

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091020\
 Data File : VX018347.D
 Acq On : 10 Sep 2020 18:19
 Operator : JC/SP
 Sample : L3952-10 10X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 16718

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_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	2.154	157	175	178	rBV2	18956697	58635796	100.00%	68.360%
2	5.465	707	718	730	rBV3	290240	840855	1.43%	0.980%
3	5.635	736	746	758	rVB2	489561	1390563	2.37%	1.621%
4	6.032	797	811	817	rBV	325008	837362	1.43%	0.976%
5	6.117	817	825	848	rVB	1253153	3324269	5.67%	3.876%
6	6.836	931	943	953	rBV	726666	1761509	3.00%	2.054%
7	8.696	1241	1248	1254	rBV	1513318	2353845	4.01%	2.744%
8	8.769	1254	1260	1275	rVB	4349988	6802359	11.60%	7.930%
9	9.604	1392	1397	1405	rBV	444407	848530	1.45%	0.989%
10	10.098	1472	1478	1484	rBV	1605356	2147988	3.66%	2.504%
11	10.342	1510	1518	1527	rBV	1391621	1816424	3.10%	2.118%
12	10.683	1569	1574	1584	rVB	707992	919148	1.57%	1.072%
13	11.122	1640	1646	1657	rBV	1454527	1788879	3.05%	2.086%
14	12.061	1795	1800	1806	rBV	1920679	2307185	3.93%	2.690%

