

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091620\
 Data File : VX018447.D
 Acq On : 16 Sep 2020 17:59
 Operator : JC/SP
 Sample : L4021-01 20X
 Misc : 9.44g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B14-1-3

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	2.995	306	313	320	rVB2	17621	44631	1.75%	0.290%
2	4.654	578	585	598	rVB2	19973	61996	2.43%	0.403%
3	5.465	706	718	731	rBV2	315552	917441	35.98%	5.963%
4	5.635	732	746	759	rVV2	531379	1515904	59.44%	9.853%
5	6.031	801	811	827	rBV	348551	912361	35.78%	5.930%
6	6.830	931	942	955	rBV	808012	1889377	74.09%	12.281%
7	8.696	1242	1248	1257	rBV	1649113	2550119	100.00%	16.576%
8	9.317	1344	1350	1358	rVB	88253	130973	5.14%	0.851%
9	9.604	1392	1397	1407	rBV	147538	317947	12.47%	2.067%
10	10.098	1472	1478	1486	rBV	1740400	2324893	91.17%	15.112%
11	11.122	1641	1646	1655	rBV	1544806	1916838	75.17%	12.459%
12	11.518	1707	1711	1716	rVB3	28899	43018	1.69%	0.280%
13	11.787	1752	1755	1761	rBV3	15571	26505	1.04%	0.172%
14	12.061	1795	1800	1806	rBV	2072790	2473540	97.00%	16.078%
15	12.360	1844	1849	1855	rBV9	16820	37009	1.45%	0.241%
16	12.463	1861	1866	1869	rBV2	24644	30706	1.20%	0.200%
17	13.335	2006	2009	2014	rVB7	25335	44596	1.75%	0.290%
18	14.268	2158	2162	2168	rVB7	27751	46167	1.81%	0.300%
19	14.658	2222	2226	2228	rBV3	25971	35221	1.38%	0.229%
20	15.225	2316	2319	2322	rBV5	18878	25842	1.01%	0.168%
21	15.536	2363	2370	2374	rVB10	19802	39687	1.56%	0.258%

