

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021524\
Data File : VX040257.D
Acq On : 15 Feb 2024 17:08
Operator : JC/MD
Sample : P1491-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
20241095-35N-N-FIELD-BLANK-1

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\82X020924W.M

Title : SW846 8260

Signal : TIC: VX040257.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.380	205	212	217	rBV2	2732	4662	1.18%	0.198%
2	3.447	377	387	402	rBV	50715	113171	28.70%	4.811%
3	4.300	518	527	536	rBV2	2529	6768	1.72%	0.288%
4	5.385	695	705	718	rBV	46673	134543	34.11%	5.720%
5	5.550	723	732	745	rVB	75873	211899	53.73%	9.008%
6	5.952	788	798	814	rBV	48807	131498	33.34%	5.590%
7	6.763	921	931	945	rBV	119403	285454	72.38%	12.135%
8	8.647	1233	1240	1257	rBV	252096	394388	100.00%	16.767%
9	10.055	1465	1471	1487	rVB	263789	369003	93.56%	15.687%
10	11.079	1634	1639	1654	rVB	228281	293018	74.30%	12.457%
11	12.018	1788	1793	1803	rBV	299512	390672	99.06%	16.609%
12	14.737	2235	2239	2246	rVB	14006	17157	4.35%	0.729%

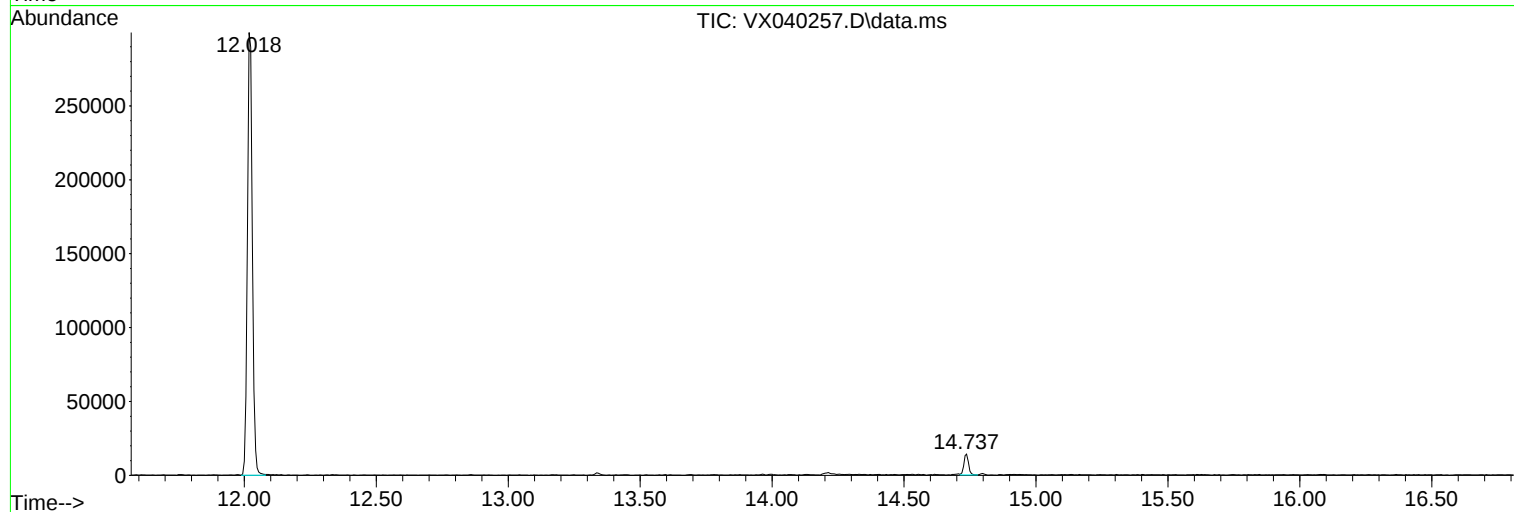
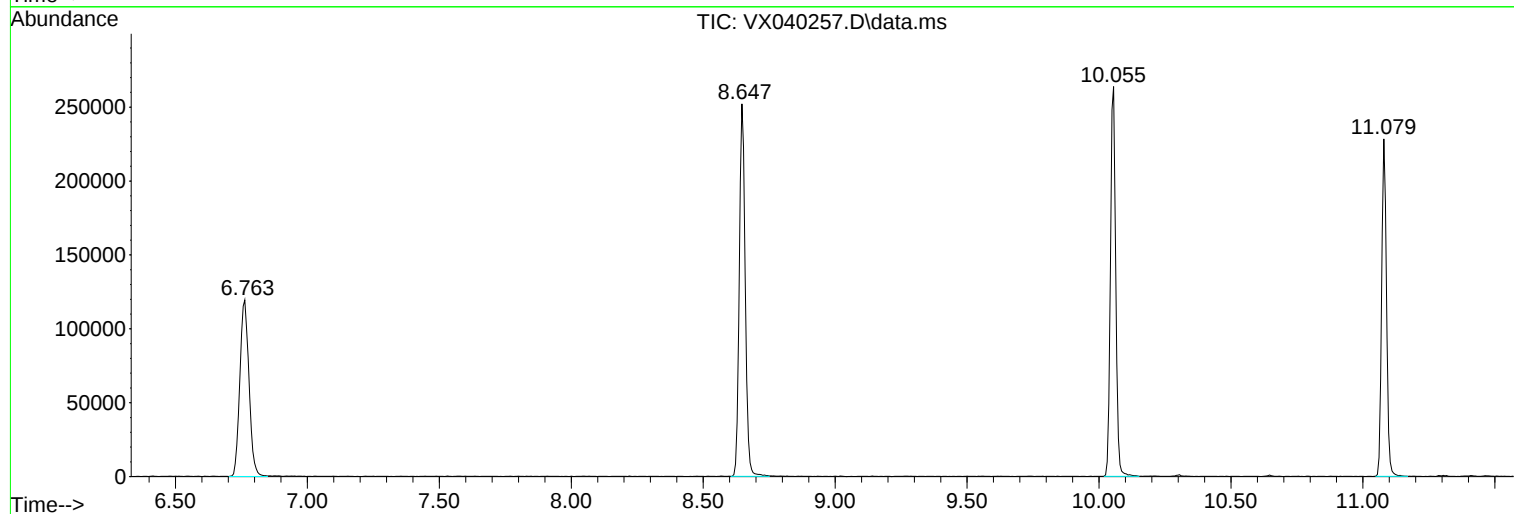
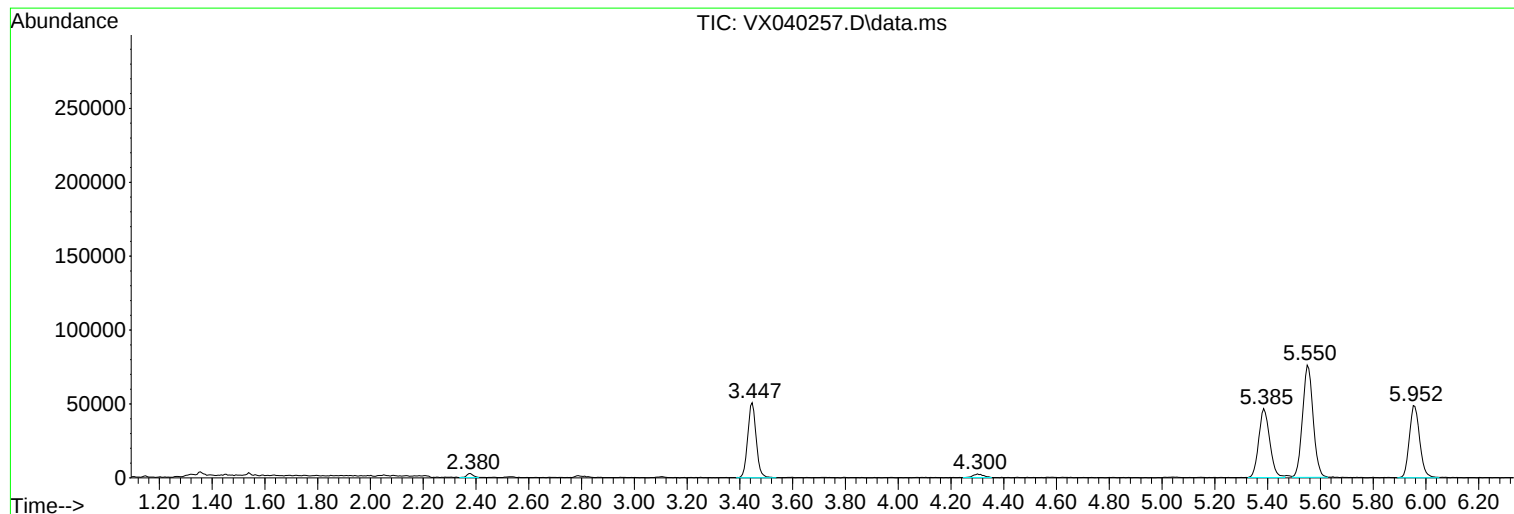
Sum of corrected areas: 2352233

