

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN072020\
 Data File : VN062558.D
 Acq On : 20 Jul 2020 13:41
 Operator : JC/MD
 Sample : L3331-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 001-WILLETS-PT-BLVD-(JULY)

Integration Parameters: RTEINT3.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0 % 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\624N071320W.M
 Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.777	598	623	670	rBV	940733	2491314	100.00%	34.731%
2	4.021	681	699	735	rBV	315089	891369	35.78%	12.426%
3	6.796	1549	1562	1587	rBV2	9413	25541	1.03%	0.356%
4	7.153	1654	1673	1696	rBV2	106111	292880	11.76%	4.083%
5	7.989	1913	1933	1956	rBV	112070	264998	10.64%	3.694%
6	8.551	2093	2108	2129	rBV	235293	492458	19.77%	6.865%
7	9.956	2532	2545	2555	rBV3	14393	27174	1.09%	0.379%
8	10.059	2562	2577	2588	rBV	373434	687058	27.58%	9.578%
9	10.124	2588	2597	2620	rVV	356086	667263	26.78%	9.302%
10	11.381	2974	2988	3006	rBV	329876	583671	23.43%	8.137%
11	12.111	3203	3215	3229	rBV5	14941	27536	1.11%	0.384%
12	12.378	3286	3298	3318	rBV	228879	388165	15.58%	5.411%
13	12.918	3457	3466	3476	rBV3	15813	26795	1.08%	0.374%
14	12.992	3477	3489	3503	rVB3	13574	28311	1.14%	0.395%
15	13.114	3507	3527	3547	rBV3	39291	105890	4.25%	1.476%
16	13.265	3564	3574	3583	rBV2	24816	40876	1.64%	0.570%
17	13.378	3597	3609	3618	rVV7	20816	51332	2.06%	0.716%
18	14.622	3985	3996	4008	rBV2	48352	80588	3.23%	1.123%

