

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN082620\
 Data File : VN063134.D
 Acq On : 26 Aug 2020 15:47
 Operator : JC/MD
 Sample : L3774-09
 Misc : 1.00g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TOTE-1-2

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\82N081820W.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	2.445	190	209	210	rBV6	12798	24939	1.72%	0.382%
2	7.548	1777	1796	1807	rBV	104622	264151	18.20%	4.046%
3	7.622	1807	1819	1843	rVB2	176314	420167	28.95%	6.435%
4	7.985	1915	1932	1952	rBV	112226	259887	17.91%	3.980%
5	8.548	2089	2107	2126	rBV	284309	588079	40.52%	9.007%
6	10.059	2563	2577	2590	rBV	444289	820042	56.51%	12.559%
7	10.124	2590	2597	2611	rVB	28943	52999	3.65%	0.812%
8	11.381	2974	2988	3011	rBV	424803	738751	50.91%	11.314%
9	12.082	3159	3206	3246	rBV5	85033	600392	41.37%	9.195%
10	12.378	3286	3298	3316	rBV	339777	573880	39.54%	8.789%
11	13.320	3577	3591	3612	rBV	386498	649084	44.73%	9.941%
12	13.931	3752	3781	3808	rBV2	397528	1451216	100.00%	22.226%
13	15.210	4171	4179	4191	rVB9	9580	20206	1.39%	0.309%
14	15.419	4231	4244	4254	rBV3	10808	28921	1.99%	0.443%
15	15.856	4370	4380	4397	rVB	17618	36752	2.53%	0.563%

