

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY080320\
Data File : VY003425.D
Acq On : 03 Aug 2020 18:27
Operator : SY/MD
Sample : L3548-01
Misc : 8.73G/5ML/MSVOA Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
CRT-025

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_Y\METHODS\82Y073120S.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	2.524	122	129	138	rBV	10048	33983	1.50%	0.272%
2	4.724	480	490	501	rBV	51274	159704	7.05%	1.277%
3	7.273	896	908	921	rBV2	85760	231014	10.20%	1.847%
4	7.724	970	982	988	rBV	313597	762538	33.68%	6.098%
5	7.797	988	994	1006	rVB	555241	1241367	54.83%	9.927%
6	8.151	1042	1052	1062	rBV	284391	639799	28.26%	5.116%
7	8.693	1131	1141	1153	rBV	831196	1612661	71.23%	12.896%
8	10.181	1376	1385	1402	rBV	1308579	2264003	100.00%	18.105%
9	11.485	1590	1599	1611	rVB	1195553	1969053	86.97%	15.746%
10	12.479	1755	1762	1772	rVB	1010098	1529722	67.57%	12.233%
11	13.083	1851	1861	1876	rVB	22291	98637	4.36%	0.789%
12	13.424	1910	1917	1924	rBV	1286700	1962311	86.67%	15.692%

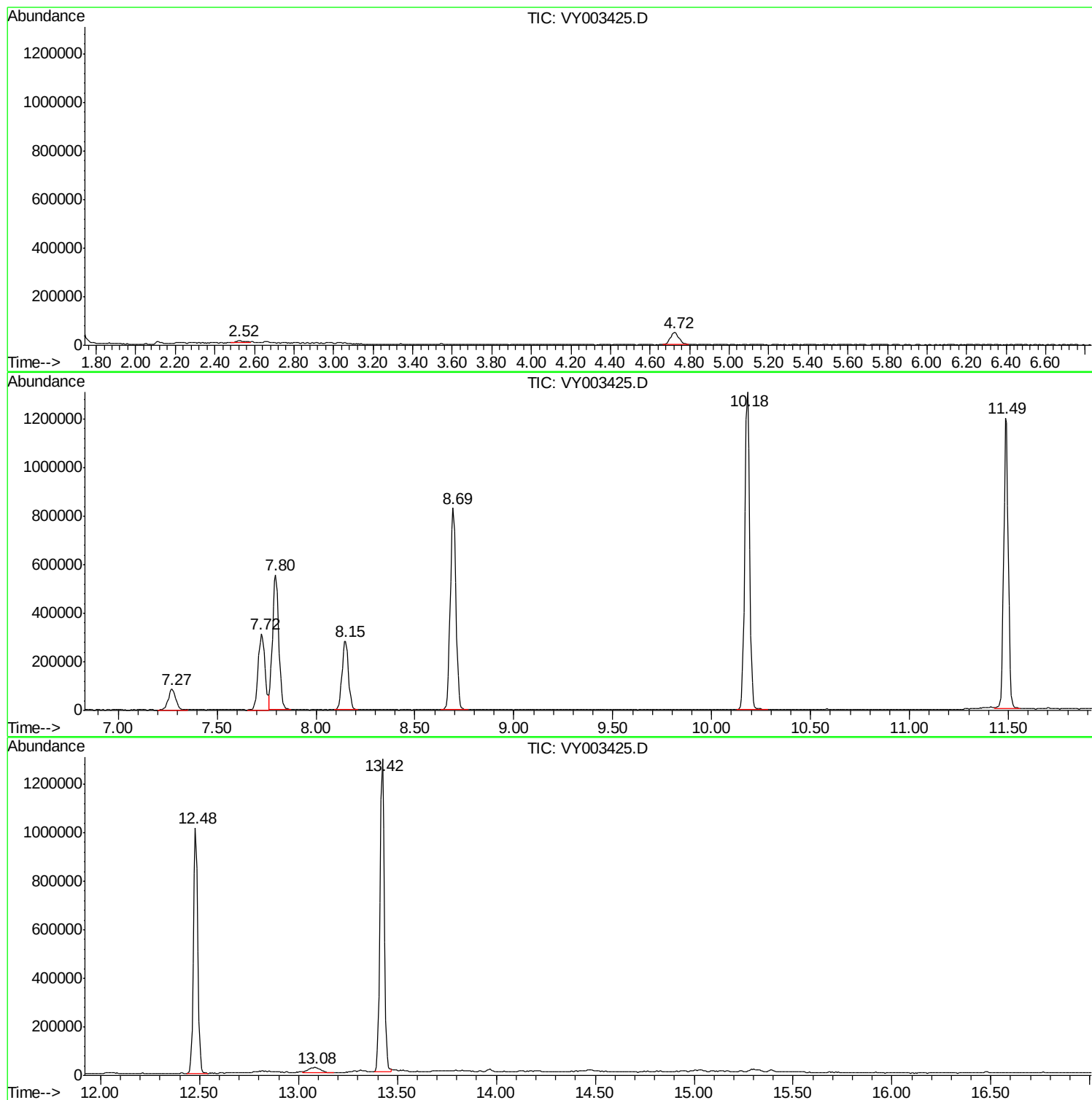
Sum of corrected areas: 12504792

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y073120S.M
Quant Title : SW846 8260

