

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090420\
 Data File : VN063248.D
 Acq On : 4 Sep 2020 15:25
 Operator : JC/MD
 Sample : L3902-18
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SA-PZ172I-20200902

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_N\METHODS\82N083120W.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	2.033	83	91	98	rBV5	10852	19371	1.04%	0.183%
2	2.422	200	212	222	rBV5	57861	126234	6.75%	1.192%
3	4.020	695	709	722	rBV2	14218	39672	2.12%	0.375%
4	4.512	844	862	865	rBV2	8137	19472	1.04%	0.184%
5	5.801	1244	1263	1281	rBV3	31813	102784	5.49%	0.971%
6	7.550	1789	1807	1818	rBV2	225282	566766	30.29%	5.352%
7	7.627	1819	1831	1849	rVB2	361122	863117	46.12%	8.150%
8	7.987	1927	1943	1966	rBV2	260756	630365	33.69%	5.952%
9	8.550	2103	2118	2135	rBV	615067	1280624	68.44%	12.093%
10	10.061	2573	2588	2613	rBV	1007655	1871295	100.00%	17.670%
11	11.383	2984	2999	3021	rBV	963459	1716299	91.72%	16.207%
12	12.379	3296	3309	3330	rBV	706868	1230694	65.77%	11.621%
13	13.318	3587	3601	3618	rBV	844670	1447258	77.34%	13.666%
14	13.942	3759	3795	3816	rBV7	156705	676078	36.13%	6.384%

