

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW070318\
 Data File : VW003706.D
 Acq On : 03 Jul 2018 18:26
 Operator : SY/AP
 Sample : J3813-02
 Misc : 5.03G/5ML/MSVOA W/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 IDW-S02-SRI3

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_W\METHOD\82W070218S.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	1.751	18	27	29	rBV	5629	11988	1.57%	0.900%
2	1.831	29	40	55	rVV	206812	765431	100.00%	57.447%
3	4.123	405	416	427	rBV	9498	26731	3.49%	2.006%
4	7.952	1037	1044	1054	rVB	30445	65117	8.51%	4.887%
5	8.311	1095	1103	1112	rBV	17456	37794	4.94%	2.836%
6	8.848	1183	1191	1200	rBV2	41137	82658	10.80%	6.204%
7	10.329	1427	1434	1441	rBV	45442	82982	10.84%	6.228%
8	11.634	1641	1648	1658	rBV	59179	99909	13.05%	7.498%
9	12.621	1803	1810	1818	rVB	34191	56459	7.38%	4.237%
10	13.566	1958	1965	1971	rVB	50985	84645	11.06%	6.353%
11	13.658	1973	1980	1986	rBV	6484	10560	1.38%	0.793%
12	14.335	2086	2091	2097	rVB2	5454	8148	1.06%	0.612%

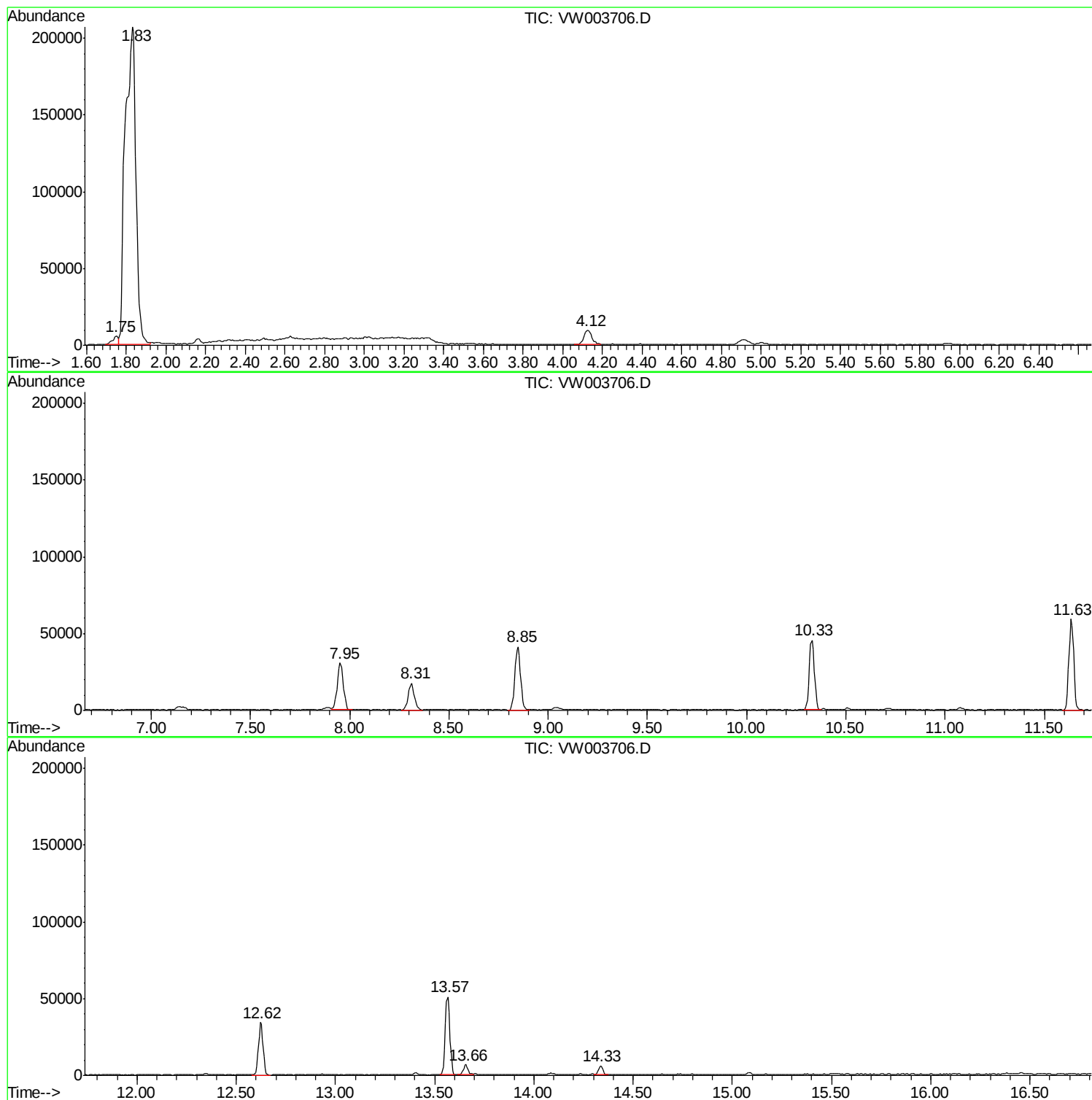
Sum of corrected areas: 1332422

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W070218S.M
Quant Title : SW846 8260