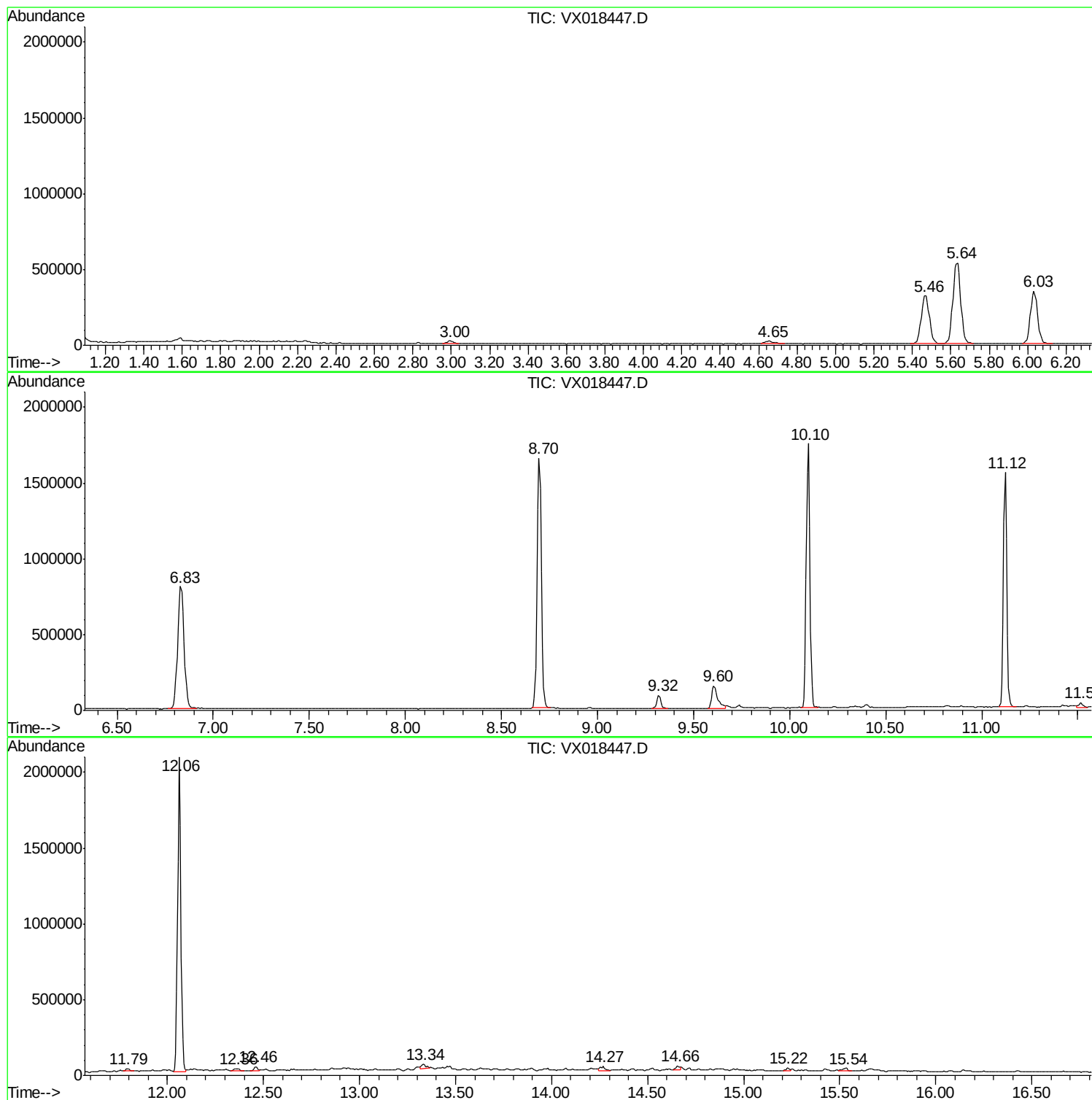
Sum of corrected areas: 15384771

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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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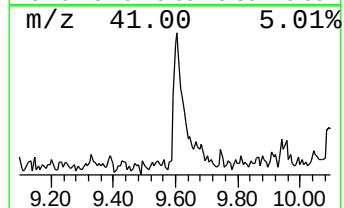
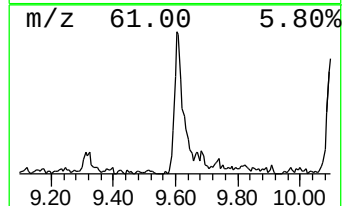
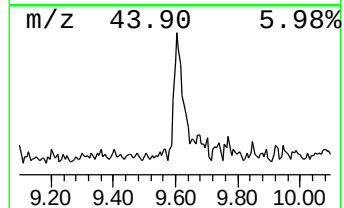
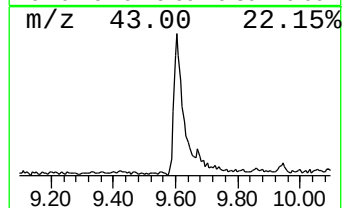
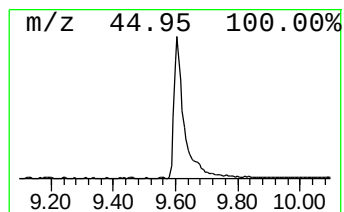
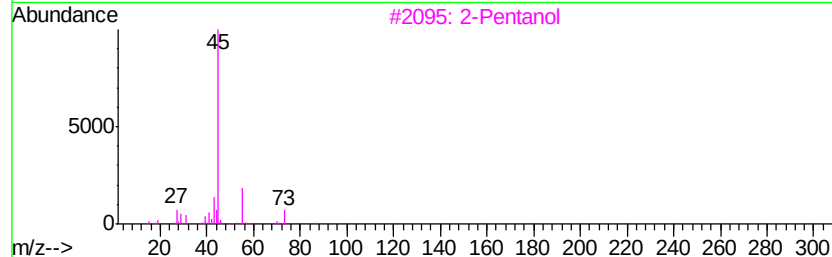
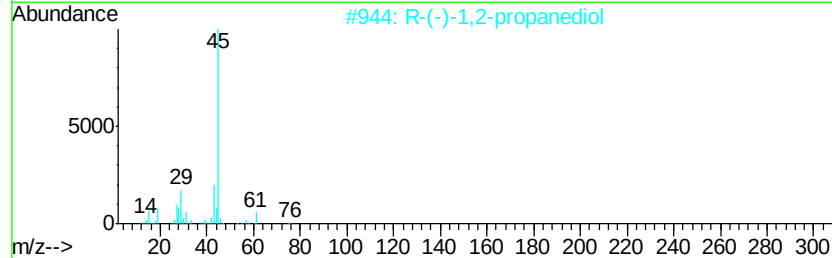
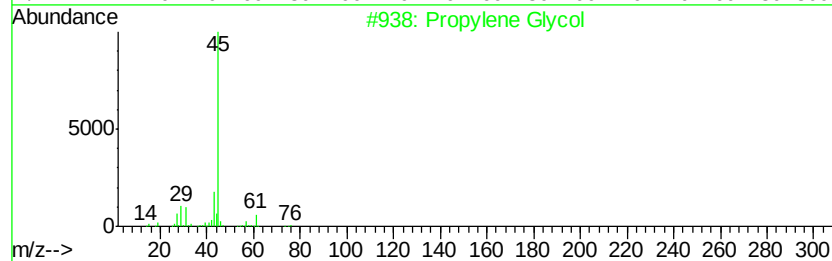
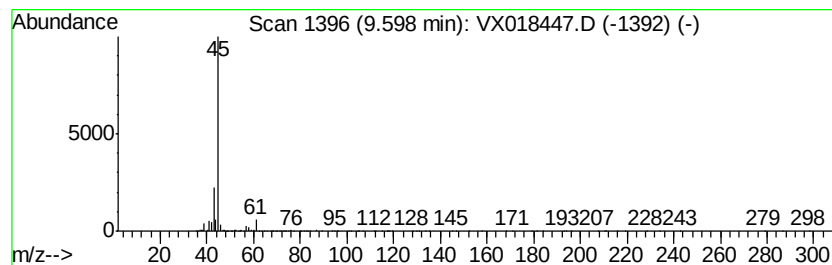
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Propylene Glycol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.60	6.84 ug/l	317947	Chlorobenzene-d5	10.10	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	86
2	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	40
3	2-Pentanol	88	C5H12O	006032-29-7	40
4	(R)-(-)-3-Methyl-2-butanol	88	C5H12O	001572-93-6	9
5	2-Heptanol, (S)-	116	C7H16O	006033-23-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propylene Glycol	9.60	6.8	ug/l	317947	3	10.10	2324890	50.0