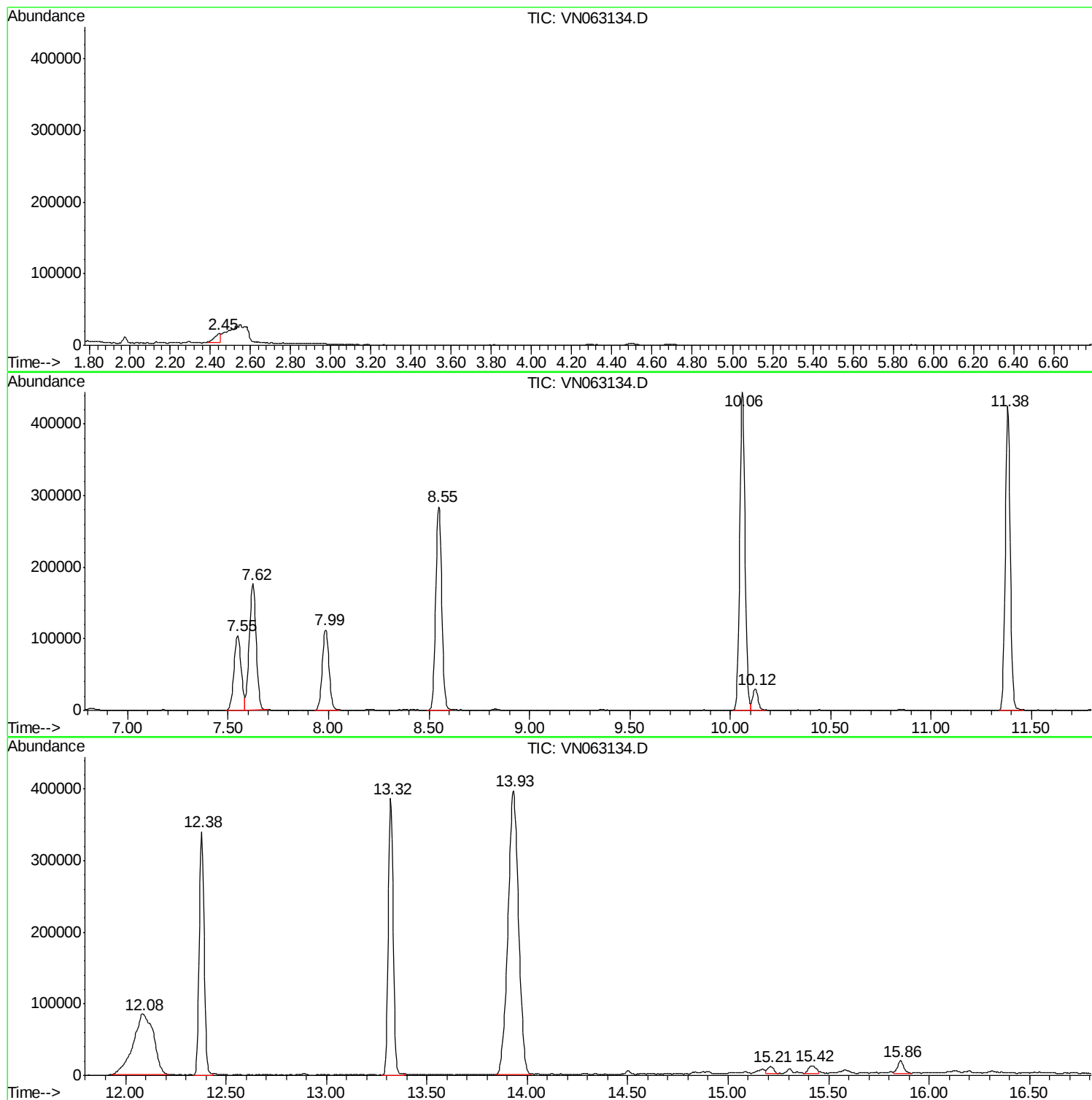
Sum of corrected areas: 6529466

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



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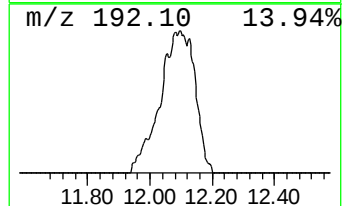
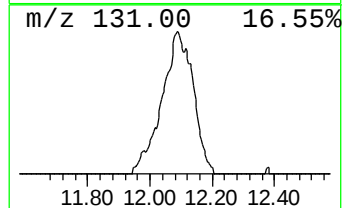
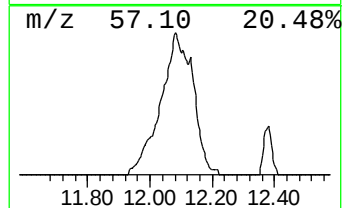
