

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN030520\
 Data File : VN060363.D
 Acq On : 5 Mar 2020 13:44
 Operator : JC/MD
 Sample : L1803-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 N38129

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_N\METHODS\82N022720W.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	7.564	1784	1801	1813	rBV	119726	298882	16.91%	4.084%
2	7.641	1813	1825	1848	rVB	240368	566120	32.02%	7.735%
3	8.005	1919	1938	1958	rBV2	143758	338525	19.15%	4.625%
4	8.567	2094	2113	2130	rBV	360662	748801	42.35%	10.231%
5	10.079	2568	2583	2596	rBV	564930	1045431	59.13%	14.284%
6	10.143	2596	2603	2616	rVB	21687	38508	2.18%	0.526%
7	11.400	2980	2994	3017	rBV	520265	919530	52.01%	12.563%
8	12.165	3181	3232	3235	rBV2	573328	1767980	100.00%	24.156%
9	12.397	3291	3304	3320	rVB	433400	723519	40.92%	9.885%
10	13.339	3584	3597	3612	rBV	536647	871771	49.31%	11.911%

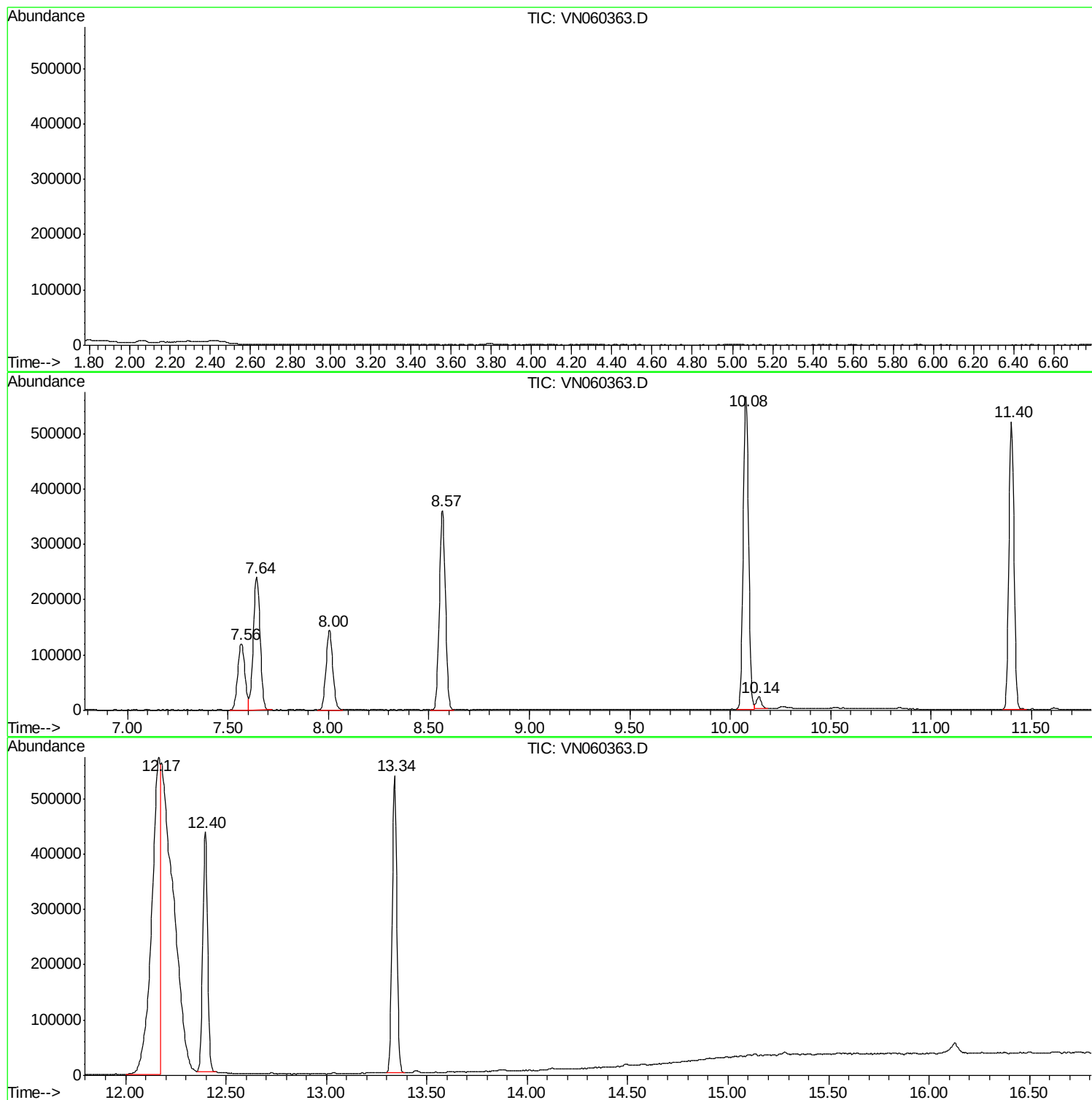
Sum of corrected areas: 7319067

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N022720W.M
Quant Title : SW846 8260

