

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092421\
 Data File : VY006148.D
 Acq On : 24 Sep 2021 19:01
 Operator : SY/MD
 Sample : M3927-21
 Misc : 6.06G/5ML/MSVOA_Y/SOIL
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 WOOSPK-BACKFILL-05

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\82Y091421S.M

Title : SW846 8260

Signal : TIC: VY006148.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.662	133	138	146	rVB3	4932	10506	1.24%	0.262%
2	3.960	340	351	360	rBV3	4007	10774	1.27%	0.268%
3	4.729	466	477	498	rBV3	24524	87084	10.25%	2.170%
4	5.186	537	552	567	rBV	75217	238140	28.02%	5.934%
5	7.722	958	968	974	rBV	80482	192847	22.69%	4.806%
6	7.795	974	980	995	rVB	141679	316631	37.26%	7.890%
7	8.149	1028	1038	1051	rBV	91436	197929	23.29%	4.932%
8	8.697	1119	1128	1139	rBV	214566	423488	49.84%	10.553%
9	10.185	1364	1372	1384	rBV	348477	591664	69.63%	14.744%
10	10.581	1428	1437	1444	rBV	27651	48239	5.68%	1.202%
11	11.490	1579	1586	1598	rVB	320118	513895	60.48%	12.806%
12	12.489	1742	1750	1759	rBV2	475183	849753	100.00%	21.176%
13	13.428	1894	1904	1912	rBV	302143	473157	55.68%	11.791%
14	13.965	1985	1992	2001	rVB	36150	58753	6.91%	1.464%

