

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW083018\
 Data File : VW005018.D
 Acq On : 30 Aug 2018 15:08
 Operator : SY/AP
 Sample : J4654-01
 Misc : 2.60G/5ML/MSVOA W/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ENY-SB-03-1.5-2

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_W\METHOD\82W082918S.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	1.794	19	34	55	rBV	3681102	6445732	100.00%	52.925%
2	4.922	534	547	562	rBV2	64745	210144	3.26%	1.725%
3	7.147	900	912	924	rBV2	28488	84165	1.31%	0.691%
4	7.891	1025	1034	1039	rBV	223768	518441	8.04%	4.257%
5	7.958	1039	1045	1057	rVB	335662	745737	11.57%	6.123%
6	8.312	1088	1103	1115	rBV	212317	469612	7.29%	3.856%
7	8.848	1181	1191	1207	rBV	390637	798627	12.39%	6.557%
8	10.329	1425	1434	1442	rBV	555114	1006336	15.61%	8.263%
9	10.750	1488	1503	1514	rBV2	106066	443670	6.88%	3.643%
10	11.640	1641	1649	1658	rBV	354322	623844	9.68%	5.122%
11	12.622	1803	1810	1818	rBV	186127	323314	5.02%	2.655%
12	13.128	1881	1893	1910	rVB4	68594	313661	4.87%	2.575%
13	13.567	1959	1965	1972	rBV	119628	195732	3.04%	1.607%

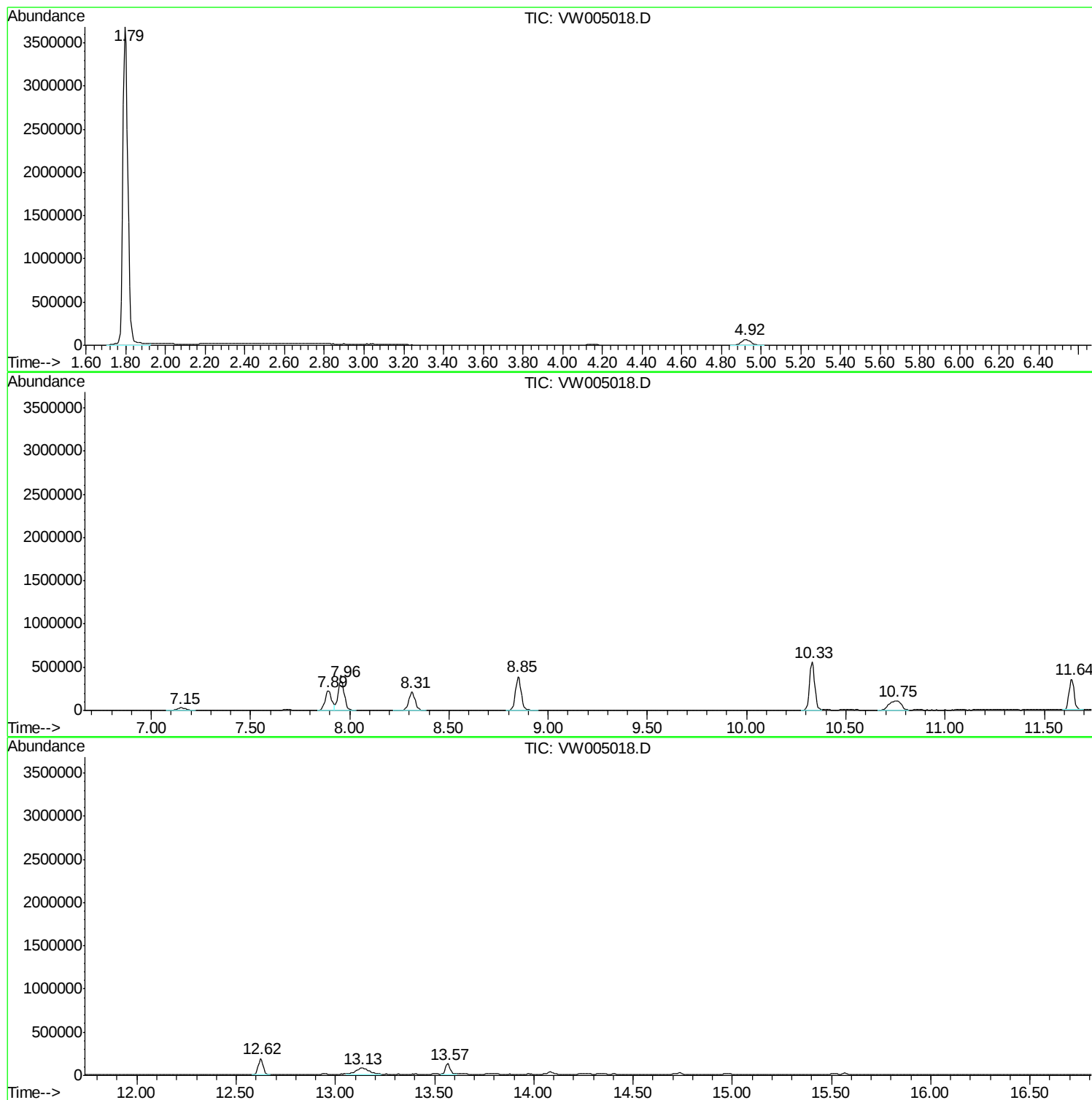
Sum of corrected areas: 12179015

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W082918S.M
Quant Title : SW846 8260