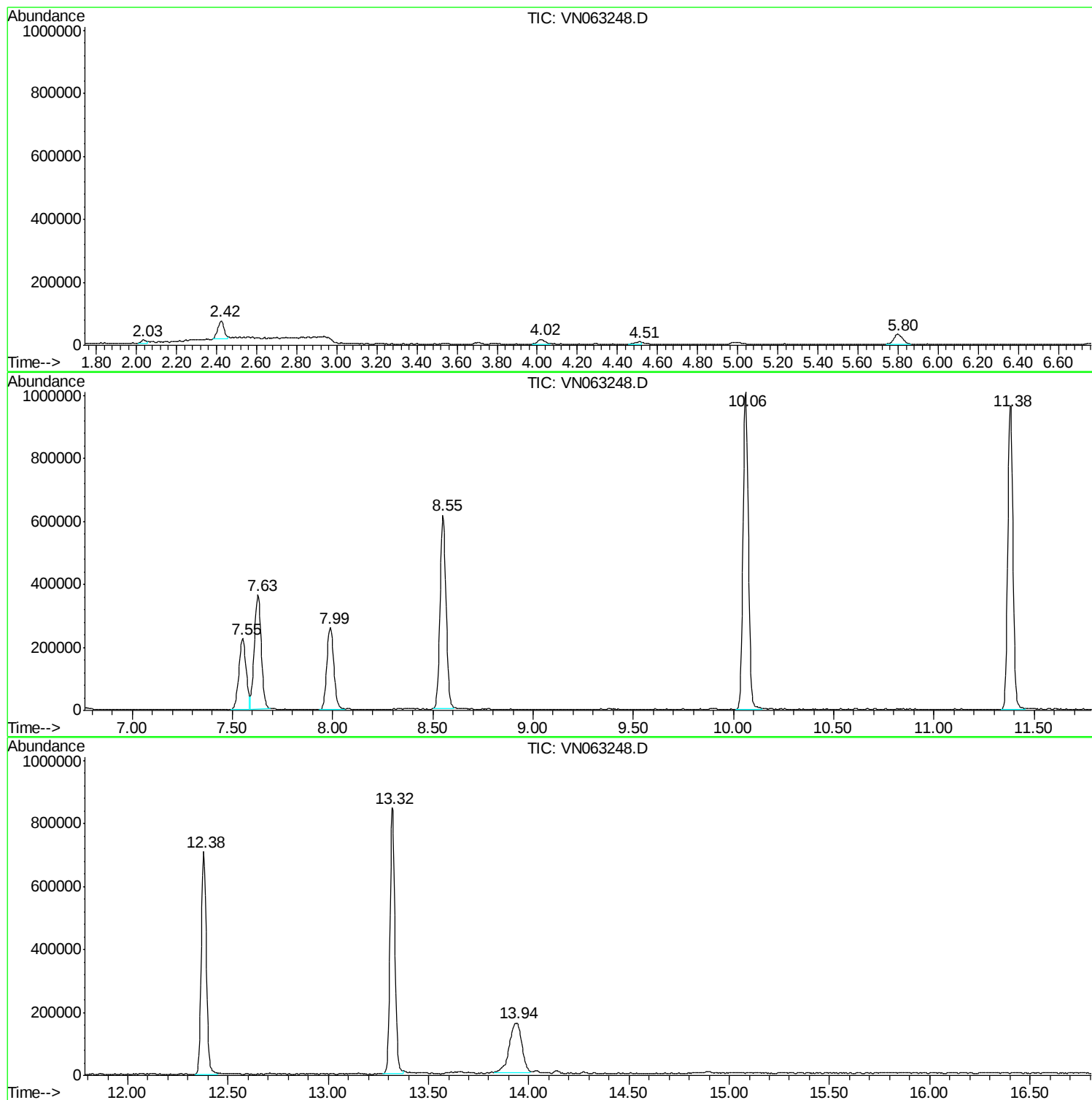
Sum of corrected areas: 10590029

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Quant Title : SW846 8260

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



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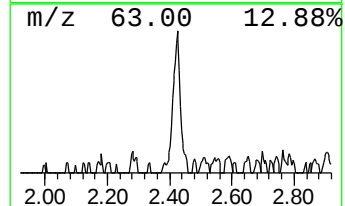
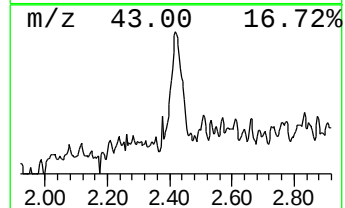
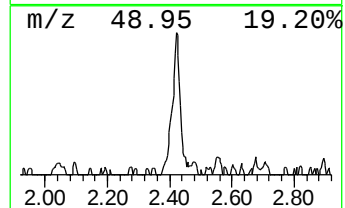
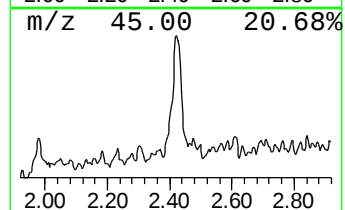
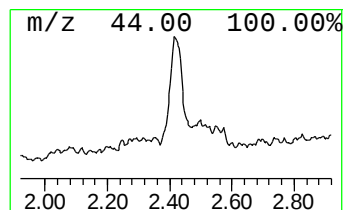
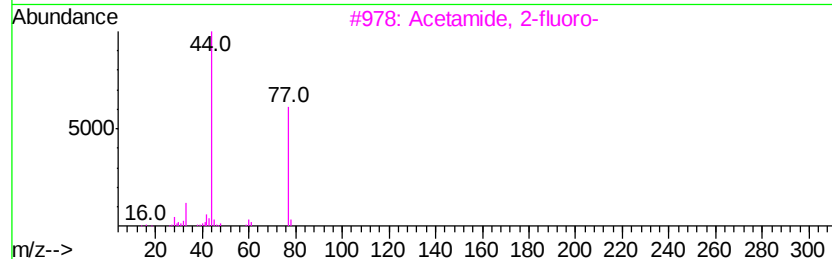
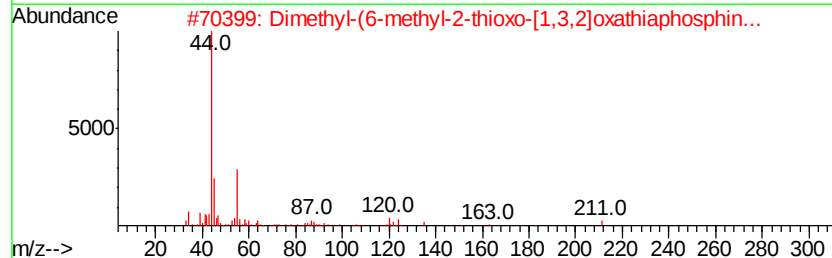
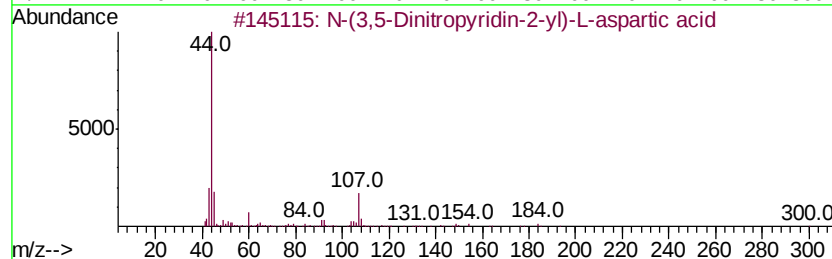