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

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



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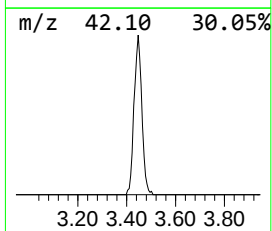
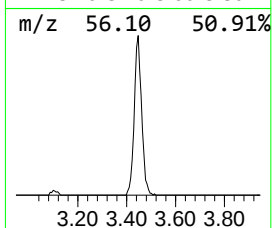
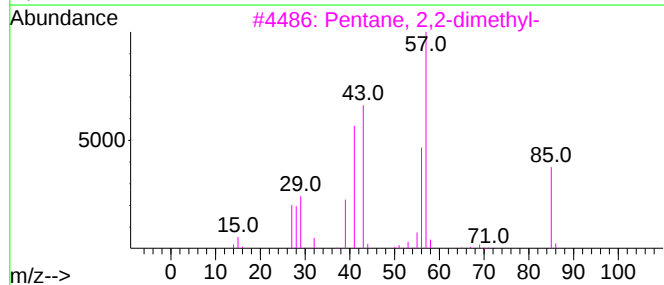
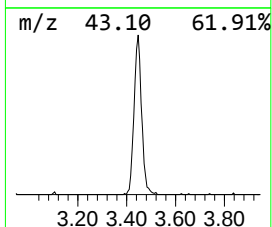
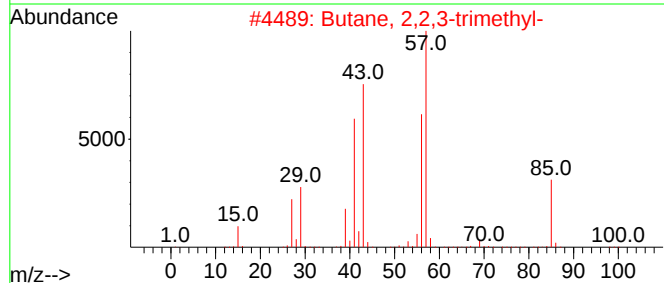
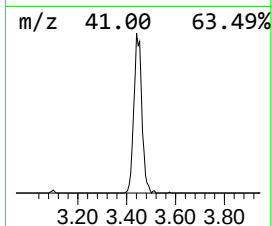
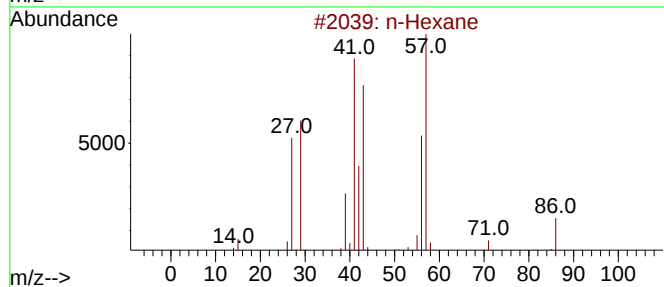
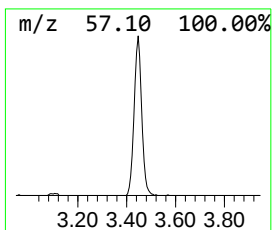
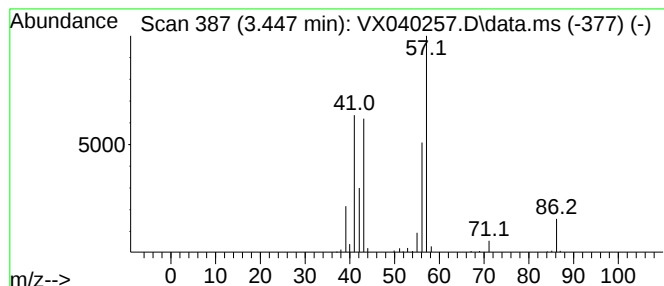
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
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TIC Integration Parameters: LSCINT.P

Peak Number 1 n-Hexane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.447	26.70 ug/l	113171	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexane	86	C6H14	000110-54-3	91
2			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	53
3			Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	50
4			1-Pentene, 4-methyl-	84	C6H12	000691-37-2	47
5			2-Ethyl-oxetane	86	C5H10O	1010386-40-2	43



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexane	3.447	26.7	ug/l	113171	1	5.556	211899	50.0