

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121918\
 Data File : VX006643.D
 Acq On : 19 Dec 2018 18:04
 Operator : JC/MD
 Sample : J6484-17
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 977-FB121818

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.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	1.611	67	75	79	rVB2	28857	53174	1.78%	0.312%
2	2.446	207	212	217	rVB	29337	48042	1.60%	0.282%
3	3.025	300	307	312	rBV	15847	38563	1.29%	0.227%
4	4.854	600	607	616	rVB	19573	52883	1.77%	0.311%
5	5.507	699	714	726	rBV	341314	986804	32.95%	5.798%
6	5.665	729	740	752	rVV2	607006	1694350	56.58%	9.954%
7	6.068	796	806	820	rBV	368910	956755	31.95%	5.621%
8	6.860	925	936	950	rBV	926369	2158748	72.08%	12.683%
9	8.713	1232	1240	1258	rBV	1926736	2994763	100.00%	17.594%
10	10.116	1463	1470	1480	rBV	1845873	2583265	86.26%	15.177%
11	11.134	1632	1637	1647	rBV	1561395	2025474	67.63%	11.900%
12	12.079	1786	1792	1799	rBV	2037435	2500686	83.50%	14.692%
13	12.274	1818	1824	1840	rBV	526140	873414	29.16%	5.131%
14	13.152	1963	1968	1973	rBV2	39875	54273	1.81%	0.319%

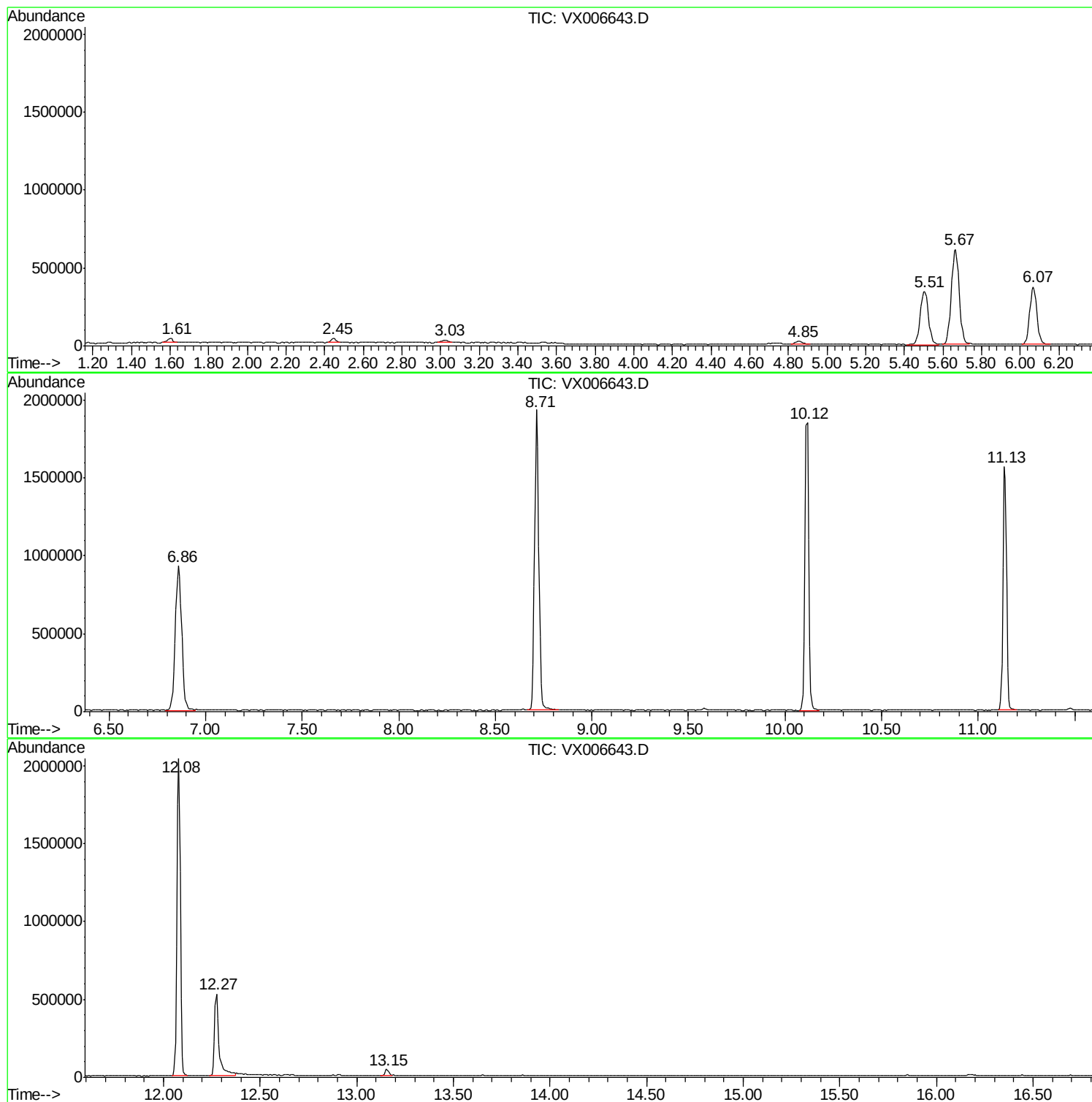
Sum of corrected areas: 17021194

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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