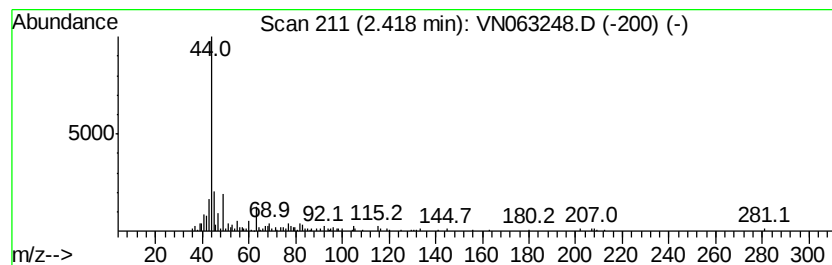
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Peak Number 1 unknown2.42 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.42	7.31 ug/l	126234	Pentafluorobenzene	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N-(3,5-Dinitropyridin-2-yl)-L-as...	300	C9H8N4O8	035899-60-6	23
2		Dimethyl-(6-methyl-2-thioxo-[1,3...	211	C6H14NOPS2	139575-19-2	23
3		Acetamide, 2-fluoro-	77	C2H4FNO	000640-19-7	22
4		2-Pentanamine	87	C5H13N	063493-28-7	22
5		Ala-gly, trimethylsilyl ester	218	C8H18N2O3Si	1000333-70-1	16



