

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111919\
 Data File : VD064285.D
 Acq On : 19 Nov 2019 13:01
 Operator : VA/SY
 Sample : K5905-01
 Misc : 10.80G/5.00ml/MSVOA D/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 RT-3037

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	2.208	43	49	50	rBV2	15052	18404	1.32%	0.219%
2	7.920	1008	1020	1025	rBV2	208658	509366	36.43%	6.062%
3	7.985	1025	1031	1040	rVB	372135	832054	59.50%	9.902%
4	8.338	1080	1091	1104	rBV	213455	484629	34.66%	5.767%
5	8.867	1173	1181	1193	rBV	512557	1039382	74.33%	12.369%
6	10.343	1424	1432	1440	rBV	774796	1398321	100.00%	16.641%
7	10.402	1440	1442	1450	rVB4	9370	15918	1.14%	0.189%
8	10.732	1492	1498	1502	rBV3	16417	28604	2.05%	0.340%
9	11.649	1646	1654	1664	rBV	703204	1249842	89.38%	14.874%
10	11.855	1686	1689	1696	rVB7	10071	19861	1.42%	0.236%
11	12.637	1815	1822	1833	rVB	598147	1002075	71.66%	11.925%
12	12.967	1872	1878	1881	rBV7	9195	16426	1.17%	0.195%
13	13.231	1906	1923	1940	rBV3	105219	524119	37.48%	6.237%
14	13.584	1976	1983	1992	rVB	777902	1264055	90.40%	15.043%