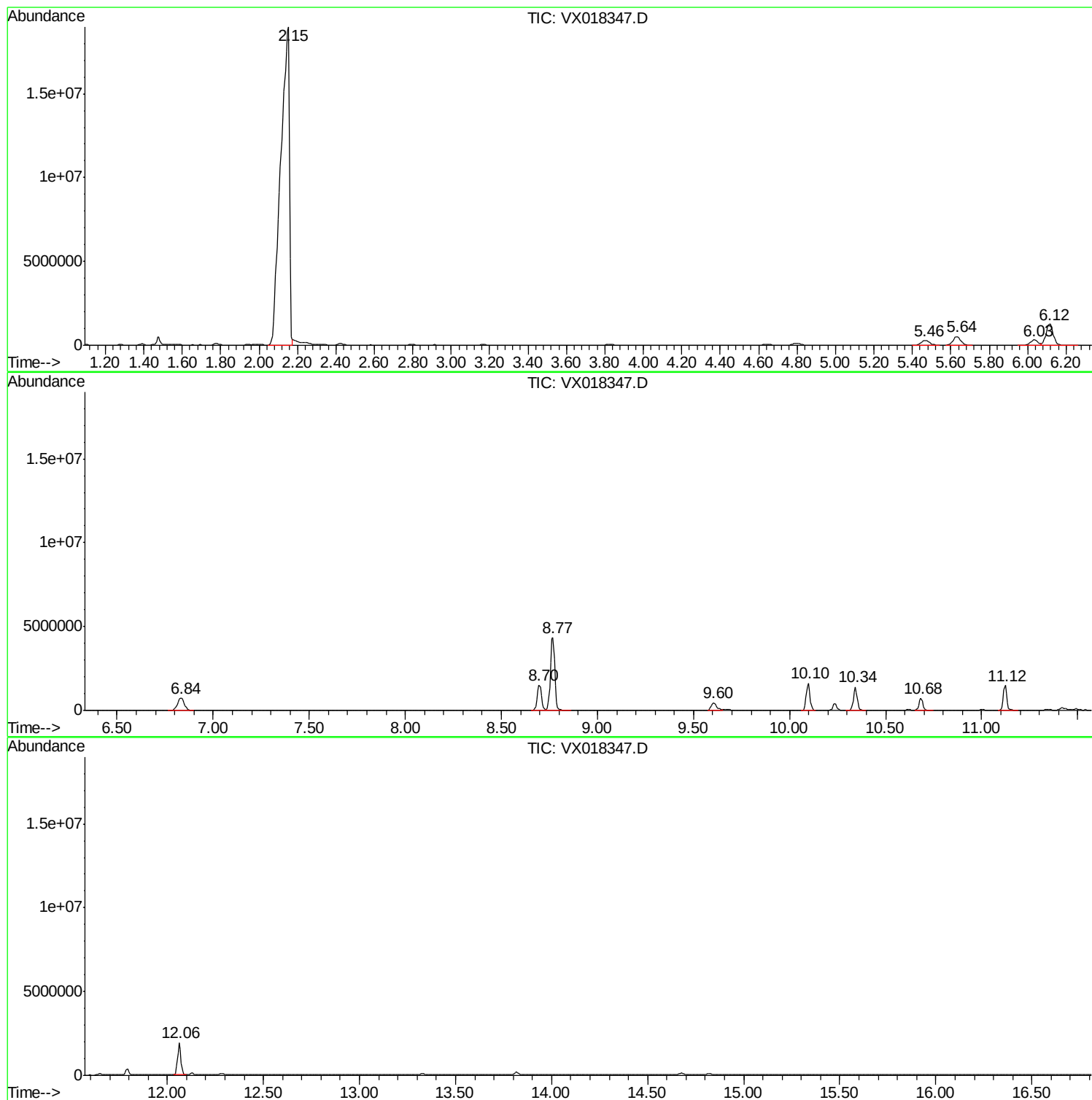
Sum of corrected areas: 85774712

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Quant Title : SW846 8260

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



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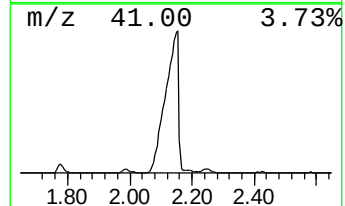
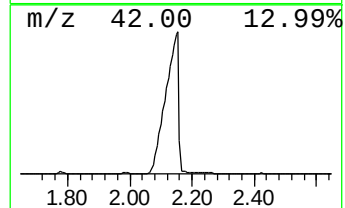
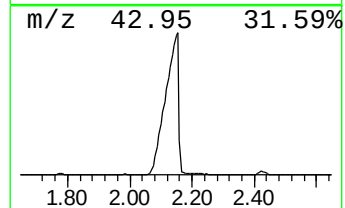
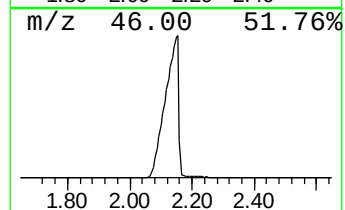
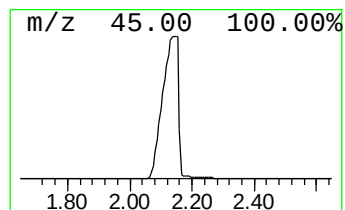
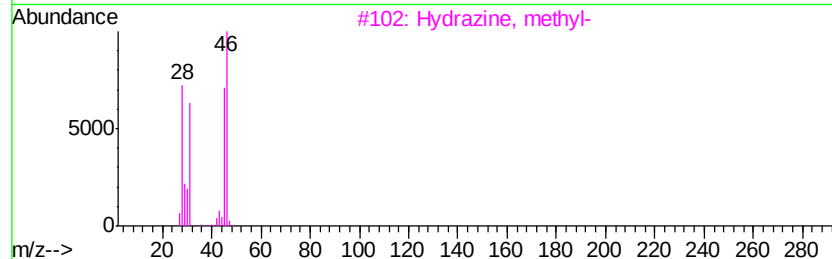
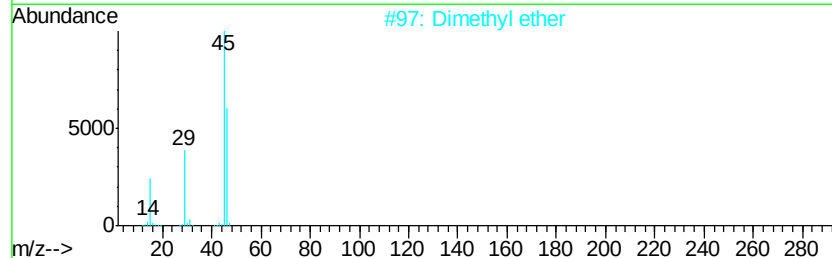
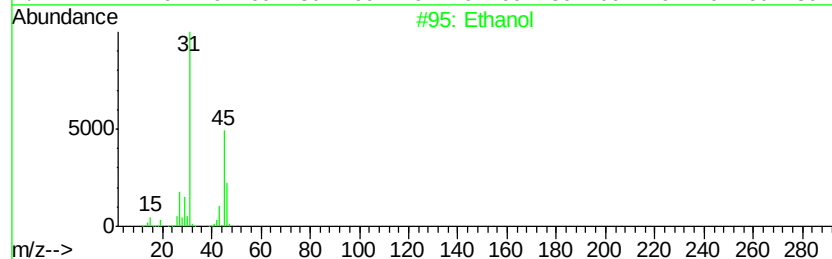
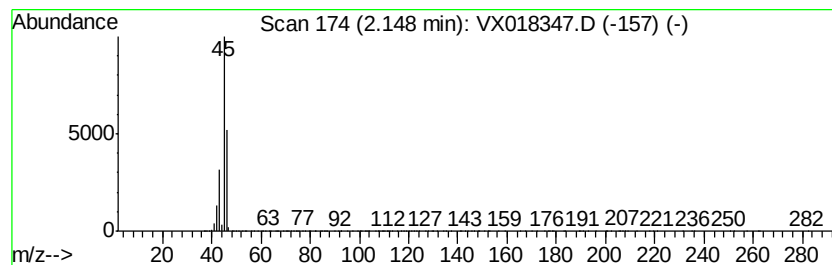
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Peak Number 1 Ethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
2.15	2108.35 ug/l	58635800	Pentafluorobenzene	5.64		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol		46	C2H6O	000064-17-5	83
2	Dimethyl ether		46	C2H6O	000115-10-6	9
3	Hvdrazine, methyl-		46	CH6N2	000060-34-4	9
4	Methane, nitroso-		45	CH3NO	000865-40-7	4
5	Ethene, fluoro-		46	C2H3F	000075-02-5	4



