

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111720\
 Data File : VD067656.D
 Acq On : 17 Nov 2020 12:21
 Operator : VA/SY
 Sample : L4769-10
 Misc : 3.86G/5.00ml/MSVOA D/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 ETGI-359

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_D\METHOD\82D110320S.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	7.920	1011	1020	1025	rBV	114523	274351	18.83%	3.388%
2	7.985	1025	1031	1044	rVB2	344841	789159	54.16%	9.746%
3	8.338	1082	1091	1104	rVB	154570	349857	24.01%	4.321%
4	8.873	1170	1182	1195	rBV	476471	990015	67.95%	12.227%
5	10.344	1424	1432	1447	rBV	756686	1395523	95.78%	17.235%
6	11.649	1643	1654	1667	rBV	799453	1365327	93.71%	16.862%
7	12.638	1813	1822	1830	rBV	674658	1097371	75.32%	13.553%
8	13.579	1975	1982	1992	rVB	880986	1457041	100.00%	17.995%
9	13.673	1992	1998	2004	rBV2	44380	70997	4.87%	0.877%
10	14.043	2055	2061	2067	rBV2	18888	28902	1.98%	0.357%
11	14.102	2067	2071	2078	rVB4	12567	20596	1.41%	0.254%
12	14.249	2090	2096	2102	rBV	163843	257977	17.71%	3.186%

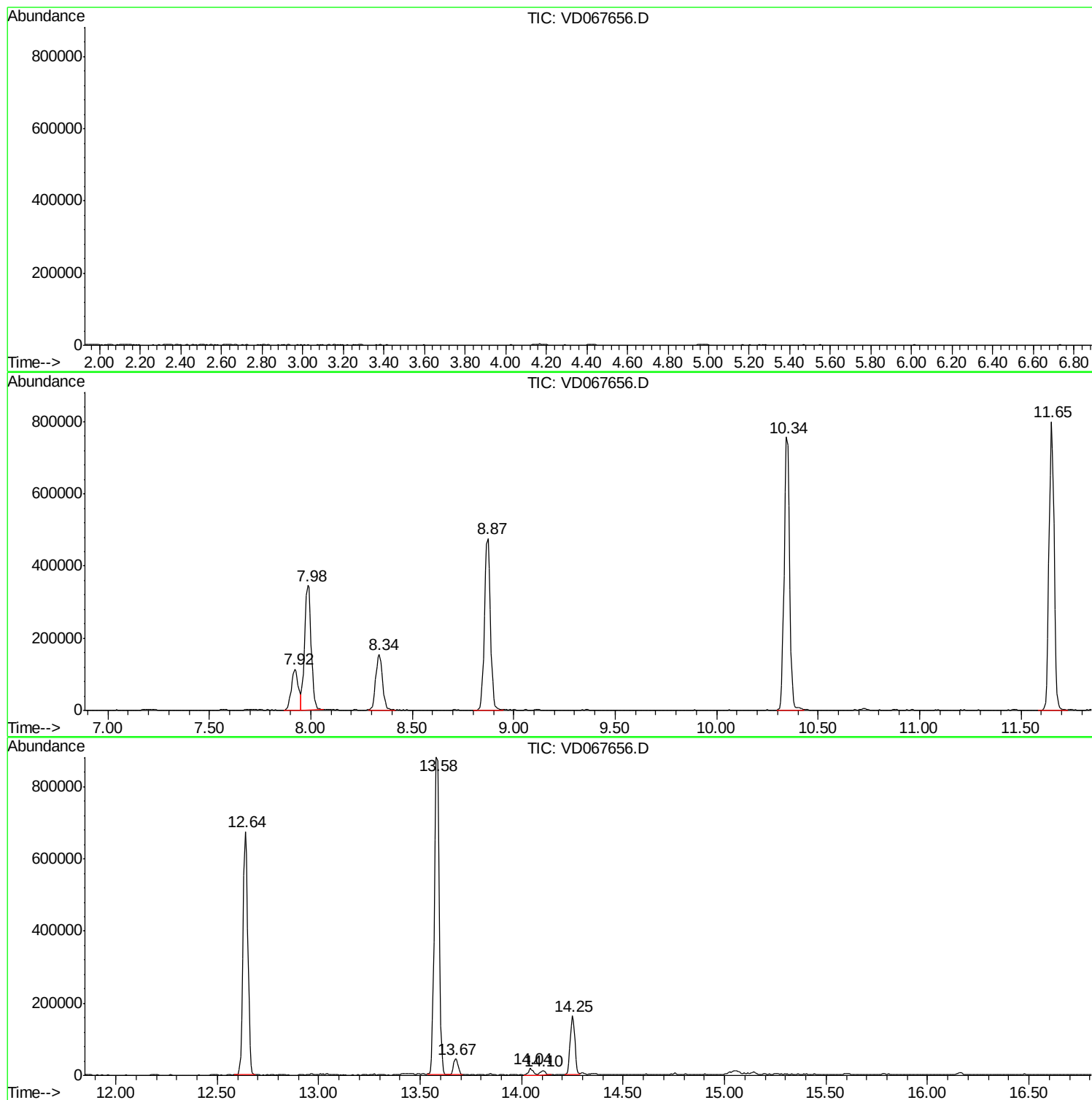
Sum of corrected areas: 8097116

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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



