

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061021\
 Data File : VX022654.D
 Acq On : 10 Jun 2021 19:39
 Operator : JC/MD
 Sample : M2644-08 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1824-25

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\82X061021W.M

Title : SW846 8260

Signal : TIC: VX022654.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.391	697	706	719	rBV2	41022	120745	1.24%	0.968%
2	5.562	722	734	750	rBV	71208	198756	2.04%	1.594%
3	5.964	790	800	815	rVB2	54042	146253	1.50%	1.173%
4	6.769	922	932	947	rBV	122514	294631	3.03%	2.363%
5	8.653	1234	1241	1247	rBV	269543	413077	4.24%	3.313%
6	8.720	1247	1252	1264	rVB	329354	519744	5.34%	4.168%
7	10.055	1466	1471	1487	rVB	258491	353024	3.63%	2.831%
8	11.085	1634	1640	1645	rBV	195379	250045	2.57%	2.005%
9	12.024	1788	1794	1801	rVB	257410	313352	3.22%	2.513%
10	12.225	1821	1827	1858	rBV	7655643	9737540	100.00%	78.092%
11	13.518	2033	2039	2047	rBV	81414	122198	1.25%	0.980%

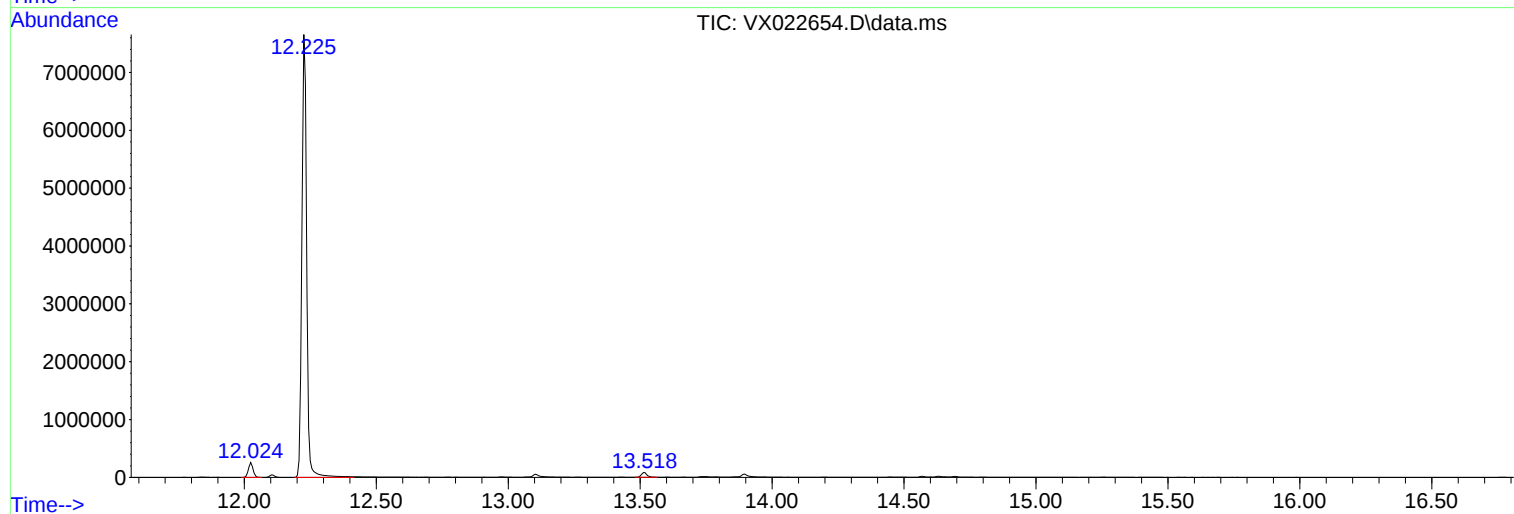
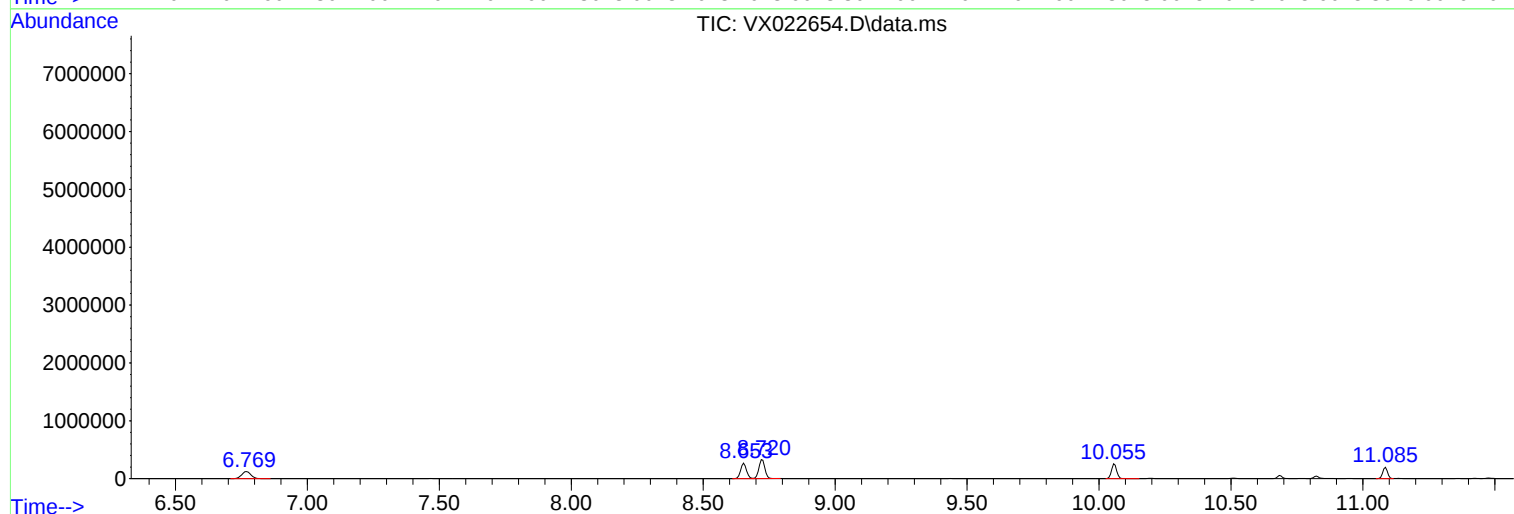
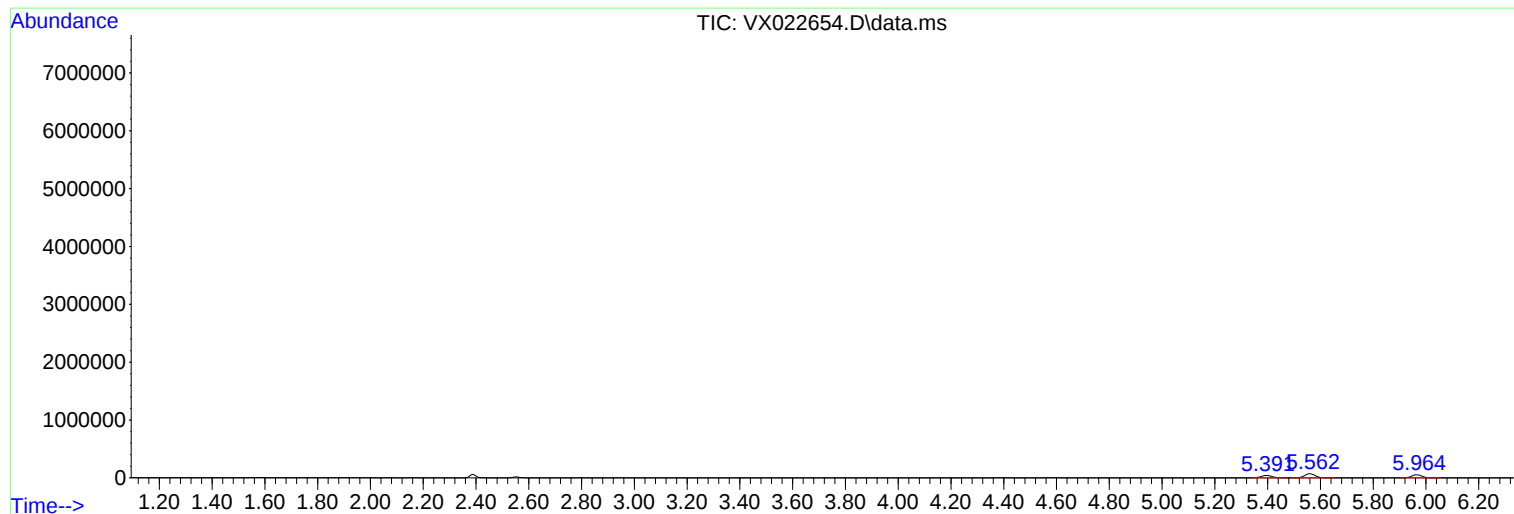
Sum of corrected areas: 12469365

