

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091620\
 Data File : VX018448.D
 Acq On : 16 Sep 2020 18:21
 Operator : JC/SP
 Sample : L4021-02 20X
 Misc : 10.47g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B14-5-7

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X082120W.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	1.593	77	83	86	rVB8	16573	30013	1.20%	0.204%
2	3.001	305	314	320	rVB2	13397	36786	1.47%	0.250%
3	4.648	577	584	589	rBV2	14462	44237	1.76%	0.301%
4	5.464	707	718	730	rBV	313577	886672	35.36%	6.033%
5	5.629	734	745	761	rVB	531767	1489935	59.42%	10.138%
6	6.031	800	811	824	rBV	344719	908643	36.24%	6.183%
7	6.830	929	942	954	rBV	798818	1877293	74.87%	12.774%
8	8.696	1241	1248	1256	rBV	1648840	2507312	100.00%	17.061%
9	9.317	1343	1350	1358	rBV	130352	199822	7.97%	1.360%
10	9.604	1393	1397	1406	rBV	114610	244292	9.74%	1.662%
11	10.098	1472	1478	1488	rVB	1691930	2264870	90.33%	15.411%
12	11.122	1640	1646	1658	rBV	1475662	1822675	72.69%	12.402%
13	11.512	1707	1710	1717	rVB	17484	27182	1.08%	0.185%
14	12.061	1795	1800	1808	rBV	1928425	2356693	93.99%	16.036%

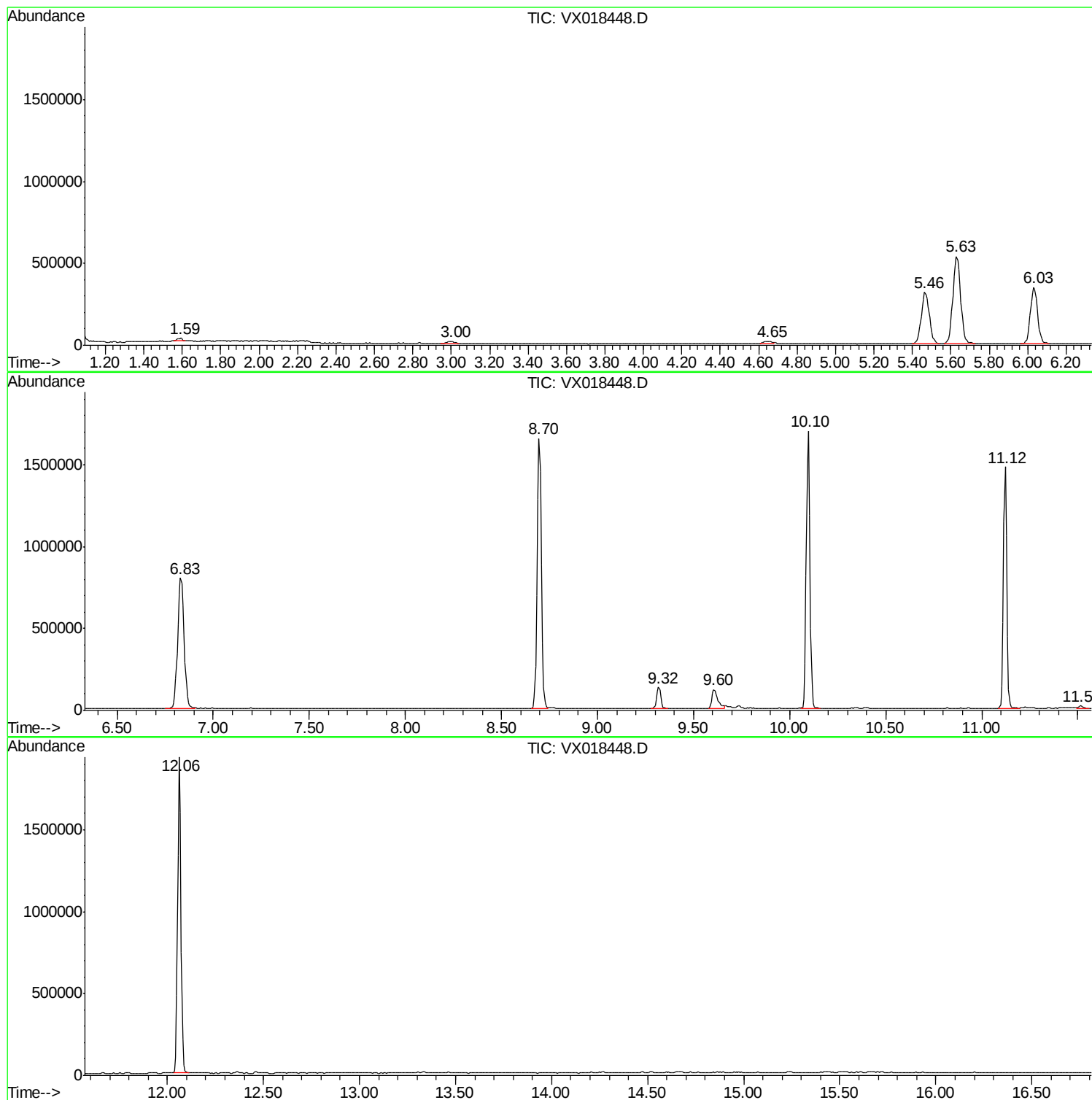
Sum of corrected areas: 14696425

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082120W.M
Quant Title : SW846 8260

