

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD112619\
 Data File : VD064382.D
 Acq On : 26 Nov 2019 14:19
 Operator : VA/SY
 Sample : K6012-01
 Misc : 8.97G/5.00ml/MSVOA D/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 211

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\82D110819S.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	1009	1020	1025	rBV	481673	1176373	33.64%	6.296%
2	7.979	1025	1030	1042	rVB	898714	2016396	57.67%	10.791%
3	8.338	1081	1091	1105	rBV	402845	900400	25.75%	4.819%
4	8.867	1170	1181	1191	rBV	1265895	2510189	71.79%	13.434%
5	10.343	1421	1432	1450	rBV	1951367	3496706	100.00%	18.713%
6	10.732	1489	1498	1505	rBV2	21356	42876	1.23%	0.229%
7	11.649	1645	1654	1666	rBV	1693217	2982070	85.28%	15.959%
8	12.637	1811	1822	1831	rBV	1400548	2318295	66.30%	12.407%
9	13.584	1974	1983	1991	rVB	1678516	2845576	81.38%	15.229%
10	14.420	2112	2125	2143	rBV5	108891	396736	11.35%	2.123%

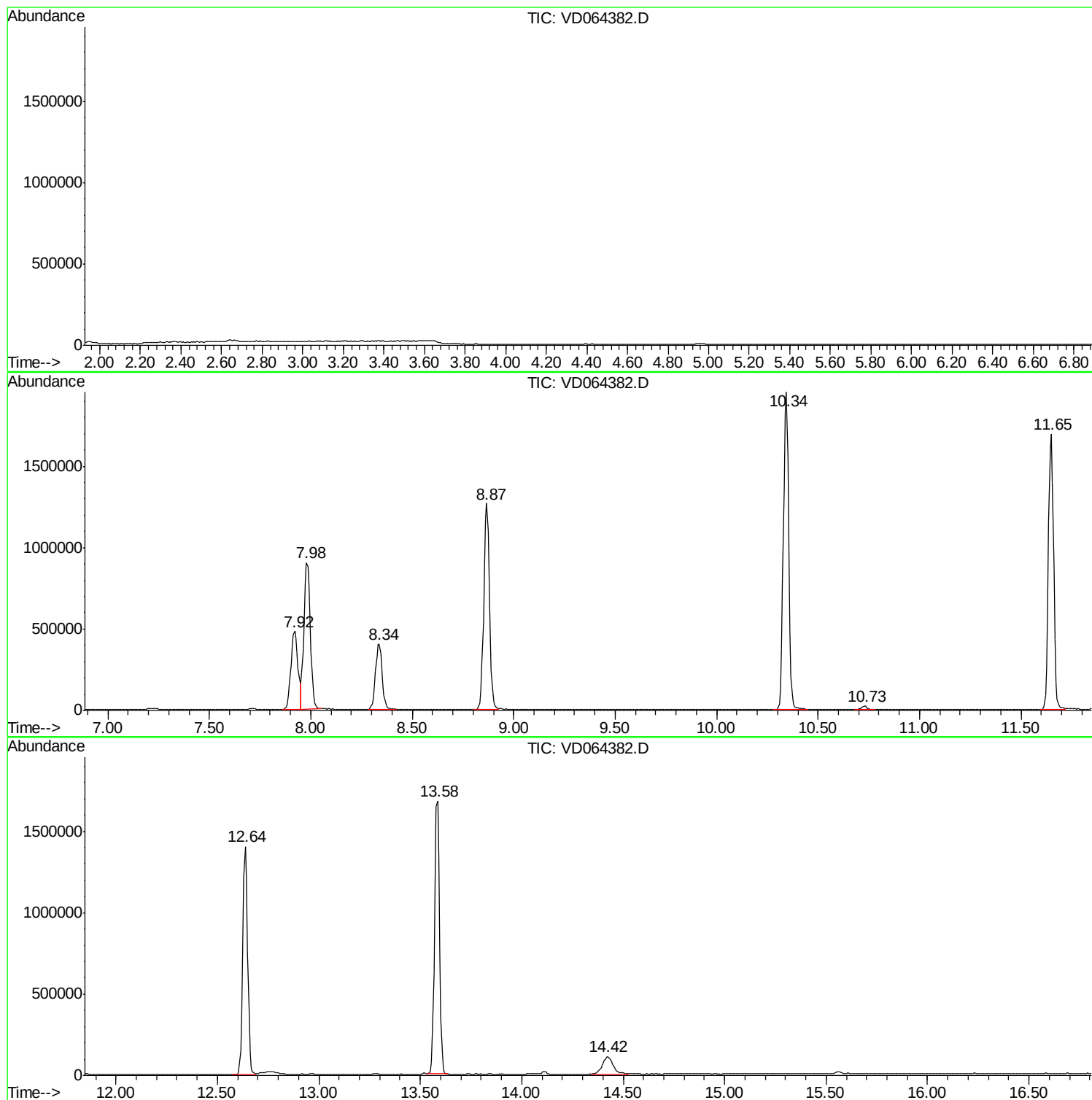
Sum of corrected areas: 18685617

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110819S.M
Quant Title : SW846 8260

