

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX040920\
 Data File : VX015602.D
 Acq On : 09 Apr 2020 17:30
 Operator : JC/SP
 Sample : L2190-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY062MW071-200406

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\82X032720W.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	1.064	23	29	43	rBV	6622342	8499856	24.88%	8.535%
2	4.563	592	603	617	rBV	427259	1150345	3.37%	1.155%
3	5.466	741	751	764	rVB	706139	2024252	5.92%	2.033%
4	5.630	764	778	791	rBV2	1228022	3441620	10.07%	3.456%
5	6.033	834	844	856	rBV	824347	2139467	6.26%	2.148%
6	6.831	963	975	988	rBV	1934600	4611601	13.50%	4.631%
7	7.191	1024	1034	1042	rBV	1822442	3882379	11.36%	3.898%
8	8.697	1274	1281	1291	rBV	4082426	6346577	18.58%	6.373%
9	9.319	1376	1383	1396	rBV	23397762	34165364	100.00%	34.306%
10	9.605	1423	1430	1457	rBV	3253734	11415609	33.41%	11.463%
11	10.099	1505	1511	1523	rVB	4120566	5803556	16.99%	5.827%
12	11.123	1673	1679	1688	rBV	3971519	5020848	14.70%	5.042%
13	12.062	1828	1833	1842	rBV	5425237	6447267	18.87%	6.474%
14	14.653	2246	2258	2270	rBV3	135743	504528	1.48%	0.507%
15	14.921	2282	2302	2337	rBV5	464369	3737938	10.94%	3.753%
16	15.299	2359	2364	2375	rVB	229386	397941	1.16%	0.400%