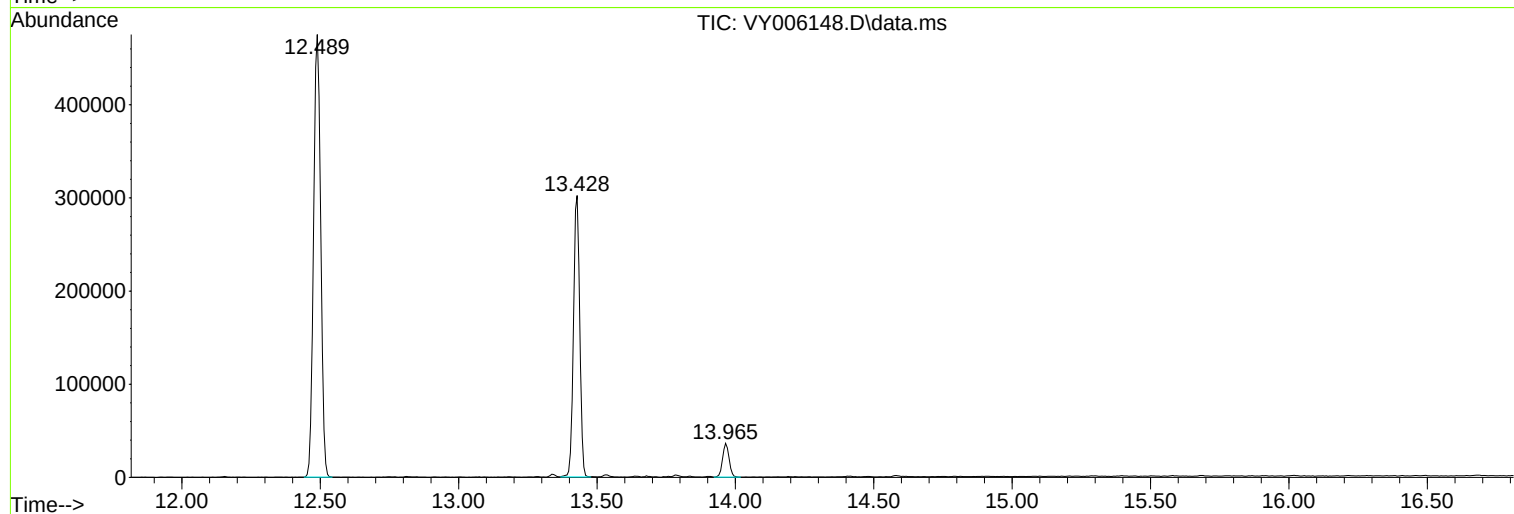
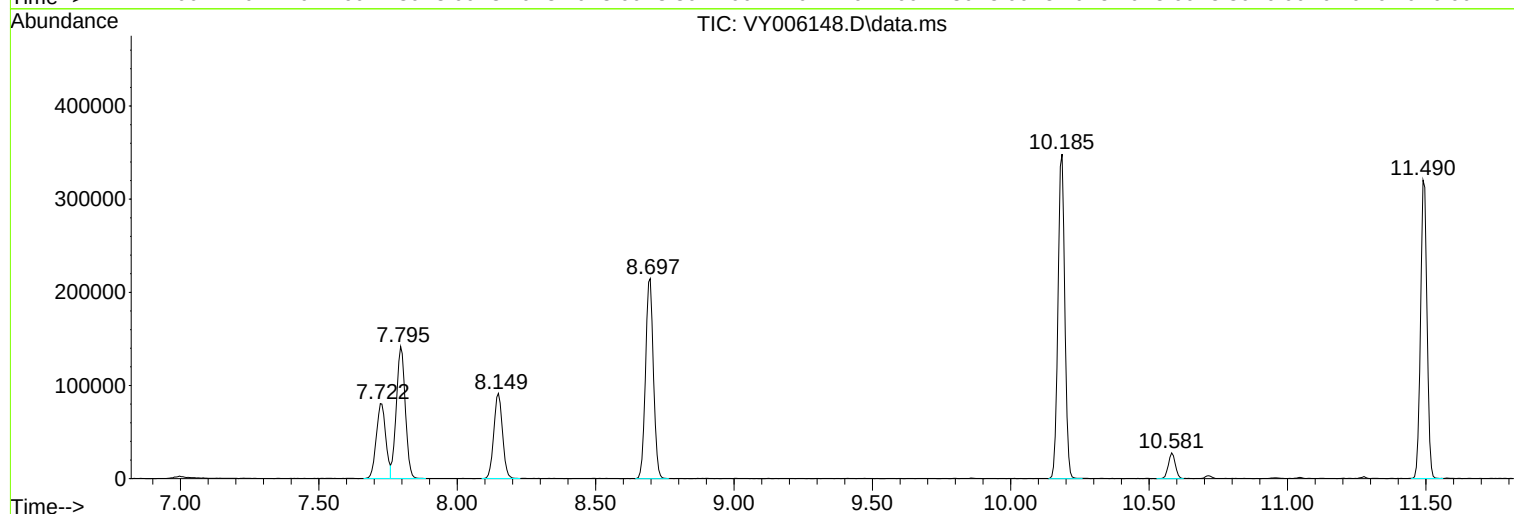
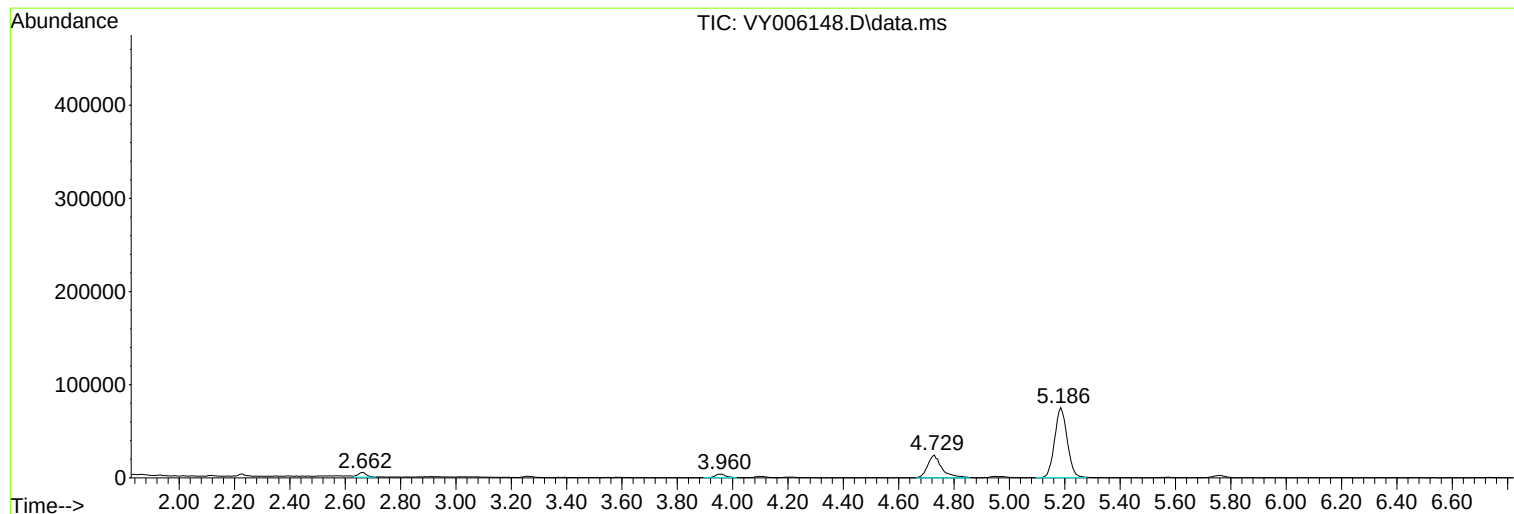
Sum of corrected areas: 4012860

