

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX063021\
 Data File : VX023037.D
 Acq On : 30 Jun 2021 16:30
 Operator : JC/MD
 Sample : M2892-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-6

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\82X062321W.M

Title : SW846 8260

Signal : TIC: VX023037.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.550	234	240	249	rVB2	2736	5148	1.35%	0.257%
2	3.123	326	334	345	rVB2	22771	50656	13.31%	2.533%
3	5.391	694	706	721	rBV	39506	115504	30.34%	5.775%
4	5.562	724	734	746	rVV	67993	184375	48.44%	9.218%
5	5.964	789	800	814	rBV	54869	144643	38.00%	7.232%
6	6.257	837	848	858	rBV2	16099	48413	12.72%	2.420%
7	6.769	922	932	942	rBV	116011	276605	72.67%	13.829%
8	7.988	1125	1132	1139	rVB2	13702	25939	6.81%	1.297%
9	8.110	1146	1152	1162	rVB2	3912	7586	1.99%	0.379%
10	8.653	1230	1241	1251	rBV	246365	380654	100.00%	19.031%
11	10.055	1465	1471	1482	rBV	235587	315340	82.84%	15.766%
12	11.085	1635	1640	1649	rBV	162279	206150	54.16%	10.307%
13	12.024	1789	1794	1802	rVB	199273	239156	62.83%	11.957%