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



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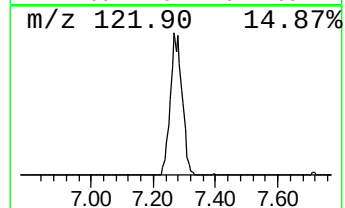
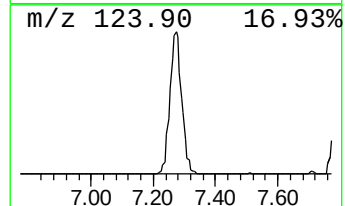
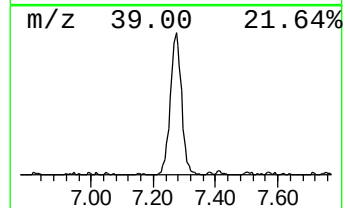
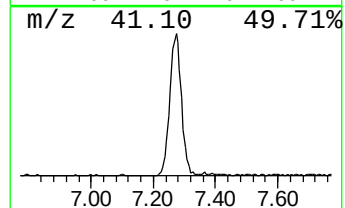
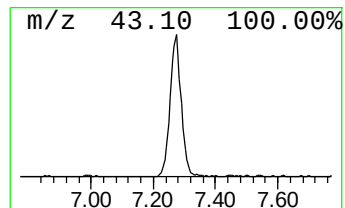
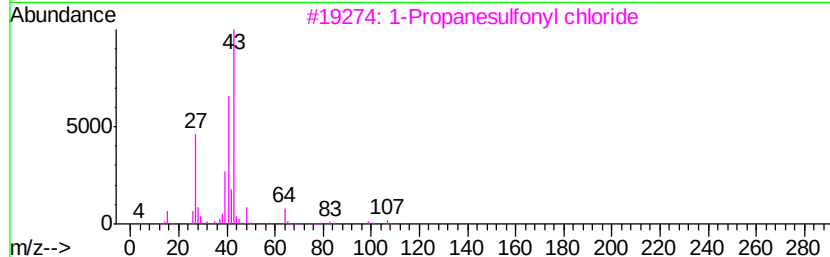
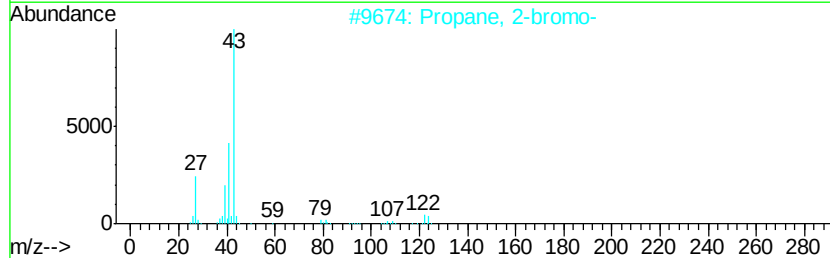
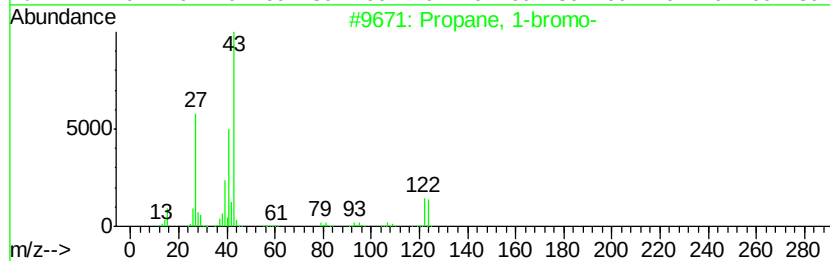
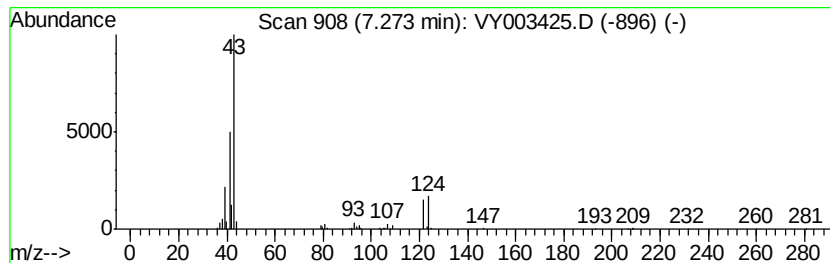
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y073120S.M
Quant Title : SW846 8260

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

Peak Number 2 Propane, 1-bromo- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.27	9.30 ug/l	231014	Pentafluorobenzene	7.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1-bromo-	122	C3H7Br	000106-94-5	83
2		Propane, 2-bromo-	122	C3H7Br	000075-26-3	47
3		1-Propanesulfonyl chloride	142	C3H7ClO2S	010147-36-1	23
4		Propane, 2-nitro-	89	C3H7NO2	000079-46-9	17
5		Propane, 1-nitro-	89	C3H7NO2	000108-03-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1-bromo-	7.27	9.3	ug/l	231014	1	7.80	1241370	50.0