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



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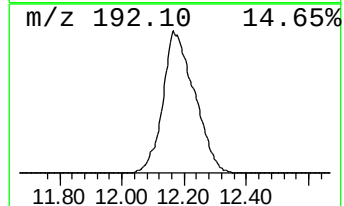
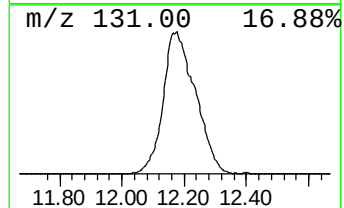
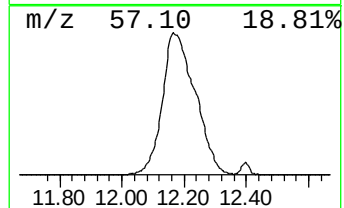
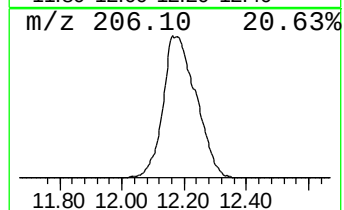
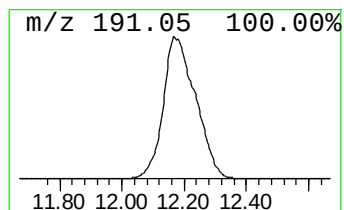
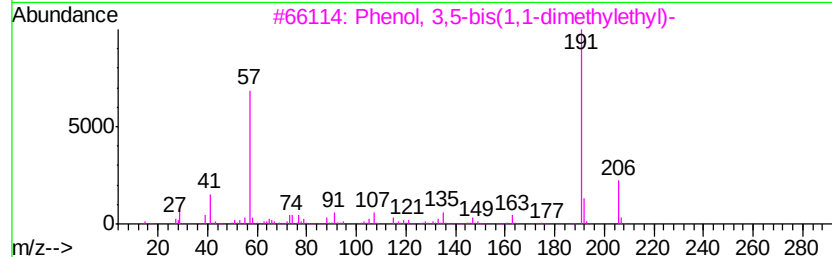
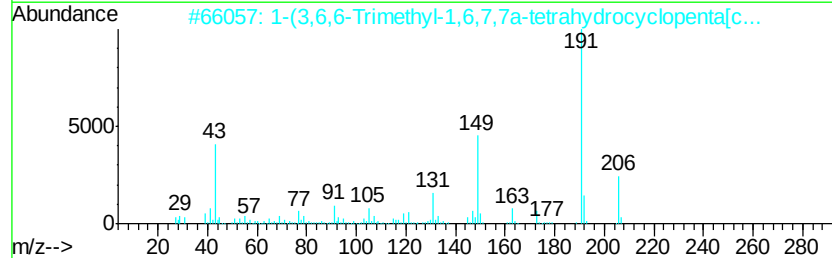
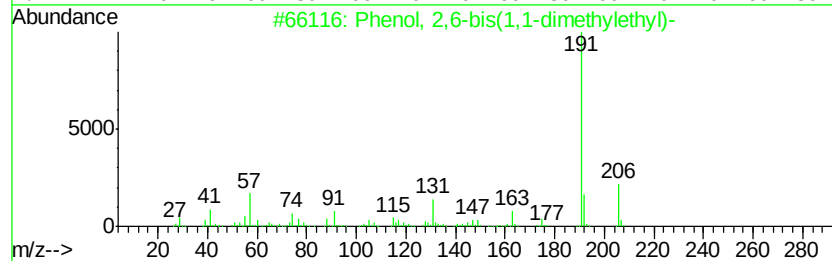
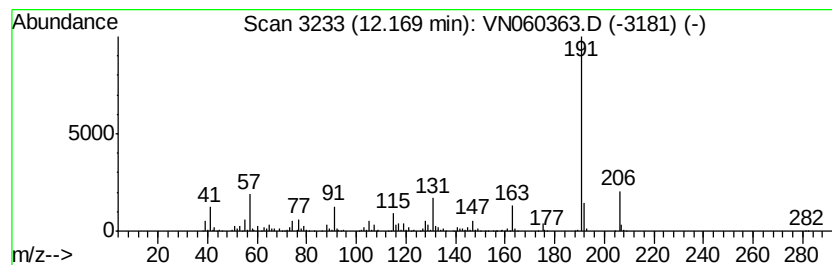
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.17	96.14 ug/l	1767980	Chlorobenzene-d5	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	97
2		1-(3,6,6-Trimethyl-1,6,7,7a-tetr...	206	C13H18O2	1000194-97-2	80
3		Phenol, 3,5-bis(1,1-dimethylethyl)-	206	C14H22O	001138-52-9	80
4		2H-Indeno[1,2-b]furan-2-one, 3,3...	206	C13H18O2	1000196-74-3	80
5		Phenol, 2,4-bis(1,1-dimethylethyl)-	206	C14H22O	000096-76-4	78



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2,6-bis(1...	12.17	96.1	ug/l	1767980	3	11.40	919530	50.0