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



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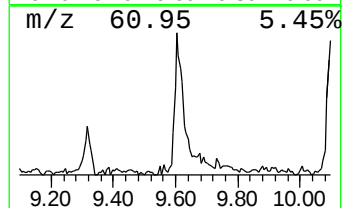
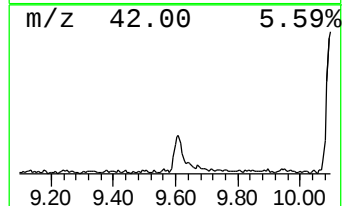
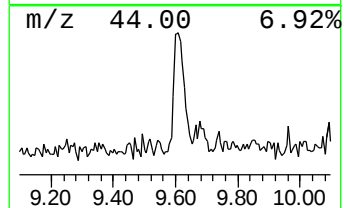
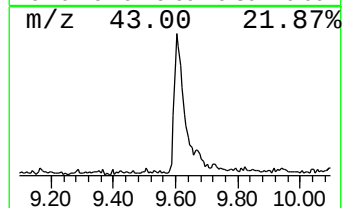
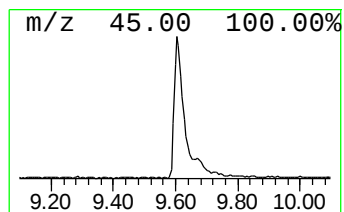
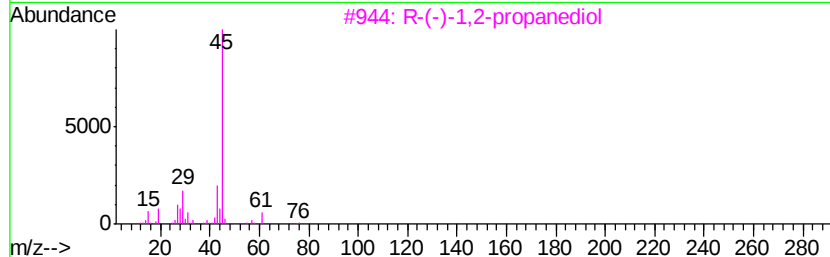
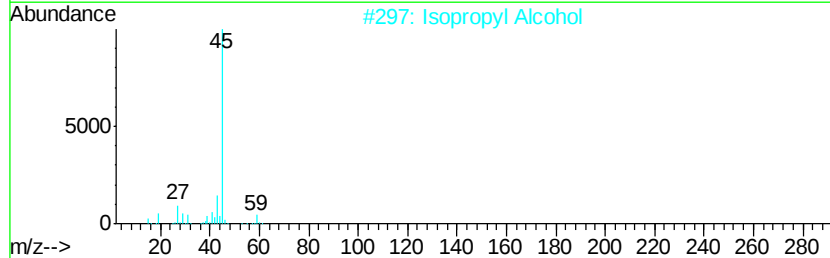
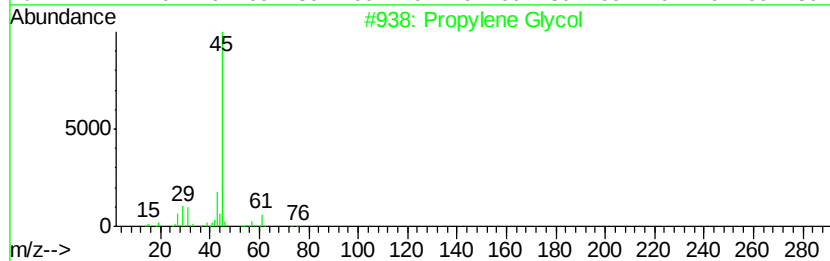
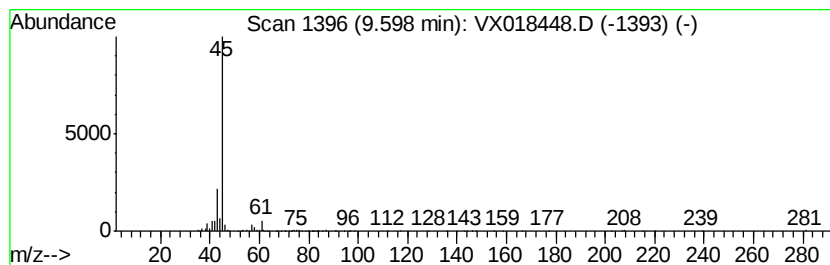
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 1 Propylene Glycol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.60	5.39 ug/l	244292	Chlorobenzene-d5	10.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propylene Glycol	76	C3H8O2	000057-55-6	86
2		Isopropyl Alcohol	60	C3H8O	000067-63-0	72
3		R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	56
4		1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	43
5		2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	39



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propylene Glycol	9.60	5.4	ug/l	244292	3	10.10	2264870	50.0