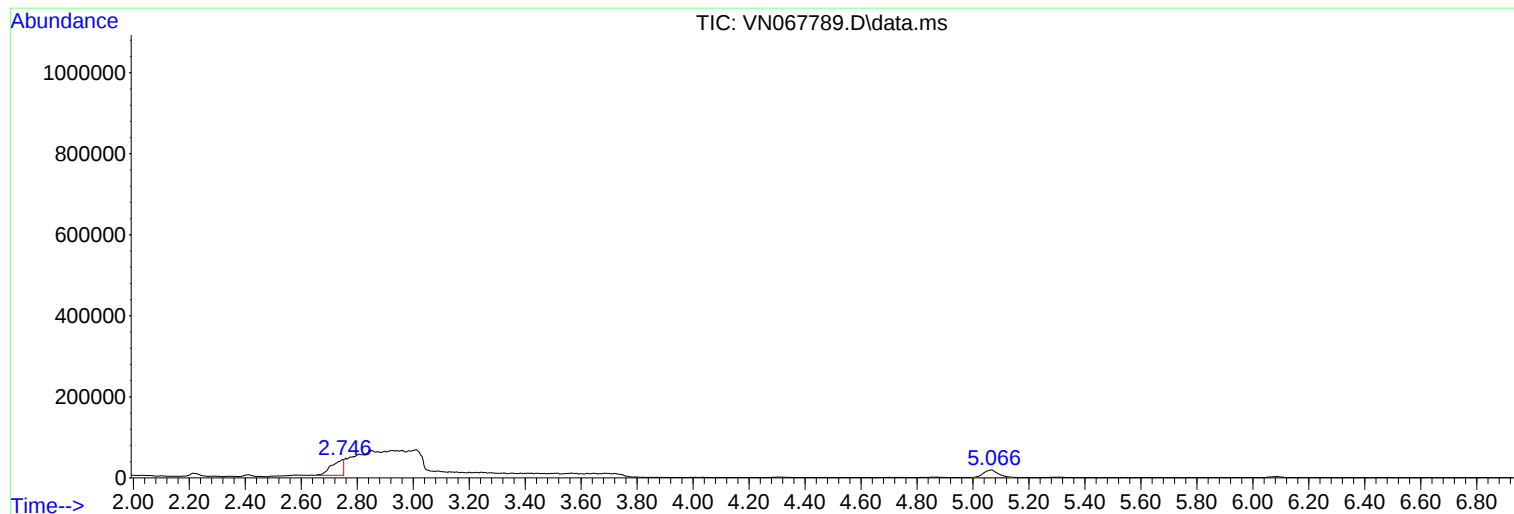
Sum of corrected areas: 10693261

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N070721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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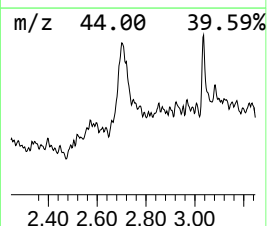
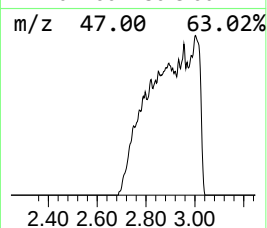
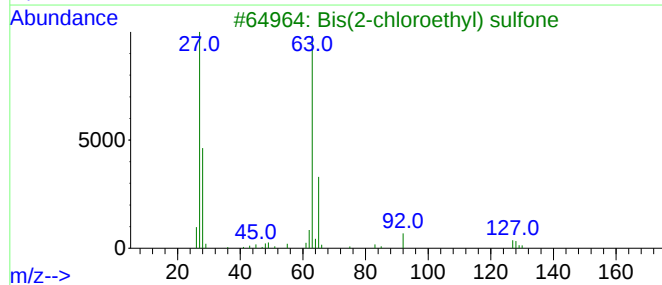
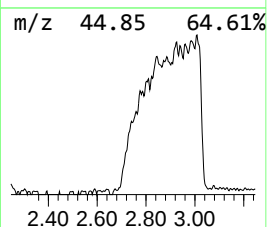
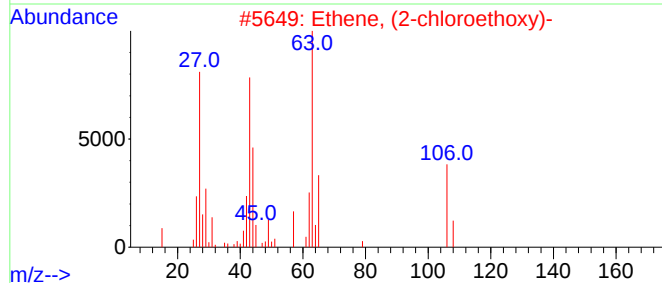
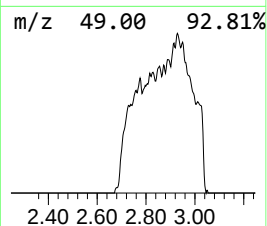
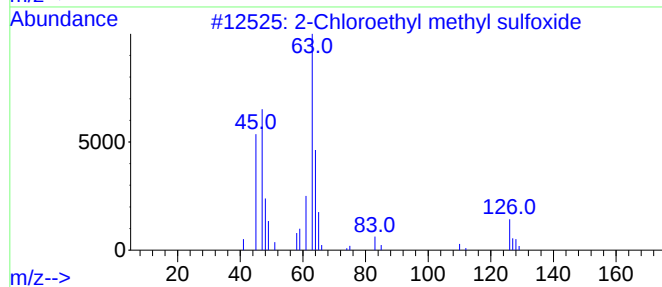
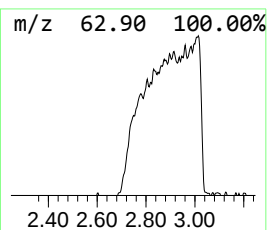
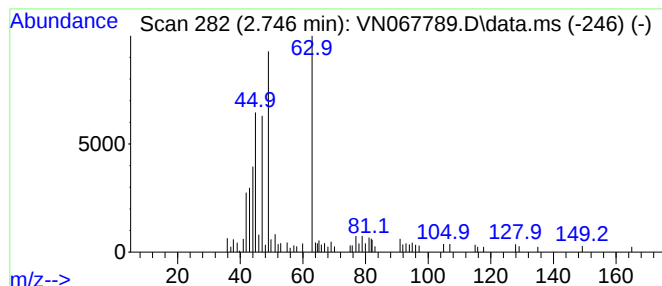
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N070721W.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown2.746 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.746	5.66 ug/l	115700	Pentafluorobenzene	8.078

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloroethyl methyl sulfoxide	126	C3H7ClOS	005331-57-7	23
2			Ethene, (2-chloroethoxy)-	106	C4H7ClO	000110-75-8	23
3			Bis(2-chloroethyl) sulfone	190	C4H8Cl2O2S	000471-03-4	12
4			Methane, bromochloro-	128	CH2BrCl	000074-97-5	9
5			Ethane, 1,1-dichloro-	98	C2H4Cl2	000075-34-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.746	2.746	5.7	ug/l	115700	1	8.078	1021700	50.0