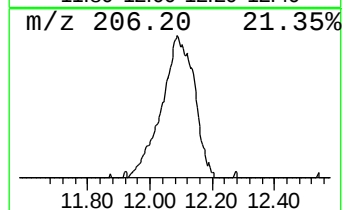
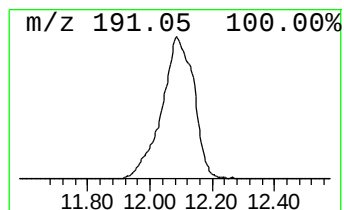
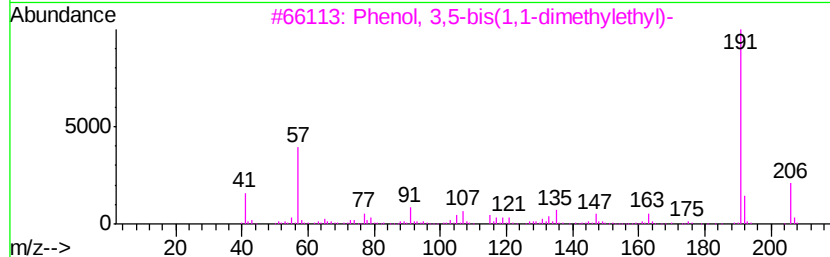
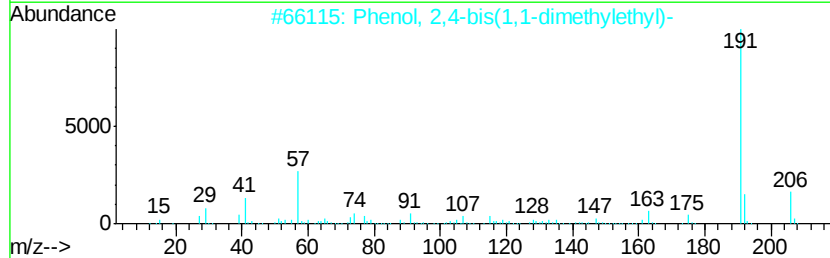
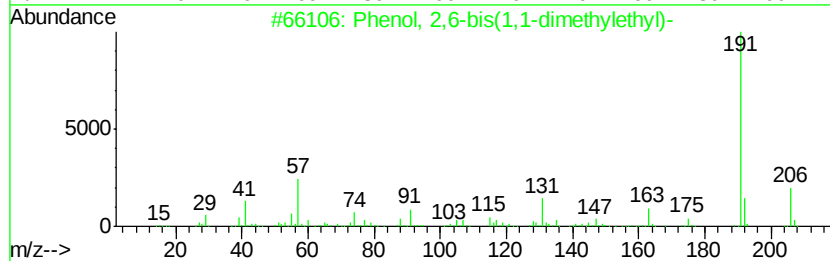
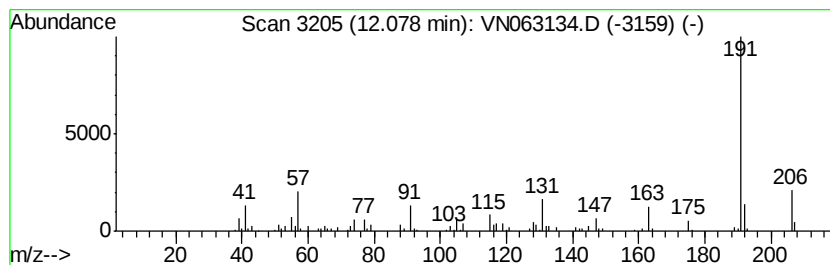
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Quant Title : SW846 8260