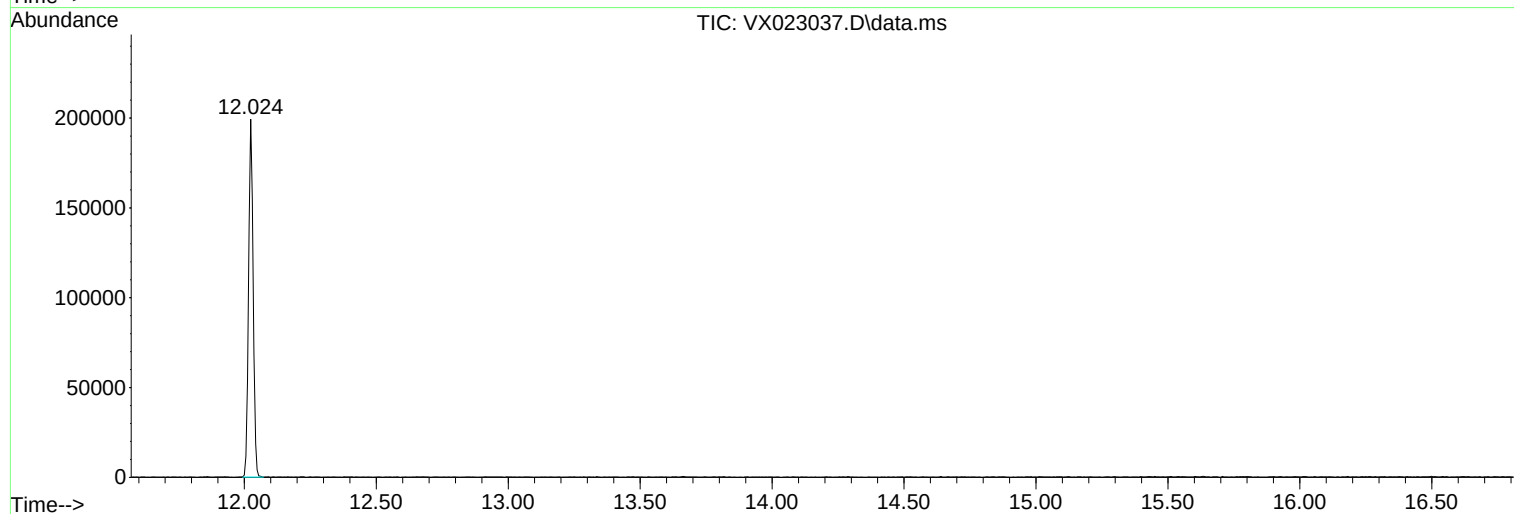
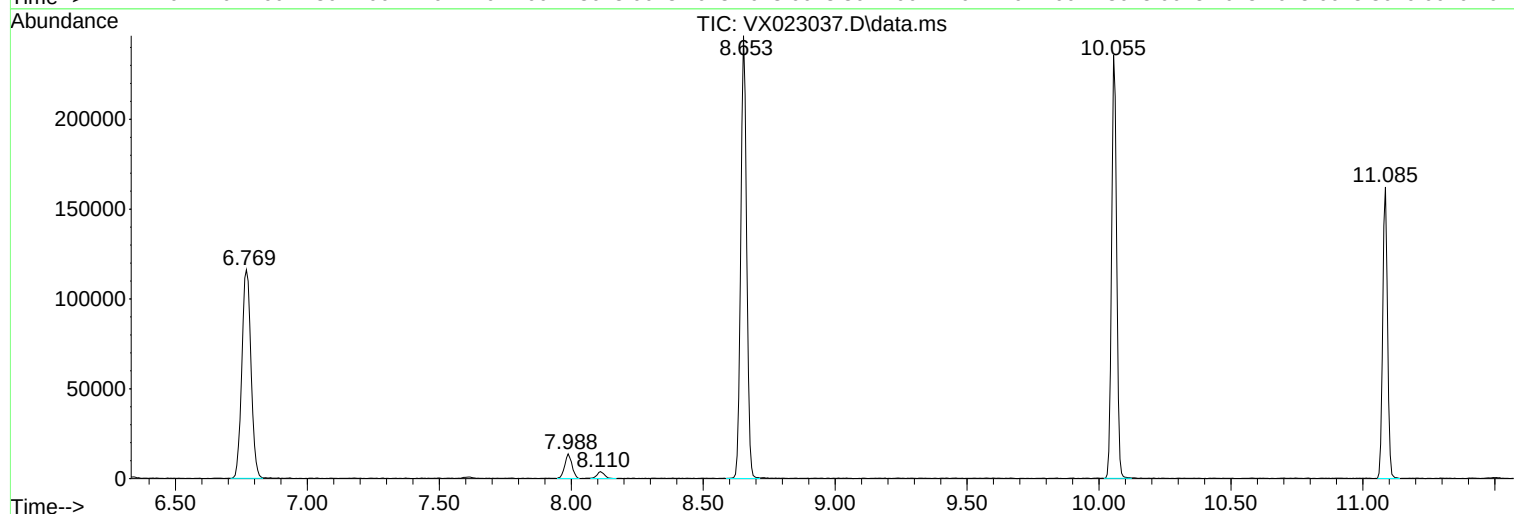
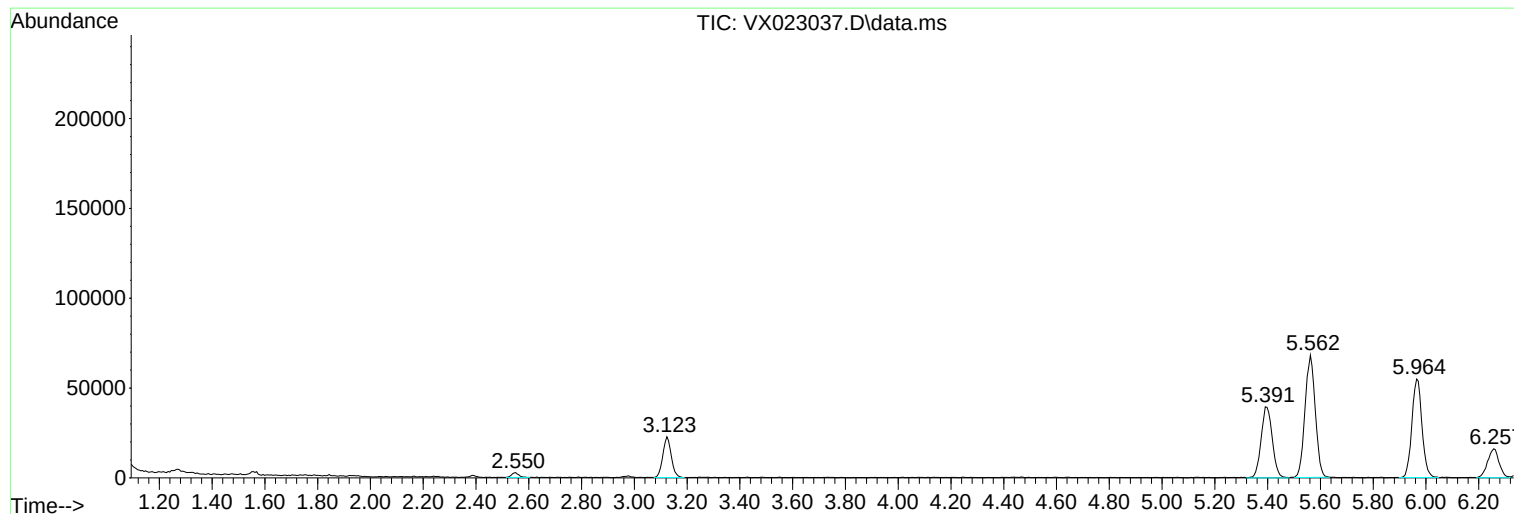
Sum of corrected areas: 2000169