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061021W.M
Quant Title : SW846 8260

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



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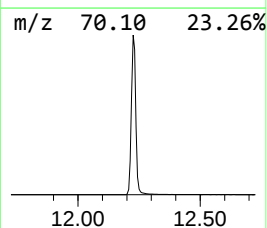
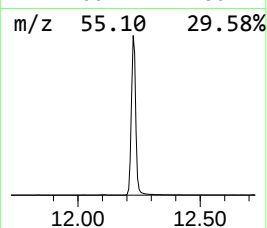
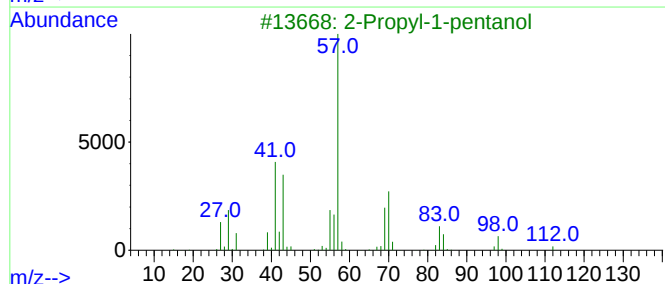
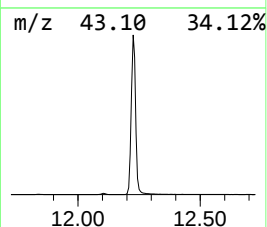
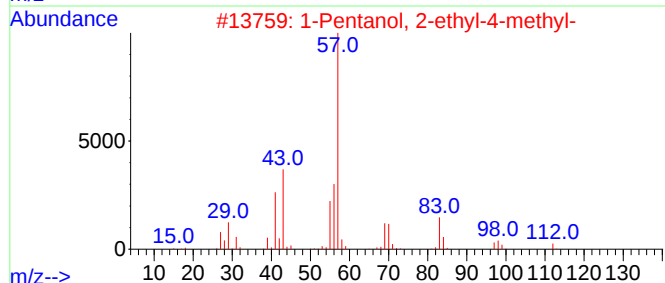
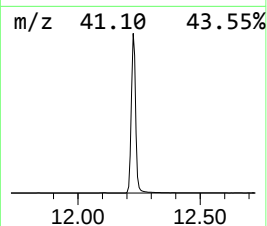
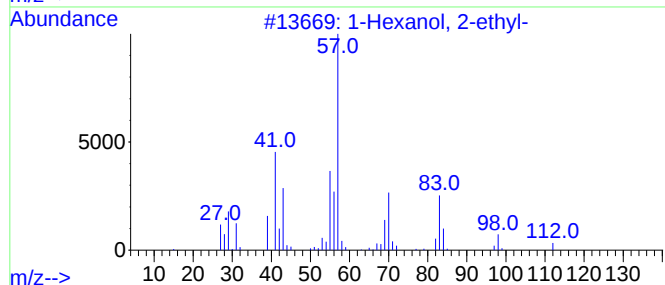
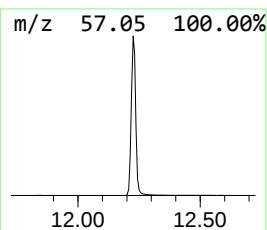
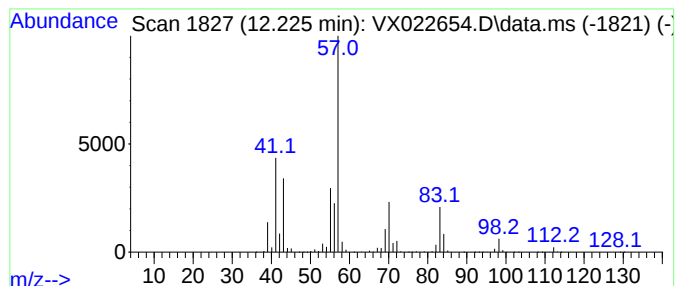
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061021W.M
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TIC Integration Parameters: LSCINT.P

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.225	1553.77 ug/l	9737540	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2		1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	56
3		2-Propyl-1-pentanol	130	C8H18O	058175-57-8	47
4		2-Ethyl-1-hexanol, methyl ether	144	C9H20O	1000365-10-2	47
5		1-Decene, 2,4-dimethyl-	168	C12H24	055170-80-4	42



