

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD091820\
 Data File : VD066765.D
 Acq On : 18 Sep 2020 14:46
 Operator : VA/SY
 Sample : L4053-04
 Misc : 5.01G/5.00ml/MSVOA D/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 VNJ-231

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_D\METHOD\82D091520S.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	7.914	1009	1019	1024	rBV	293868	700547	8.98%	3.698%
2	7.979	1024	1030	1040	rVB	509944	1147194	14.70%	6.056%
3	8.326	1077	1089	1102	rBV	264489	608265	7.79%	3.211%
4	8.861	1171	1180	1191	rBV	808387	1570592	20.13%	8.291%
5	10.338	1423	1431	1438	rBV	1237197	2153026	27.59%	11.366%
6	11.644	1645	1653	1665	rBV	1062802	1812765	23.23%	9.570%
7	12.461	1784	1792	1814	rBV	4523666	7803532	100.00%	41.195%
8	12.632	1814	1821	1829	rVB	879547	1486821	19.05%	7.849%
9	13.579	1975	1982	1990	rBV	908709	1491730	19.12%	7.875%
10	14.044	2056	2061	2067	rBV2	107318	168623	2.16%	0.890%

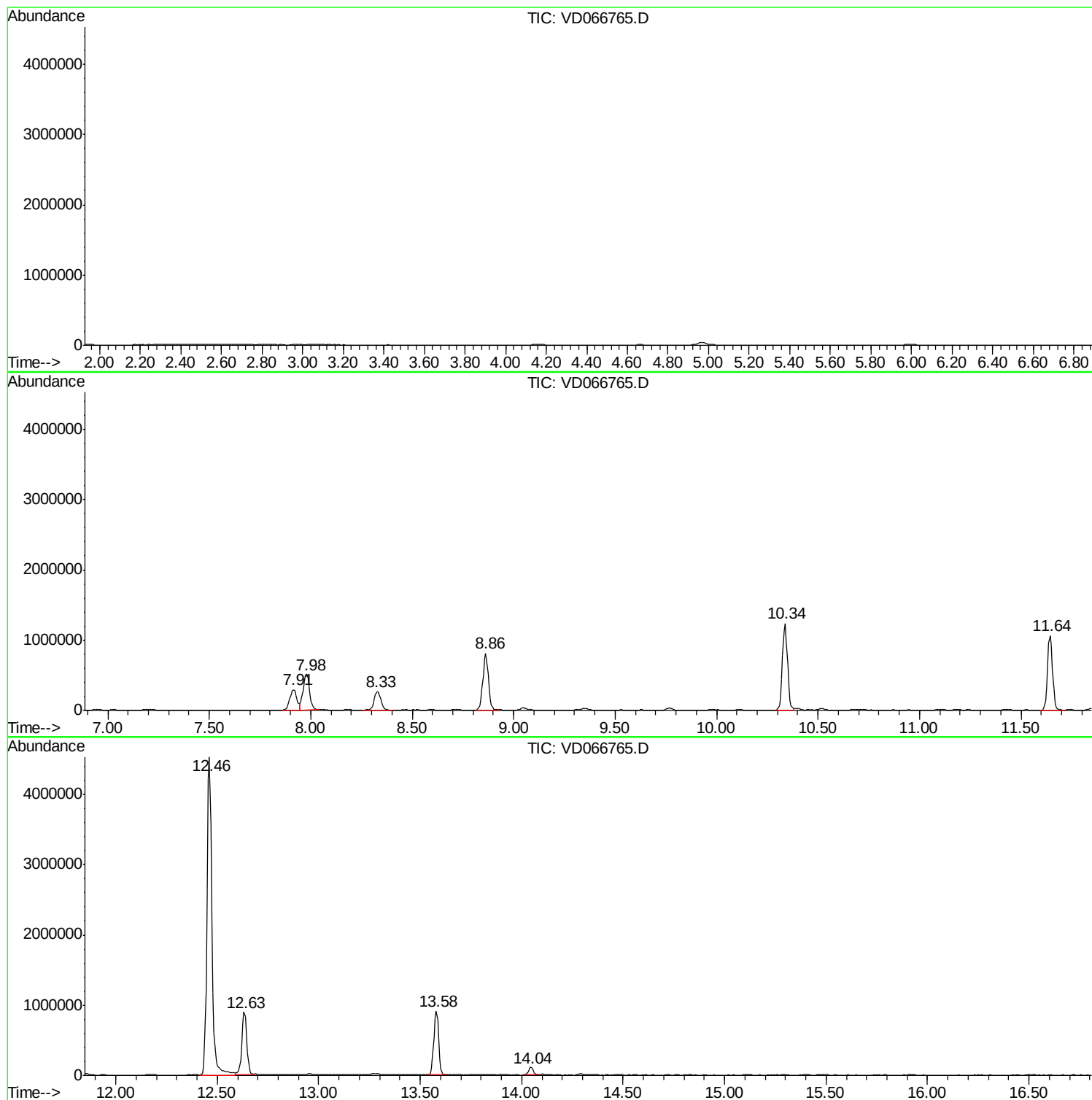
Sum of corrected areas: 18943095

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D091520S.M
Quant Title : SW846 8260