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

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



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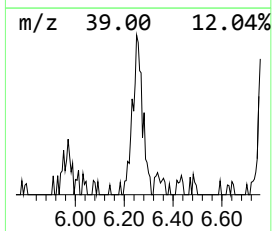
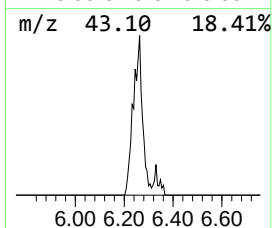
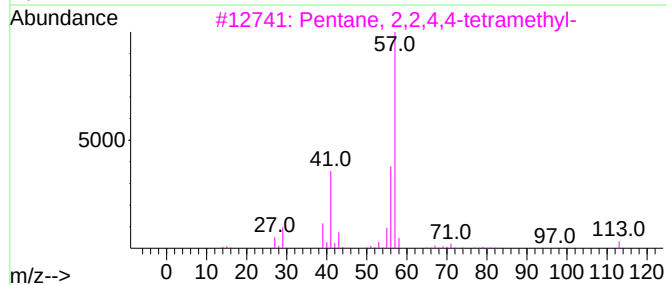
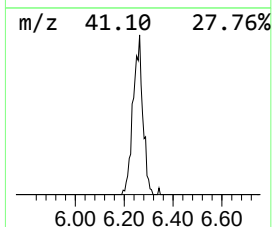
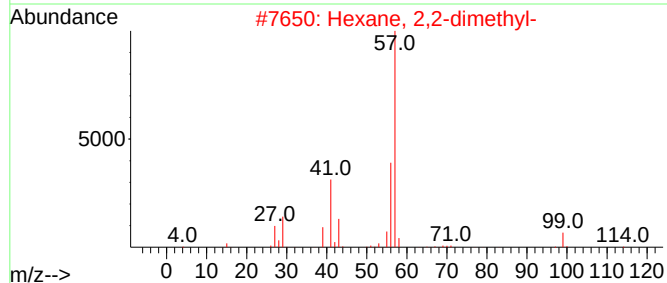
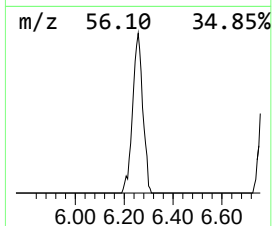
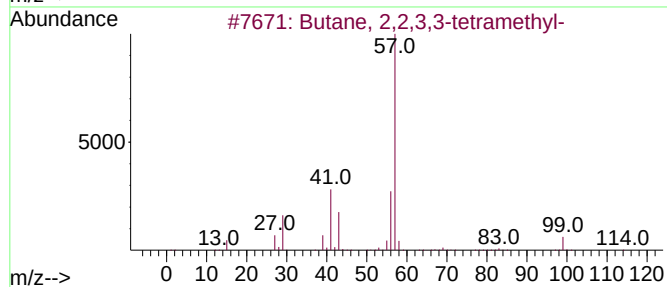
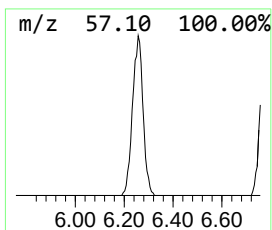
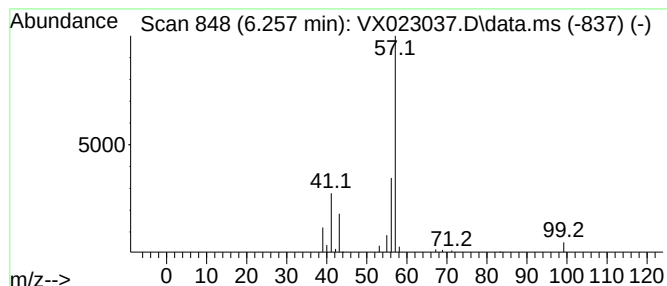
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2,2,3,3-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.257	8.75 ug/l	48413	1,4-Difluorobenzene	6.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
2			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	56
3			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	56
4			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	45
5			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	45



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2,2,3,3...	6.257	8.8	ug/l	48413	2	6.769	276605	50.0