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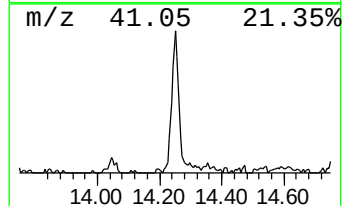
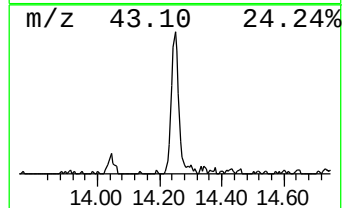
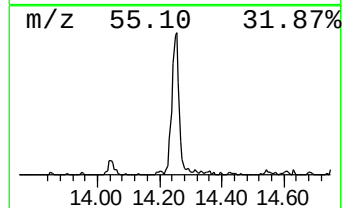
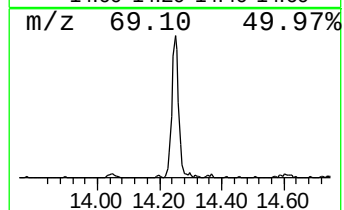
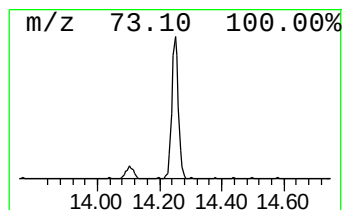
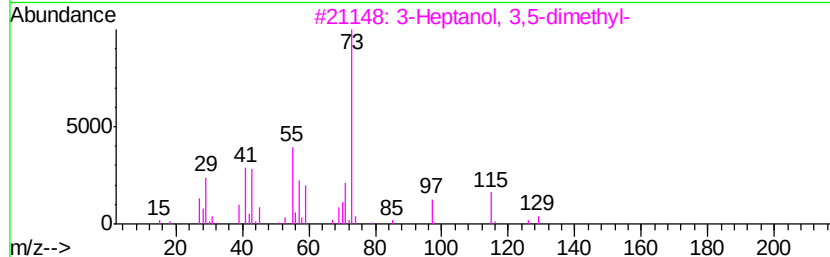
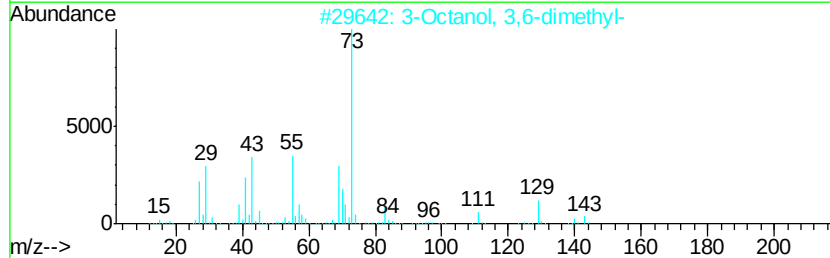
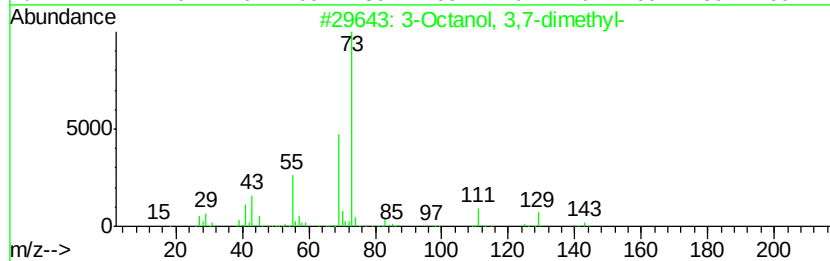
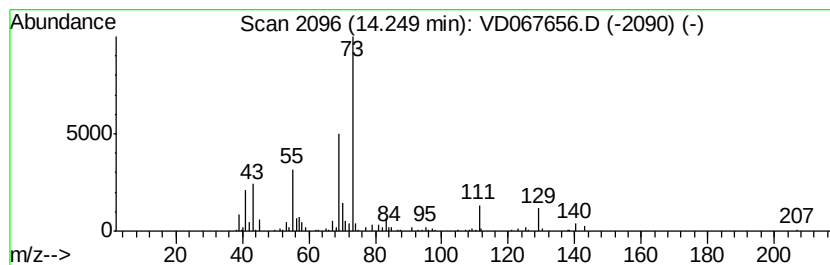
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Quant Title : SW846 8260

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

Peak Number 1 3-Octanol, 3,7-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.25	8.85 ug/l	257977	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	78
2		3-Octanol, 3,6-dimethyl-	158	C10H22O	000151-19-9	32
3		3-Heptanol, 3,5-dimethyl-	144	C9H20O	019549-74-7	23
4		3-Dodecanol, 3,7,11-trimethyl-	228	C15H32O	007278-65-1	12
5		Isopropyl 5,11-dihydroxy-3,7,11-...	314	C18H34O4	1000223-34-1	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Octanol, 3,7-di...	14.25	8.8	ug/l	257977	4	13.58	1457040	50.0