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



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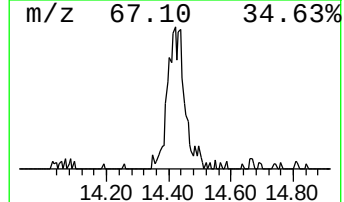
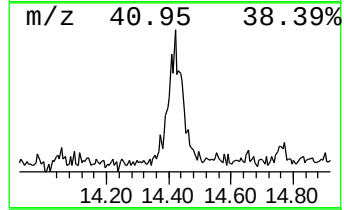
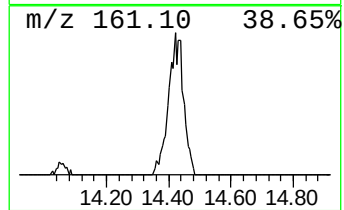
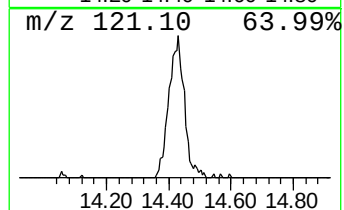
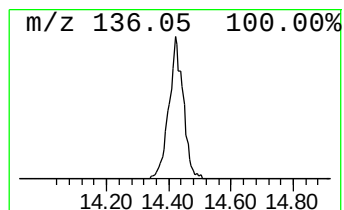
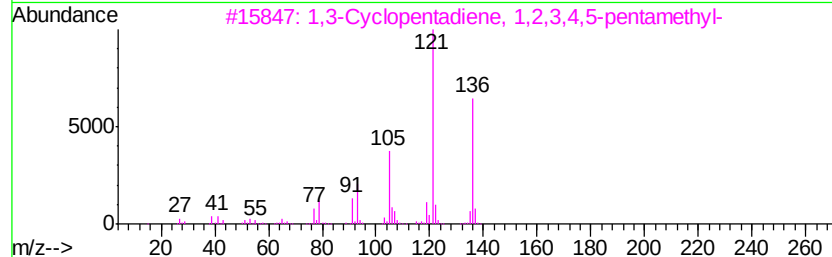
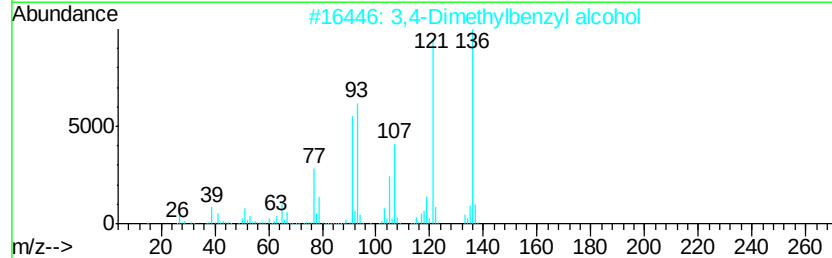
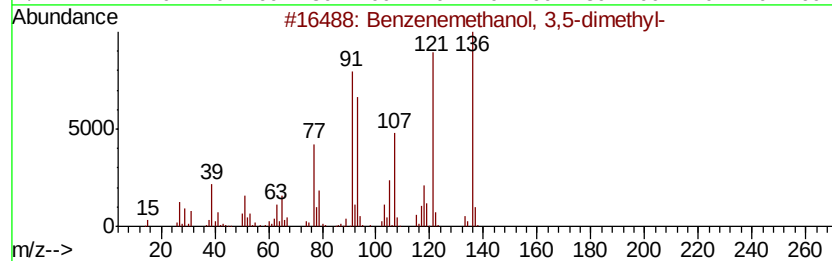
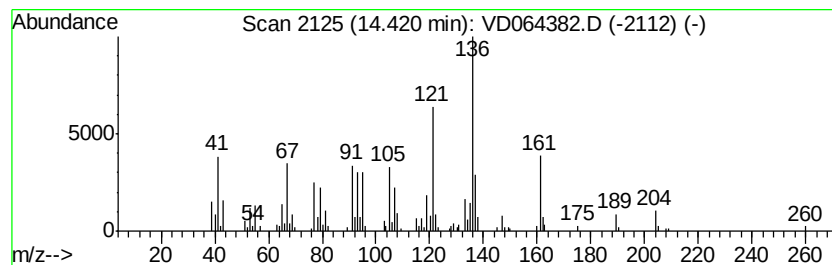
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Quant Title : SW846 8260

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

Peak Number 1 unknown14.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	6.97 ug/l	396736	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenemethanol, 3,5-dimethyl-	136	C9H12O	027129-87-9	49
2		3,4-Dimethylbenzyl alcohol	136	C9H12O	006966-10-5	46
3		1,3-Cyclopentadiene, 1,2,3,4,5-p...	136	C10H16	004045-44-7	43
4		2,4-Dimethylanisole	136	C9H12O	006738-23-4	41
5		2,3-Dimethylanisole	136	C9H12O	002944-49-2	41



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown14.42	14.42	7.0	ug/l	396736	4	13.58	2845580	50.0