

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX092619\
 Data File : VX012685.D
 Acq On : 26 Sep 2019 19:34
 Operator : JC/SP
 Sample : K5049-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-7

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\82X092319W.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	1.070	26	30	43	rBV	767205	934100	18.83%	7.502%
2	3.027	342	351	366	rBV	733128	1601230	32.28%	12.860%
3	3.186	366	377	400	rVB	2116007	4960519	100.00%	39.840%
4	3.844	473	485	496	rBV	43350	115001	2.32%	0.924%
5	5.490	743	755	767	rBV	98597	281414	5.67%	2.260%
6	5.649	771	781	798	rVV	186640	518466	10.45%	4.164%
7	6.057	837	848	861	rBV	122635	321244	6.48%	2.580%
8	6.337	883	894	903	rBV3	45582	135950	2.74%	1.092%
9	6.850	968	978	994	rBV	289595	689006	13.89%	5.534%
10	8.051	1167	1175	1182	rBV	28670	55393	1.12%	0.445%
11	8.709	1276	1283	1298	rBV	575520	913083	18.41%	7.333%
12	10.111	1507	1513	1519	rBV	533685	753017	15.18%	6.048%
13	11.135	1675	1681	1690	rBV	411917	546693	11.02%	4.391%
14	12.074	1829	1835	1843	rBV	513683	626033	12.62%	5.028%

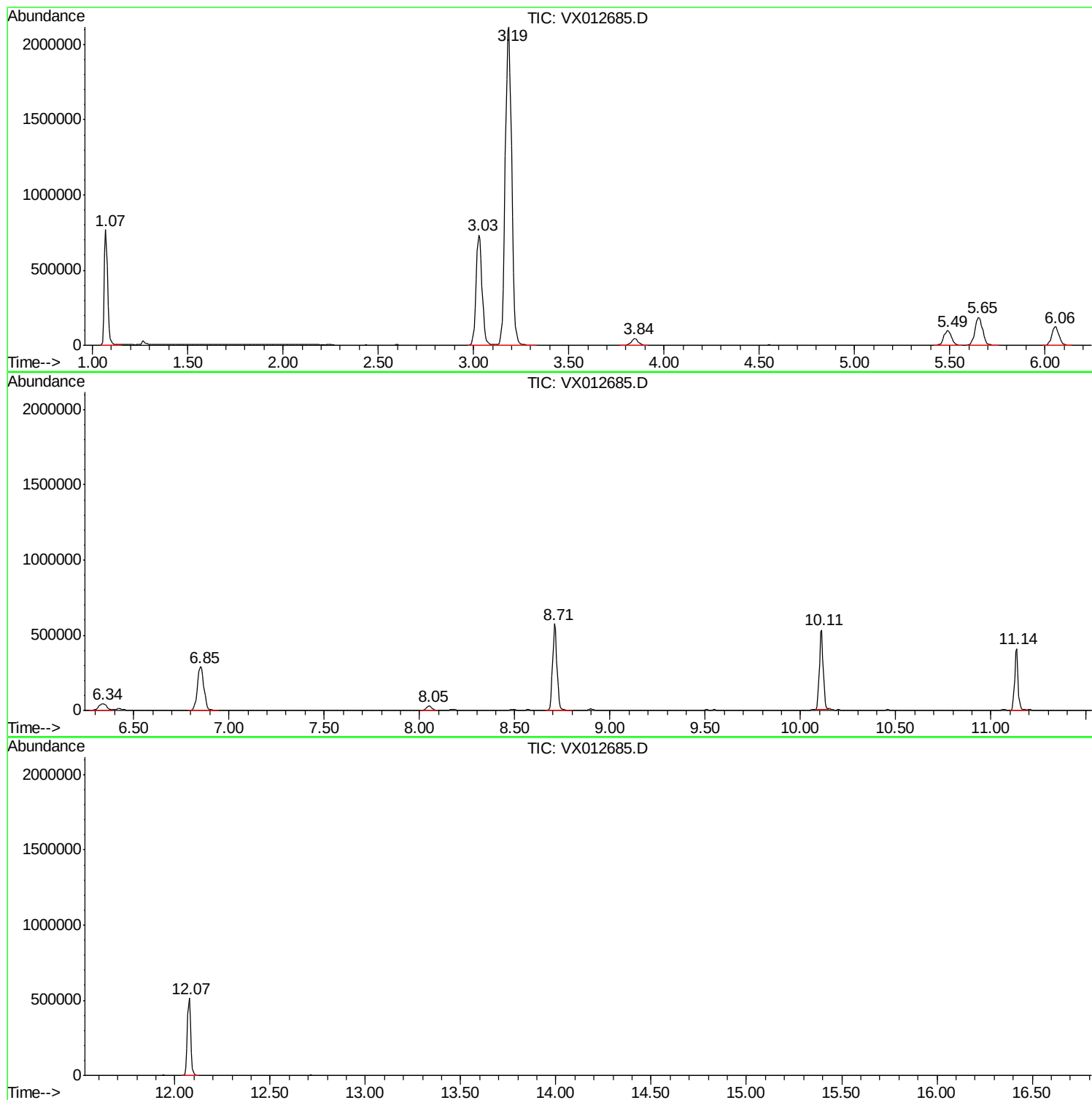
Sum of corrected areas: 12451149

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X092319W.M
Quant Title : SW846 8260