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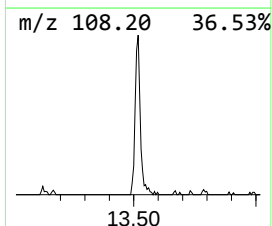
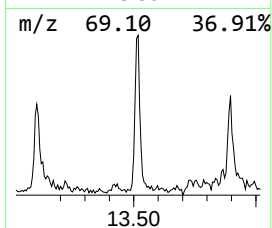
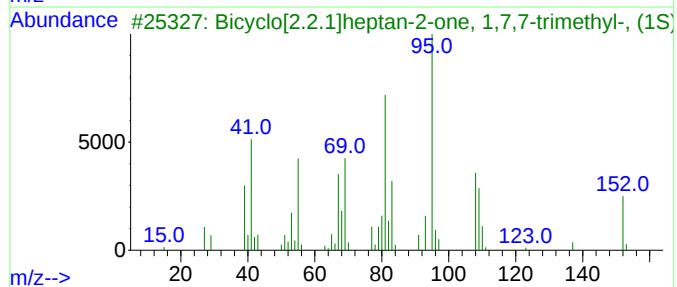
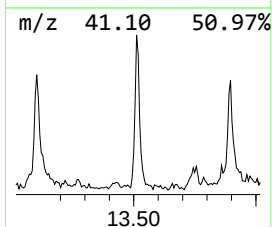
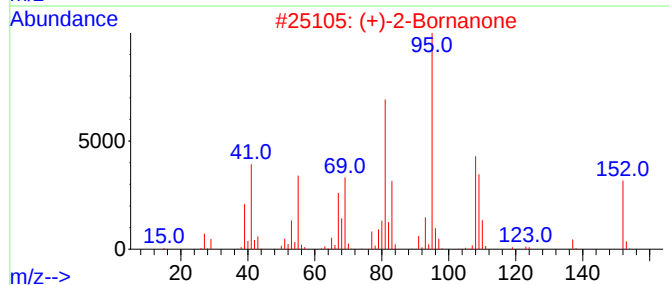
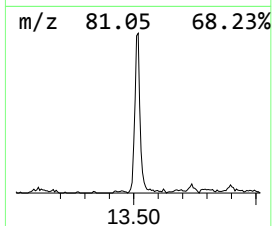
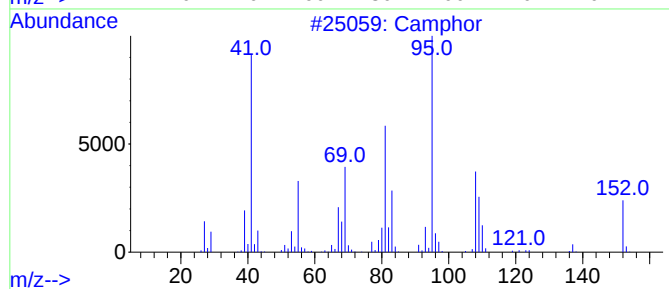
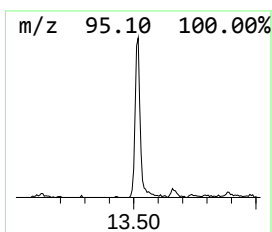
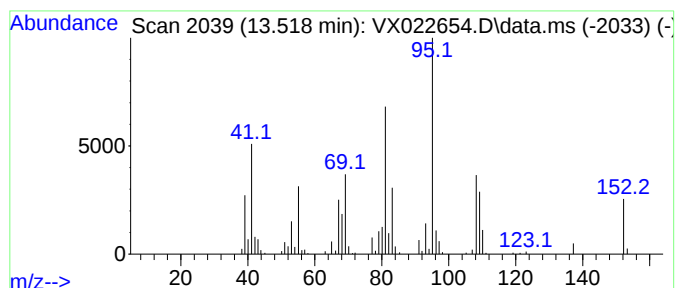
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Peak Number 2 Camphor Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.518	19.50 ug/l	122198	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Camphor	152	C10H16O	000076-22-2	98
2			(+)-2-Bornanone	152	C10H16O	000464-49-3	98
3			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	98
4			1-Adamantanol	152	C10H16O	000768-95-6	43
5			6-Hepten-1-ol, 2-methyl-	128	C8H16O	1000132-12-0	43



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Hexanol, 2-et...	12.225	1553.8	ug/l	9737540	4	12.024	313352	50.0
Camphor	13.518	19.5	ug/l	122198	4	12.024	313352	50.0