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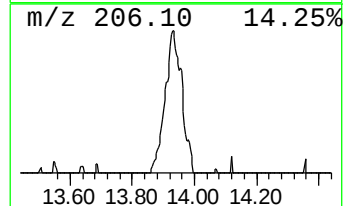
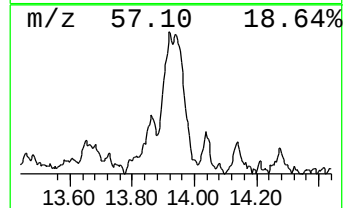
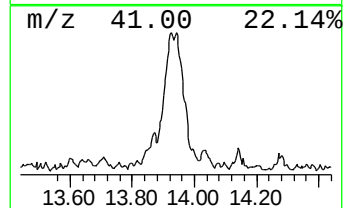
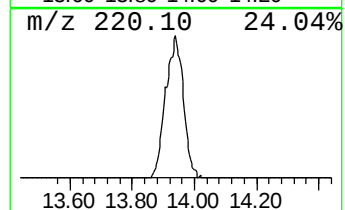
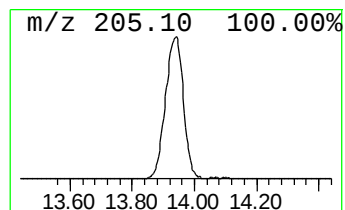
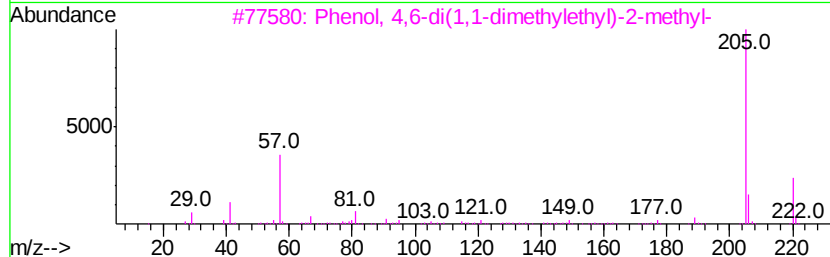
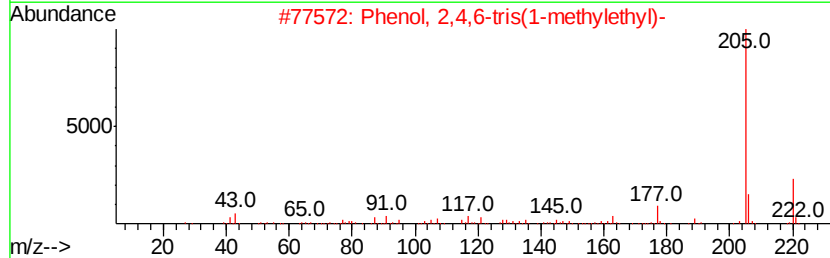
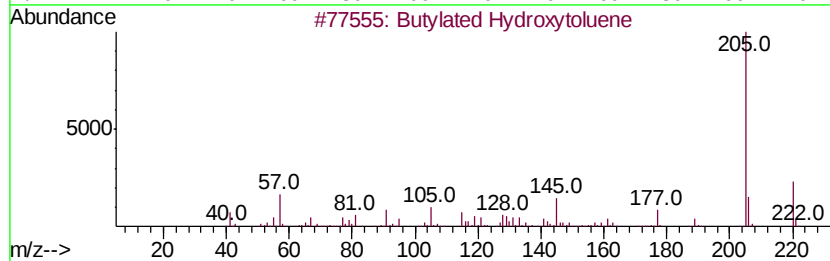
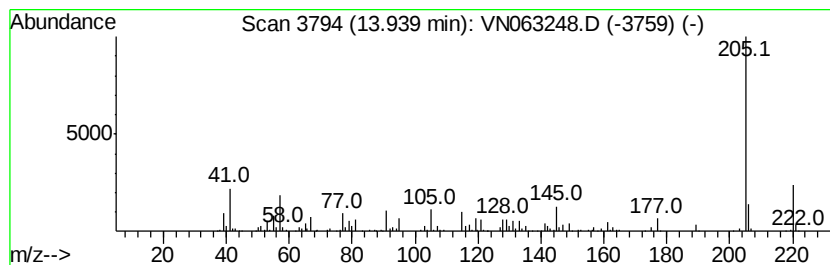
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Peak Number 2 Butylated Hydroxytoluene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.94	23.36 ug/l	676078	1,4-Dichlorobenzene-d4	13.32		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butylated Hydroxytoluene	220	C15H24O	000128-37-0	98
2		Phenol, 2,4,6-tris(1-methylethyl)-	220	C15H24O	002934-07-8	83
3		Phenol, 4,6-di(1,1-dimethylethyl)-	220	C15H24O	000616-55-7	83
4		1-Butene, 3-phenyl-1-(trimethylsilyl)-	220	C13H20OSi	055543-95-8	59
5		Phenol, 2,6-bis(1,1-dimethylethyl)-	277	C17H27NO2	001918-11-2	58



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.42	2.42	7.3	ug/l	126234	1	7.63	863117	50.0
Butylated Hydroxy...	13.94	23.4	ug/l	676078	4	13.32	1447260	50.0