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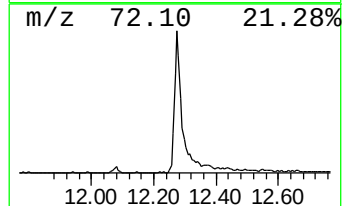
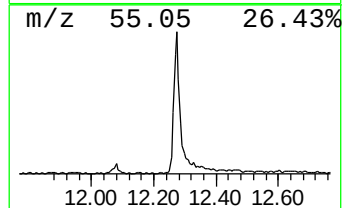
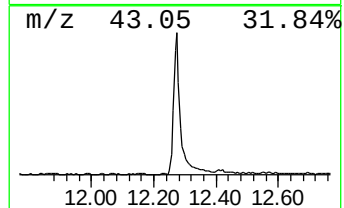
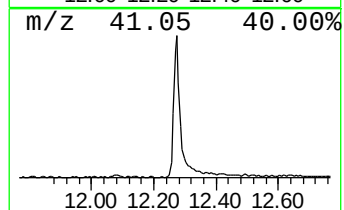
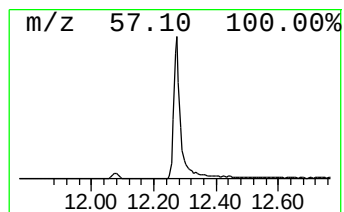
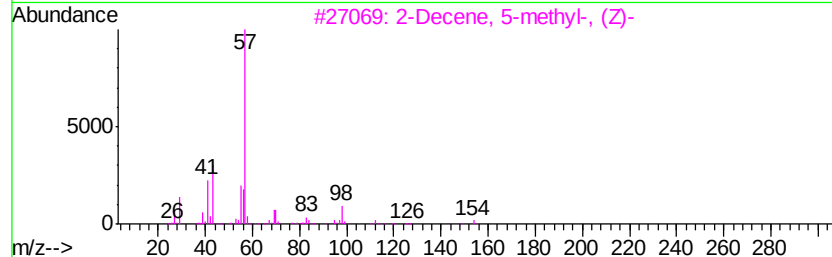
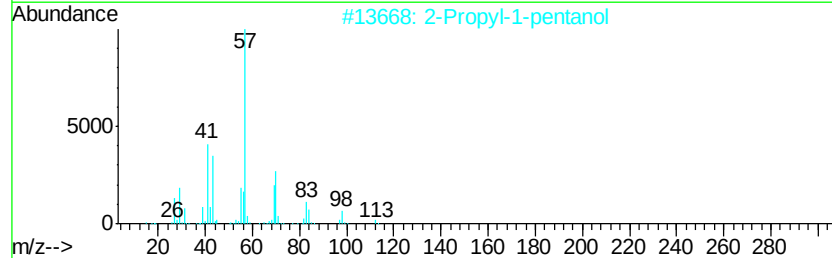
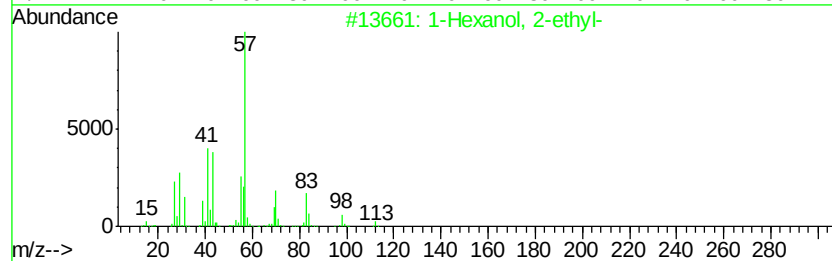
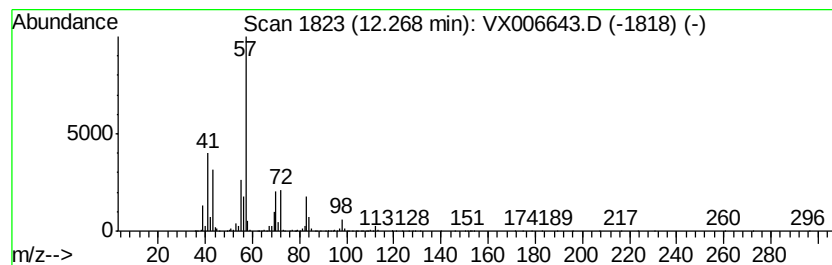
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	17.46 ug/l	873414	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
2		2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53
3		2-Decene, 5-methyl-, (Z)-	154	C11H22	074645-86-6	38
4		1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	38
5		1-Decene, 2,4-dimethyl-	168	C12H24	055170-80-4	38



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Hexanol, 2-ethyl-	12.27	17.5	ug/l	873414	4	12.08	2500690	50.0