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

Peak Number 1 Phenol, 2,6-bis(1,1-dimethylethyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.08	40.64 ug/l	600392	Chlorobenzene-d5	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	90
3		Phenol, 3,5-bis(1,1-dimethylethyl)-	206	C14H22O	001138-52-9	81
4		Ethyl 4-t-butylbenzoate	206	C13H18O2	005406-57-5	81
5		Phenol, 2,5-bis(1,1-dimethylethyl)-	206	C14H22O	005875-45-6	74



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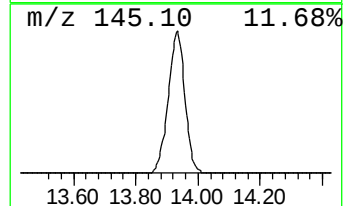
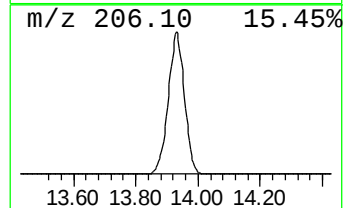
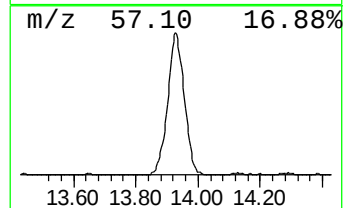
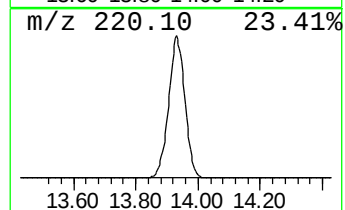
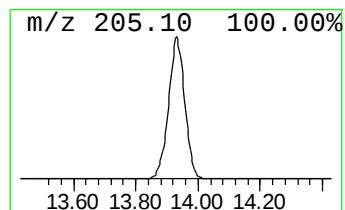
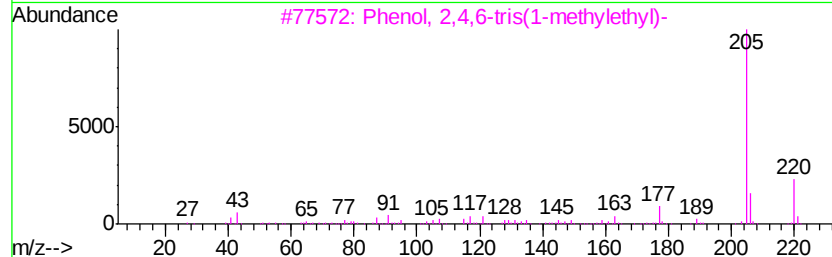
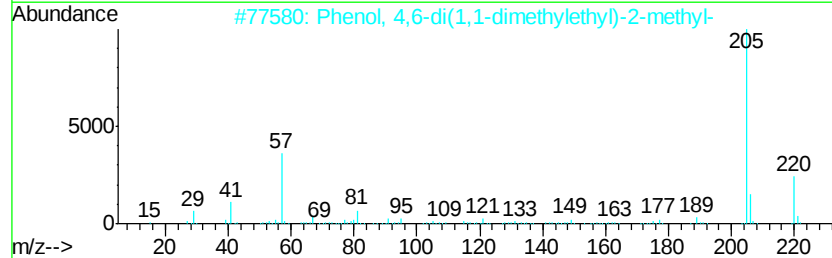
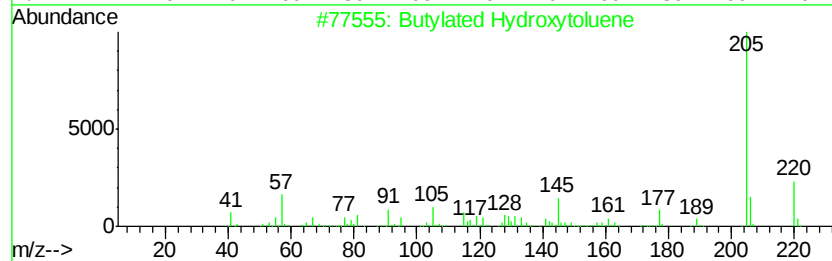
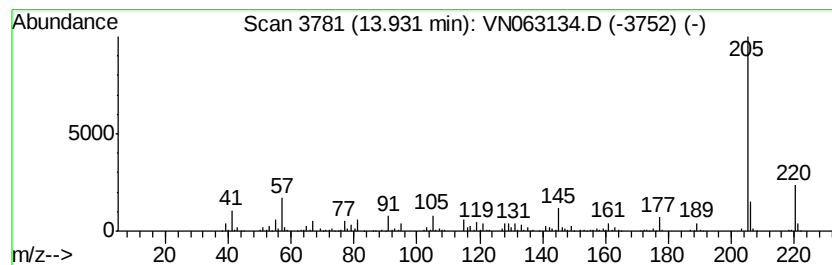
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Peak Number 2 Butylated Hydroxytoluene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.93	111.79 ug/l	1451220	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	99
2		Phenol, 4,6-di(1,1-dimethylethyl)-	220	C15H24O	000616-55-7	93
3		Phenol, 2,4,6-tris(1-methylethyl)-	220	C15H24O	002934-07-8	87
4		Propanoic acid, 2-methyl-3-[4-t-...	220	C14H20O2	1000131-87-0	72
5		2,4,6-Tris(1,1-dimethylethyl)-4-...	276	C19H32O	019687-22-0	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2,6-bis(1...	12.08	40.6	ug/l	600392	3	11.38	738751	50.0
Butylated Hydroxy...	13.93	111.8	ug/l	1451220	4	13.32	649084	50.0