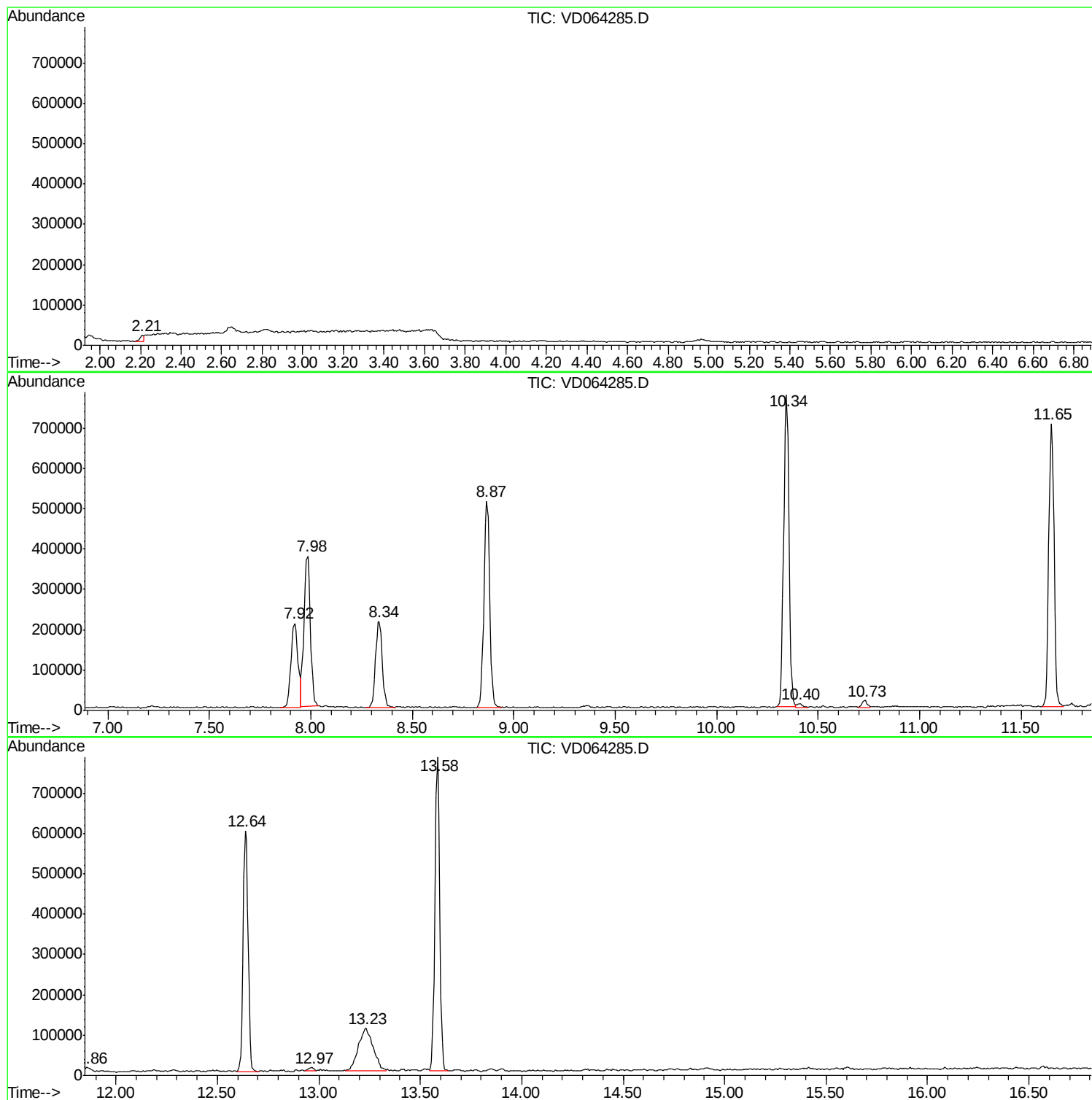
Sum of corrected areas: 8403056

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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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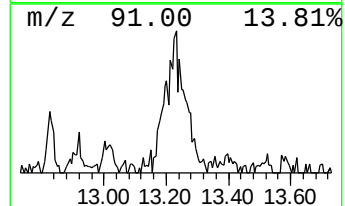
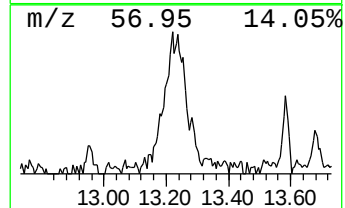
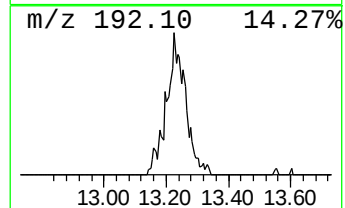
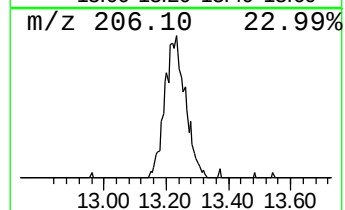
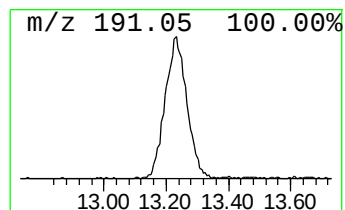
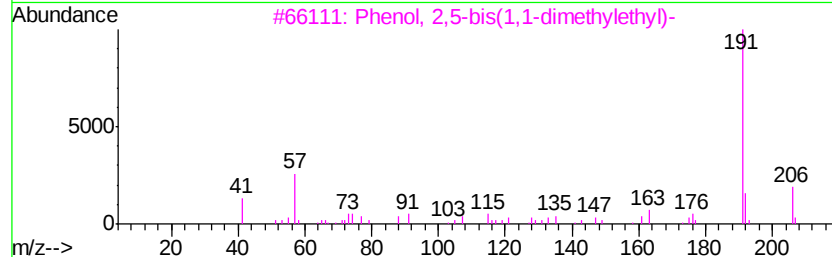
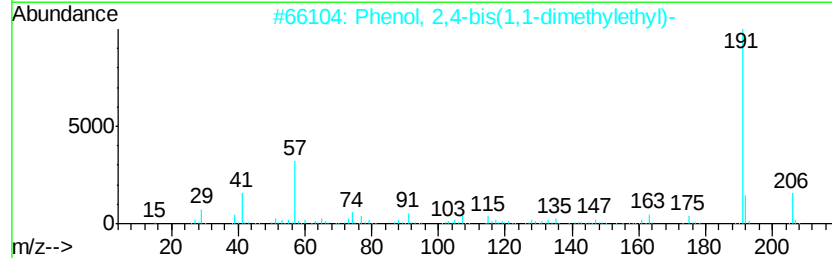
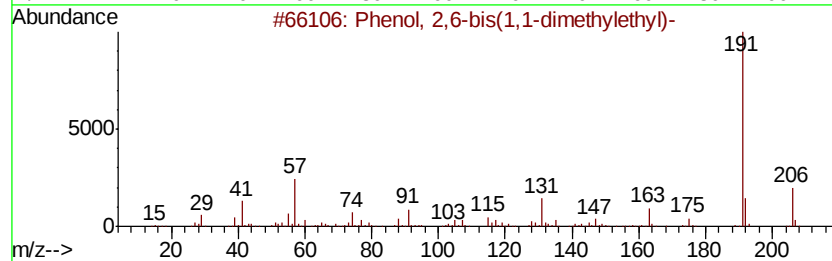
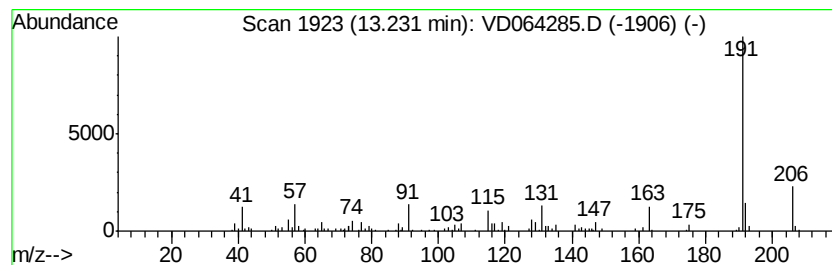
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.23	20.73 ug/l	524119	1,4-Dichlorobenzene-d4	13.58		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol,	2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	93
2	Phenol,	2,4-bis(1,1-dimethylethyl)-	206	C14H22O	000096-76-4	87
3	Phenol,	2,5-bis(1,1-dimethylethyl)-	206	C14H22O	005875-45-6	81
4	Phenol,	3,5-bis(1,1-dimethylethyl)-	206	C14H22O	001138-52-9	81
5	1-(3,6,6-Trimethyl-1,6,7,7a-tetr...		206	C13H18O2	1000194-97-2	81



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2,6-bis(1...	13.23	20.7	ug/l	524119	4	13.58	1264060	50.0