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



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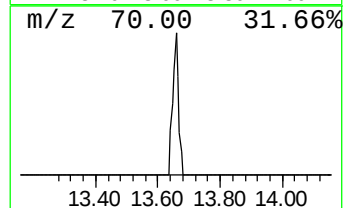
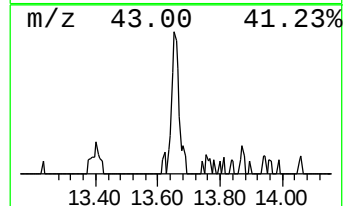
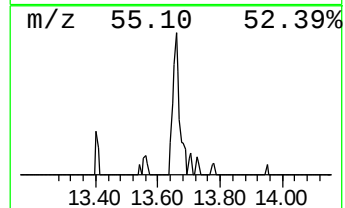
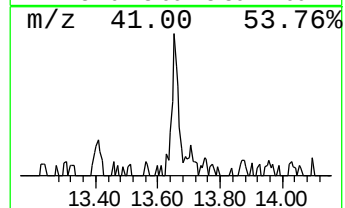
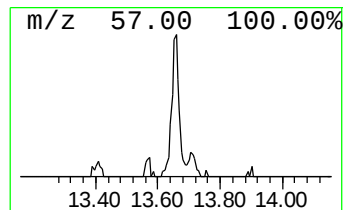
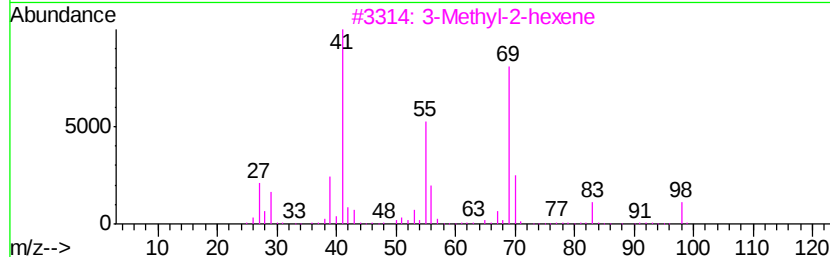
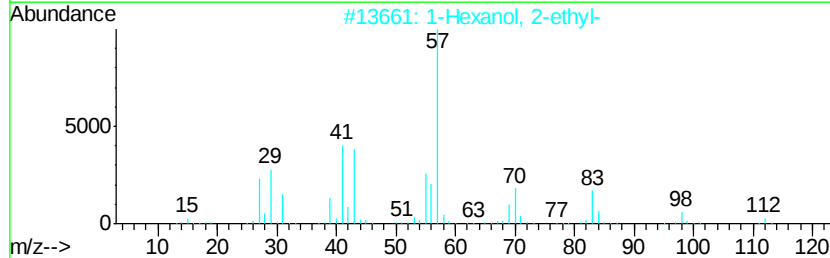
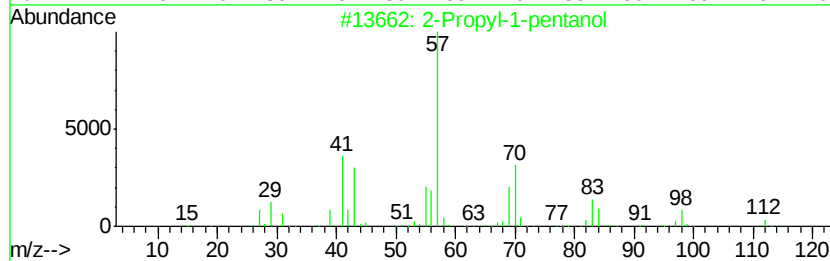
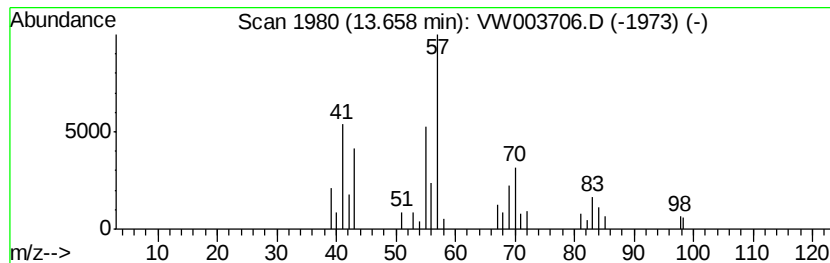
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Propyl-1-pentanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.66	6.24 ug/l	10560	1,4-Dichlorobenzene-d4	13.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	59
2			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	50
3			3-Methyl-2-hexene	98	C7H14	017618-77-8	47
4			2-Hexene, 3-methyl-, (Z)-	98	C7H14	010574-36-4	47
5			1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	43



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Propyl-1-pentanol	13.66	6.2	ug/l	10560	4	13.57	84645	50.0