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



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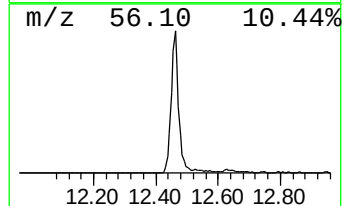
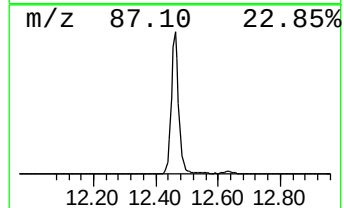
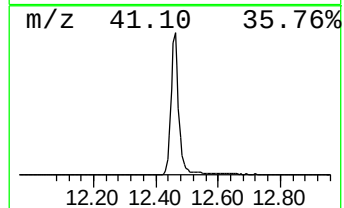
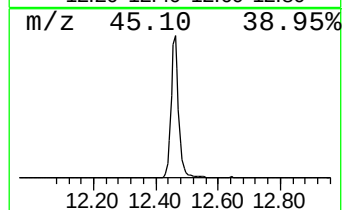
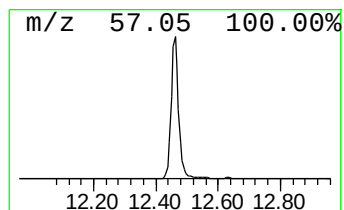
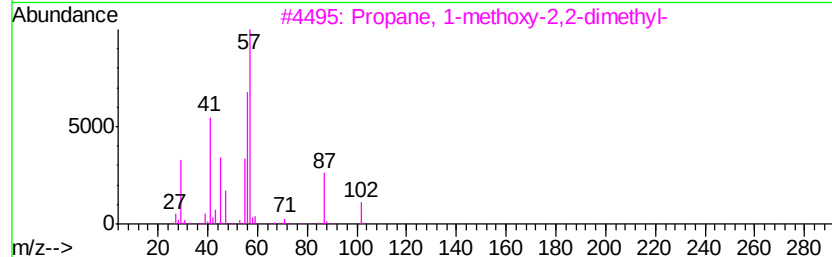
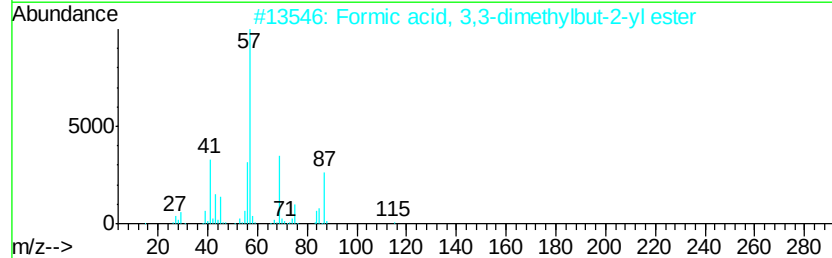
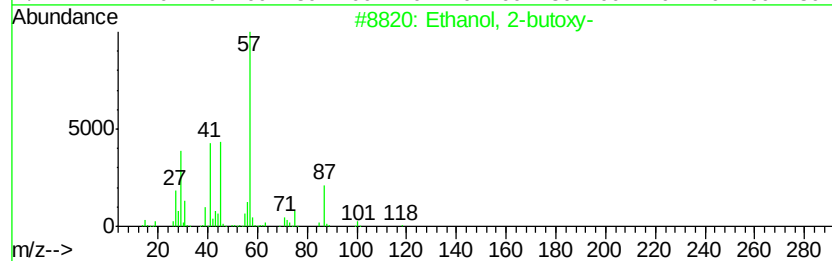
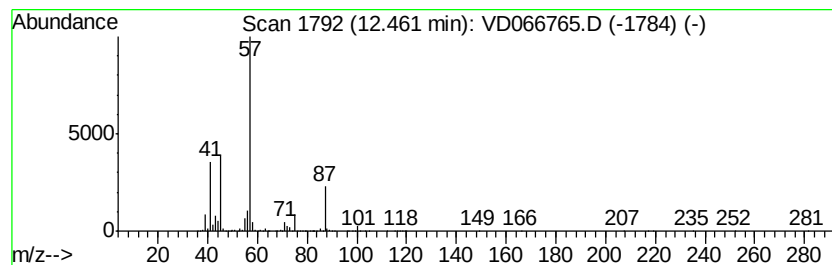
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D091520S.M
Quant Title : SW846 8260

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

Peak Number 1 Ethanol, 2-butoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.46	215.24 ug/l	7803530	Chlorobenzene-d5	11.64		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	91
2		Formic acid, 3,3-dimethylbut-2-yl...	130	C7H14O2	1000368-20-5	64
3		Propane, 1-methoxy-2,2-dimethyl-	102	C6H14O	001118-00-9	50
4		3-Buten-2-ol	72	C4H8O	000598-32-3	43
5		Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	33