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



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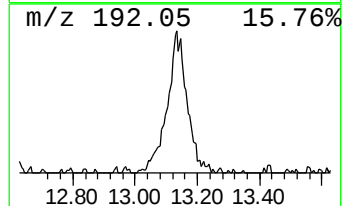
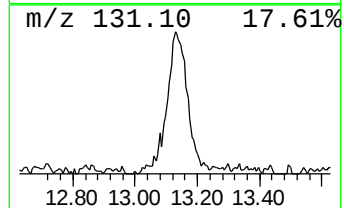
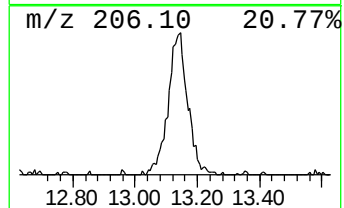
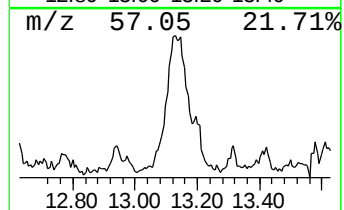
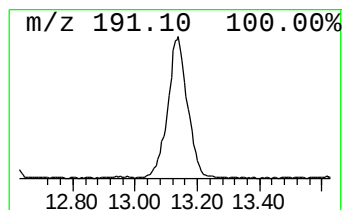
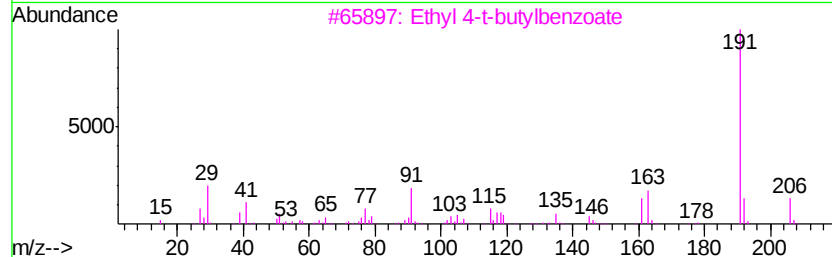
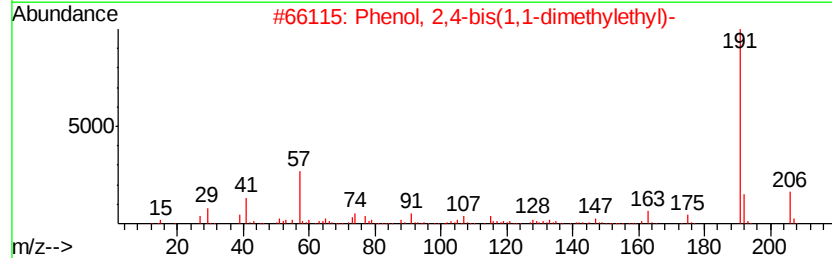
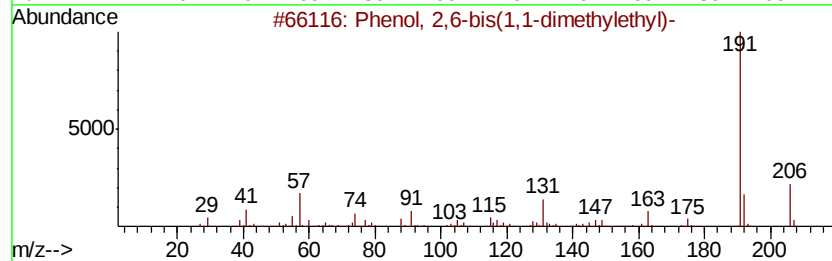
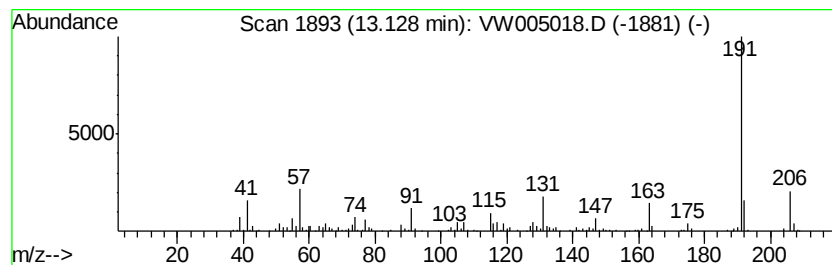
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Quant Title : SW846 8260

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

Peak Number 4 Phenol, 2,6-bis(1,1-dimethy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	80.13 ug/l	313661	1,4-Dichlorobenzene-d4	13.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	96	
2	Phenol, 2,4-bis(1,1-dimethylethyl)-	206	C14H22O	000096-76-4	87	
3	Ethyl 4-t-butylbenzoate	206	C13H18O2	005406-57-5	81	
4	Phenol, 2,5-bis(1,1-dimethylethyl)-	206	C14H22O	005875-45-6	80	
5	Phenol, 3,5-bis(1,1-dimethylethyl)-	206	C14H22O	001138-52-9	74	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2,6-bis(1...	13.13	80.1	ug/l	313661	4	13.57	195732	50.0