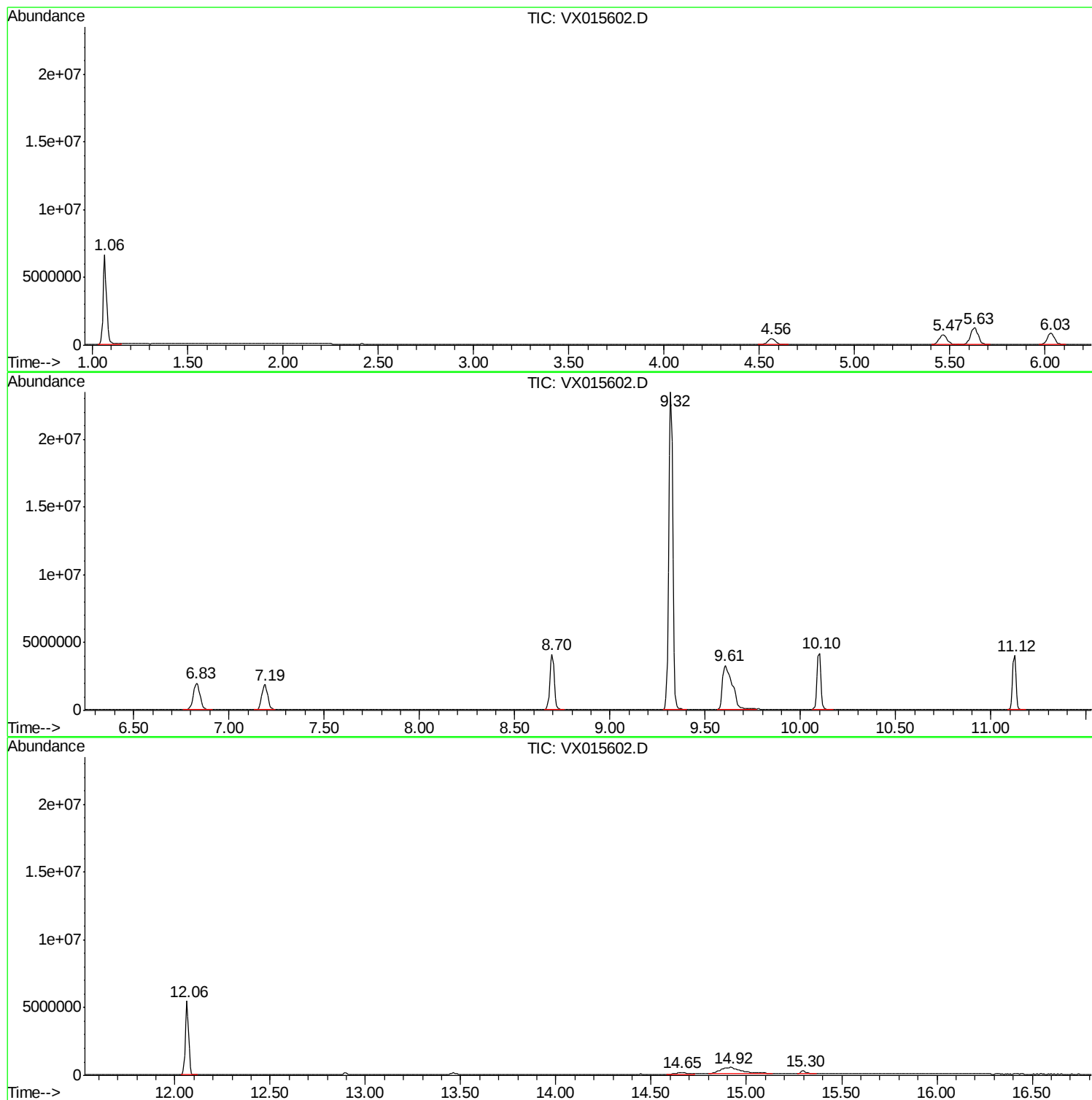
Sum of corrected areas: 99589148

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Quant Title : SW846 8260

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



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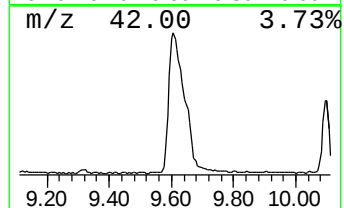
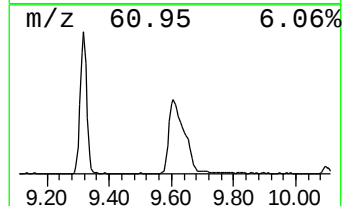
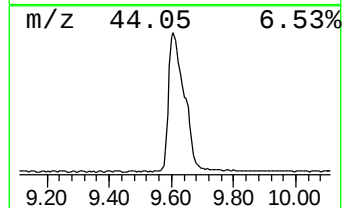
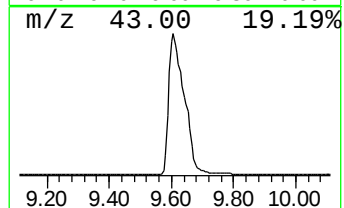
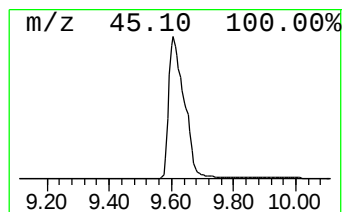
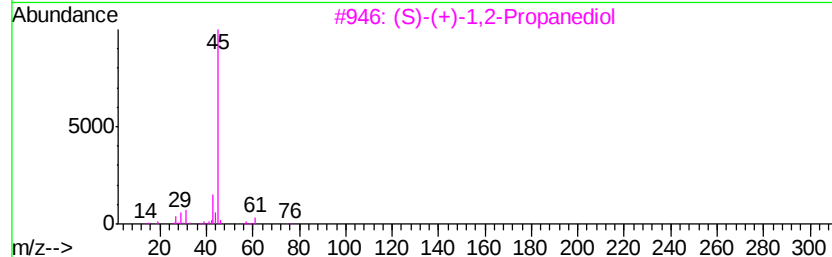
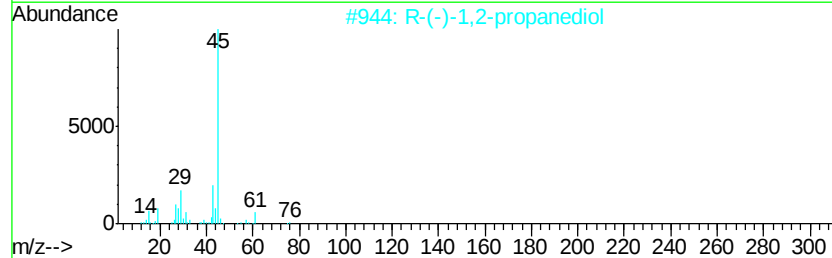
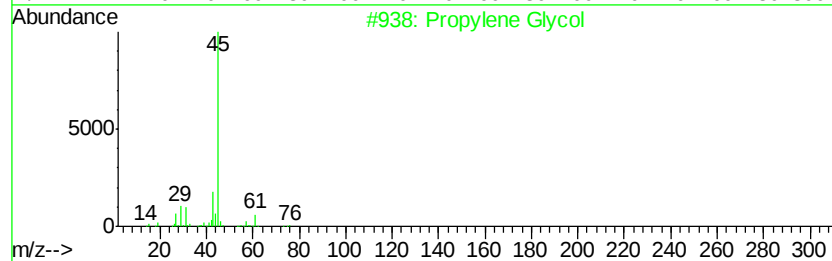
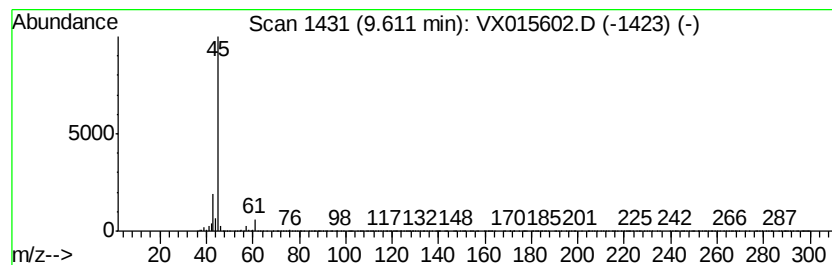
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Propylene Glycol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.61	98.35 ug/l	11415600	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		(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	64
4		Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5		2-Pentanol	88	C5H12O	006032-29-7	9