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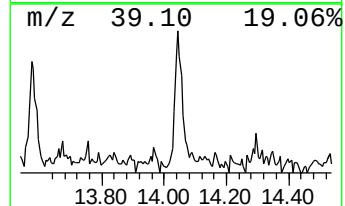
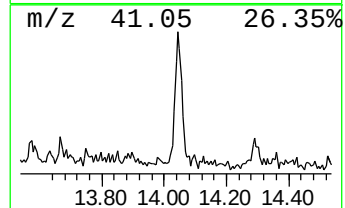
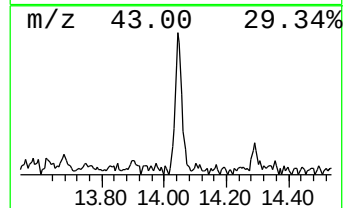
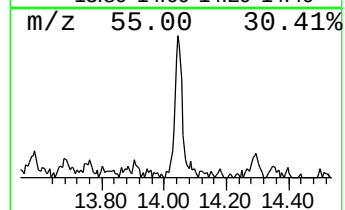
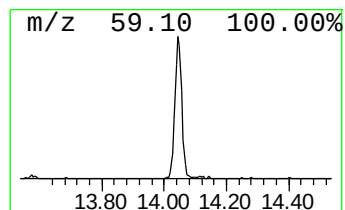
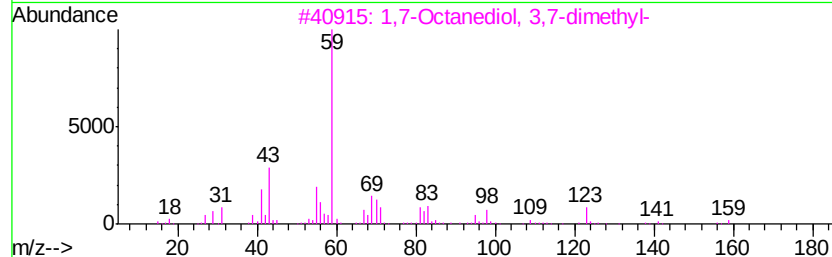
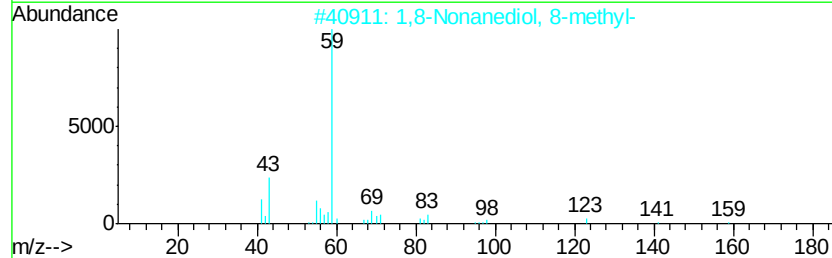
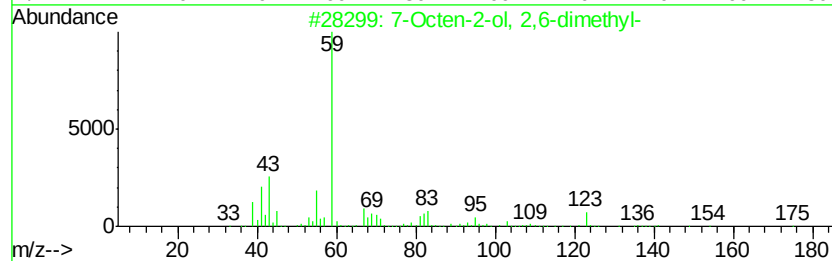
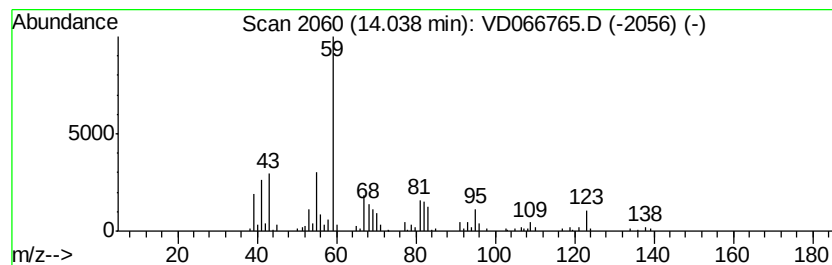
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TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown14.04 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.04	5.65 ug/l	168623	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	7-Octen-2-ol, 2,6-dimethyl-	156	C10H20O	018479-58-8	45
2		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	43
3		1,7-Octanediol, 3,7-dimethyl-	174	C10H22O2	000107-74-4	40
4		1-Cyclohexyl-2-methyl-2-propanol	156	C10H20O	005531-30-6	38
5		Acetic acid, cyanohydroxyimino-,...	128	C4H4N2O3	061295-92-9	38



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-butoxy-	12.46	215.2	ug/l	7803530	3	11.64	1812770	50.0
unknown14.04	14.04	5.7	ug/l	168623	4	13.58	1491730	50.0