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091421S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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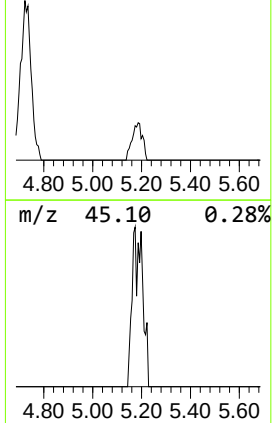
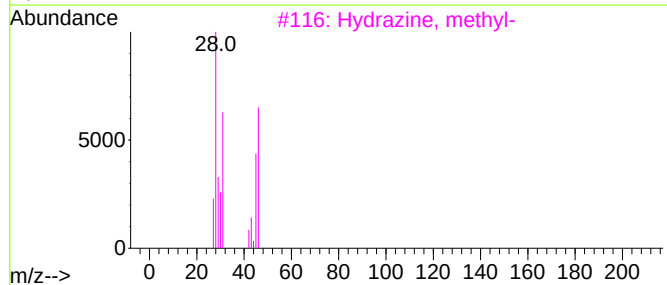
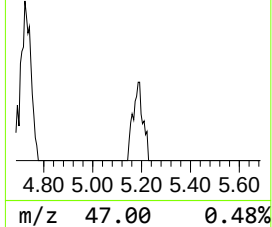
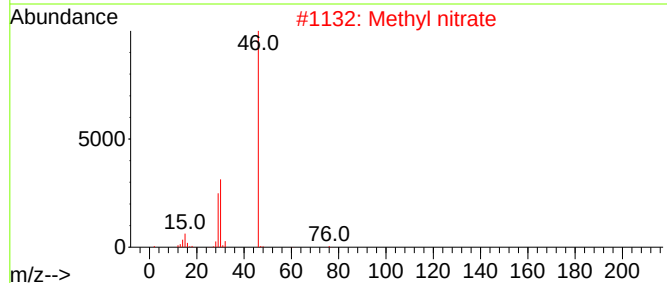
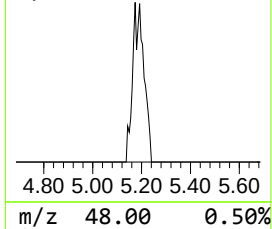
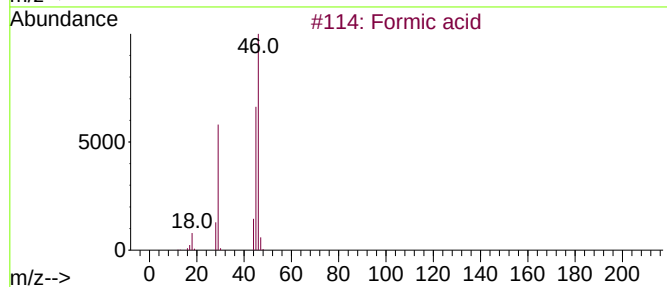
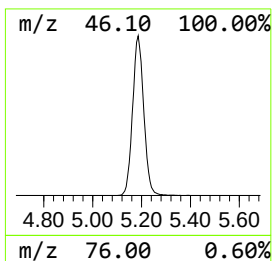
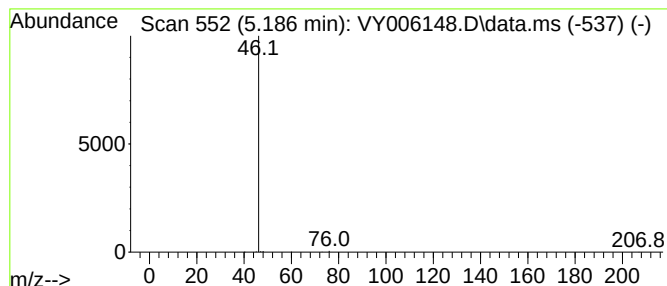
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Peak Number 1 unknown5.186 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.186	37.61 ug/l	238140	Pentafluorobenzene	7.795

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Formic acid	46	CH2O2	000064-18-6	4
2			Methyl nitrate	77	CH3NO3	000598-58-3	4
3			Hydrazine, methyl-	46	CH6N2	000060-34-4	3
4			Ethene, fluoro-	46	C2H3F	000075-02-5	3
5			Nitric acid, ethyl ester	91	C2H5NO3	000625-58-1	2



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown5.186	5.186	37.6	ug/l	238140	1	7.795	316631	50.0