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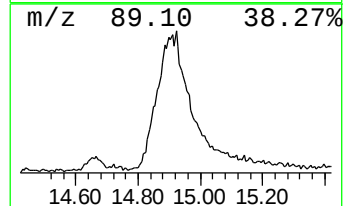
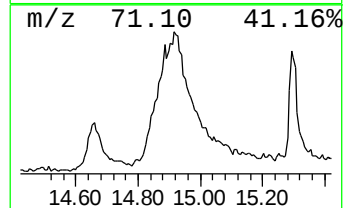
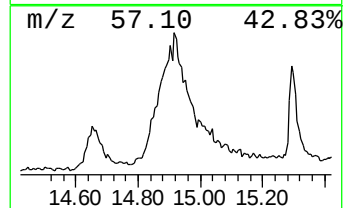
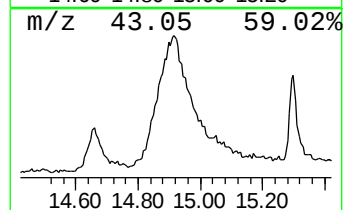
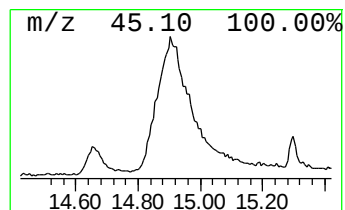
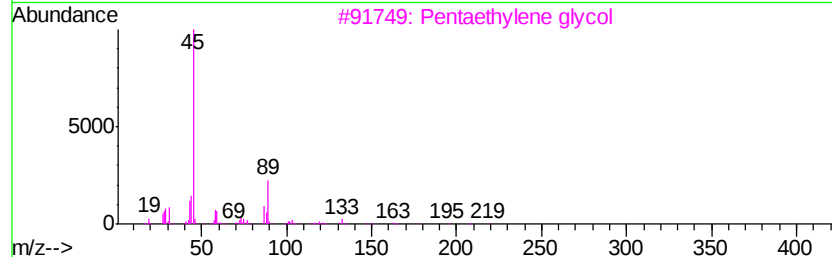
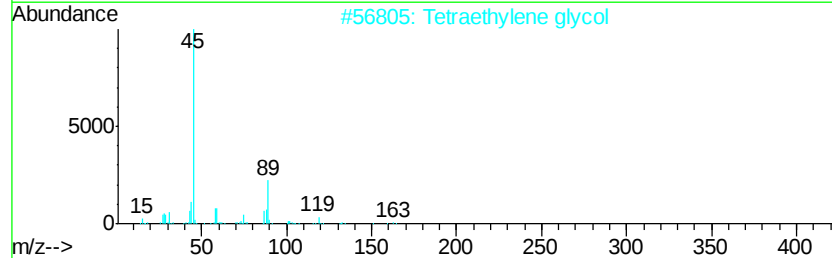
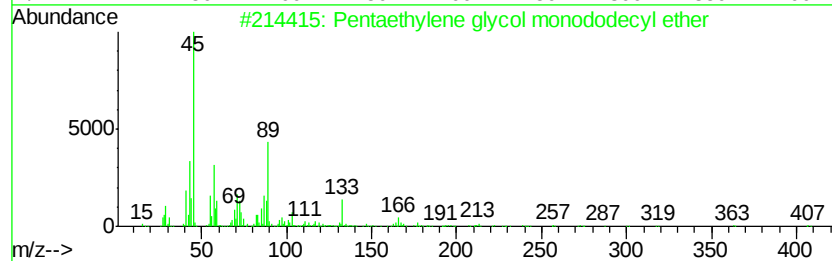
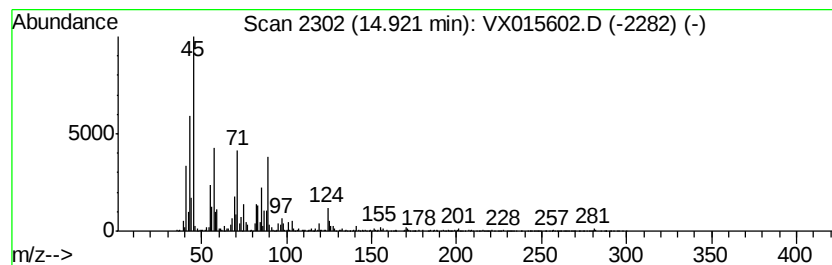
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Peak Number 3 Pentaethylene glycol monodo... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	28.99 ug/l	3737940	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pentaethylene glycol monododecyl...	406 C22H46O6	003055-95-6 53
2		Tetraethylene glycol	194 C8H18O5	000112-60-7 50
3		Pentaethylene glycol	238 C10H22O6	004792-15-8 50
4		Hexaethylene glycol	282 C12H26O7	002615-15-8 50
5		Ethanol, 2-[2-(2-methoxyethoxy)e...	164 C7H16O4	000112-35-6 47



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propylene Glycol	9.61	98.3	ug/l	11415600	3	10.10	5803560	50.0
Pentaethylene gly...	14.92	29.0	ug/l	3737940	4	12.06	6447270	50.0