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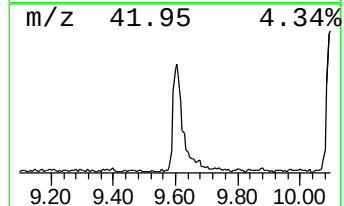
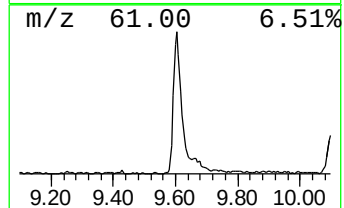
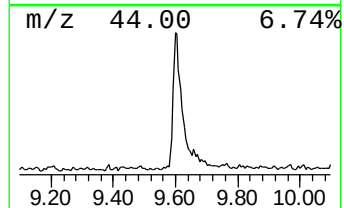
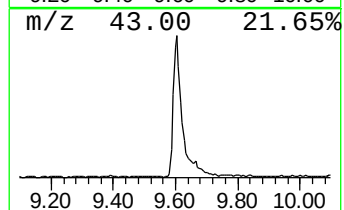
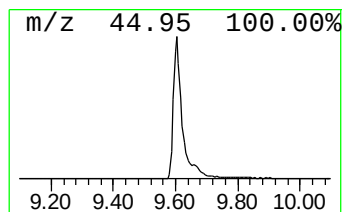
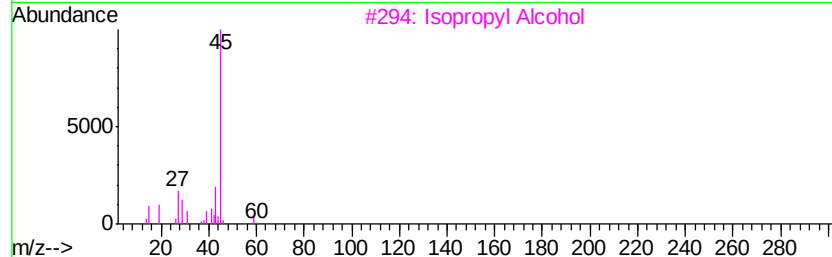
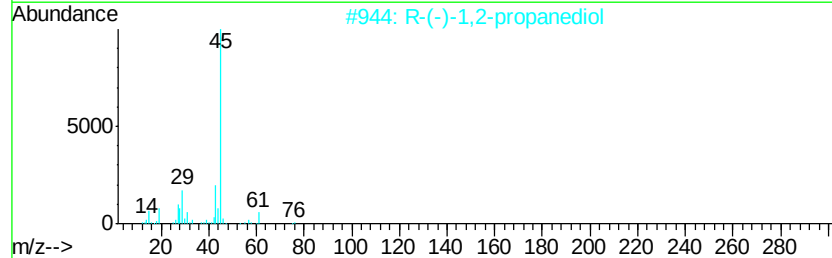
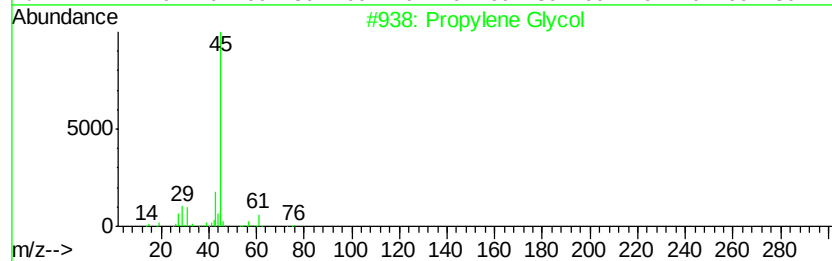
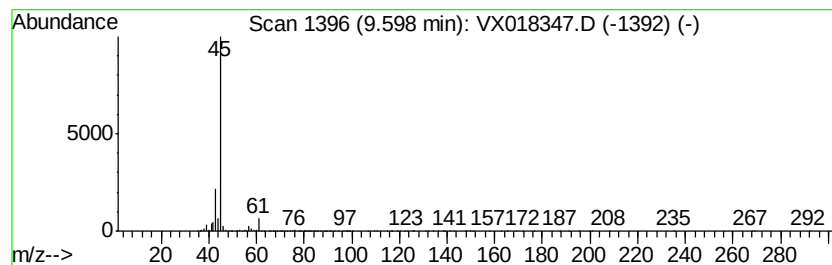
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Peak Number 2 Propylene Glycol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.60	19.75 ug/l	848530	Chlorobenzene-d5	10.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propylene Glycol	76	C3H8O2	000057-55-6	90
2		R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3		Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4		(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40
5		(R)-(-)-3-Methyl-2-butanol	88	C5H12O	001572-93-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol	2.15	2108.3	ug/l	58635800	1	5.64	1390560	50.0
Propylene Glycol	9.60	19.8	ug/l	848530	3	10.10	2147990	50.0