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



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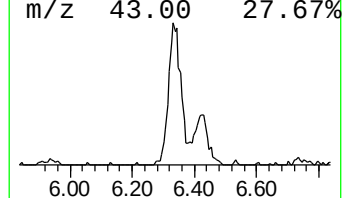
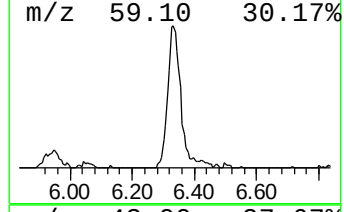
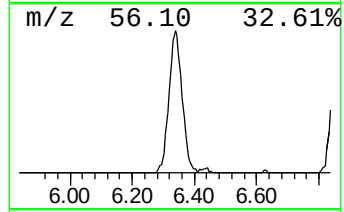
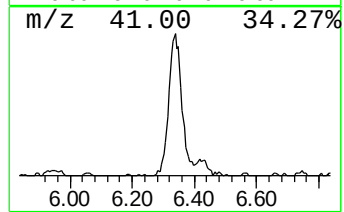
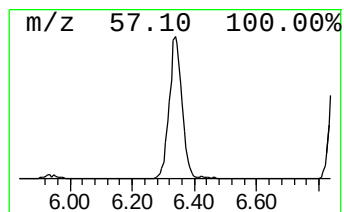
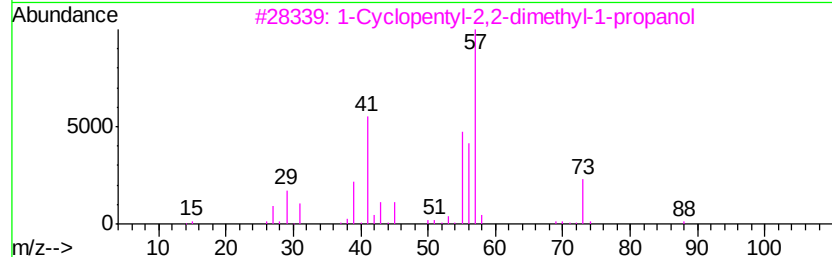
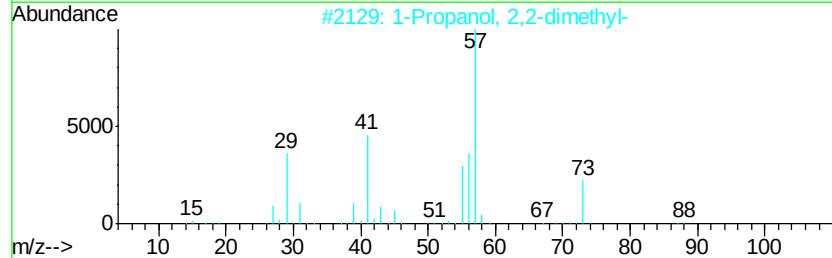
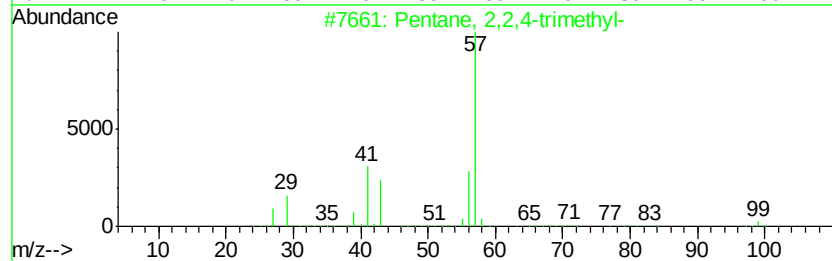
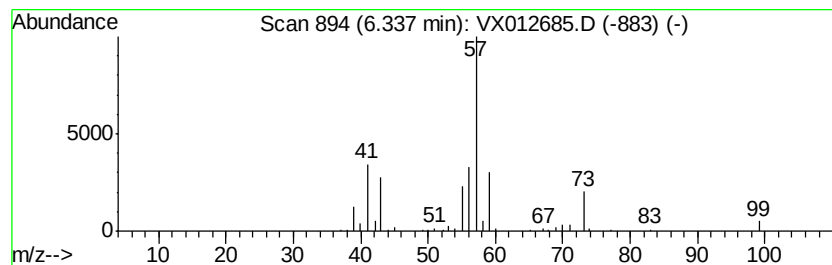
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TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.34 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.34	9.87 ug/l	135950	1,4-Difluorobenzene	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	43
2		1-Propanol, 2,2-dimethyl-	88	C5H12O	000075-84-3	42
3		1-Cyclopentyl-2,2-dimethyl-1-pro...	156	C10H20O	337966-85-5	40
4		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	38
5		Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	38



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.34	6.34	9.9	ug/l	135950	2	6.85	689006	50.0