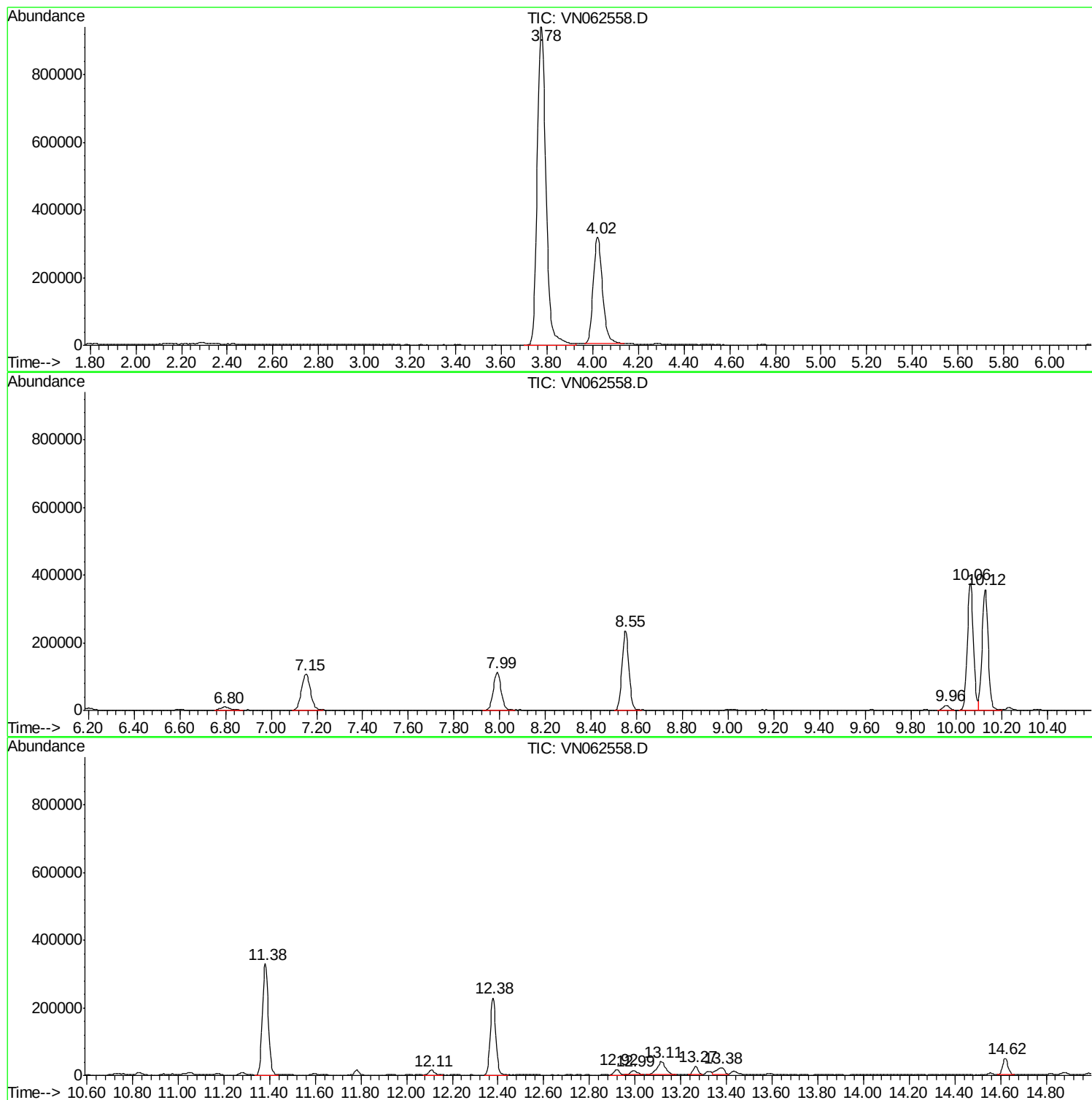
Sum of corrected areas: 7173219

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA N\METHODS\624N071320W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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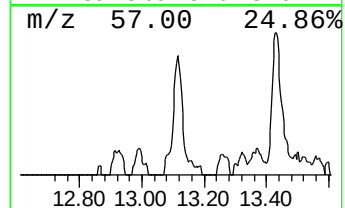
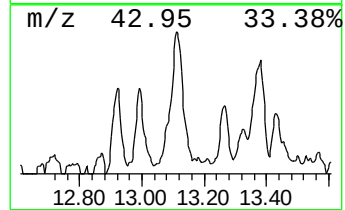
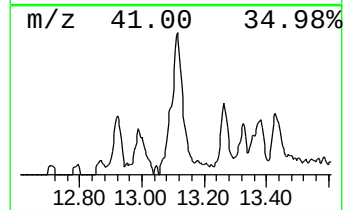
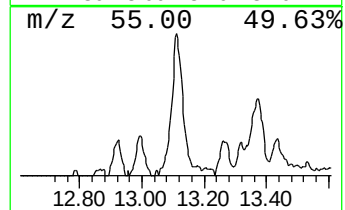
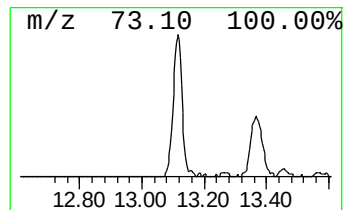
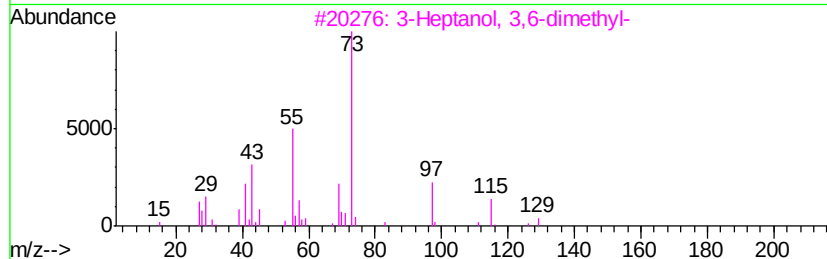
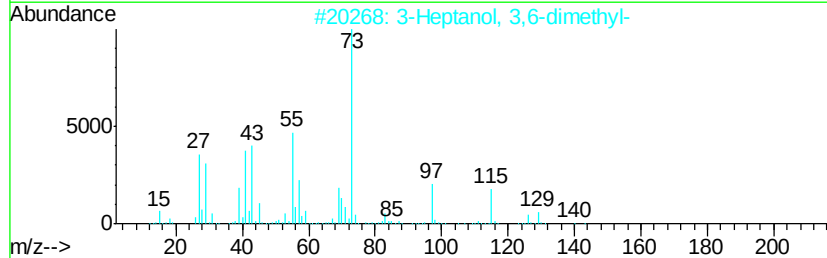
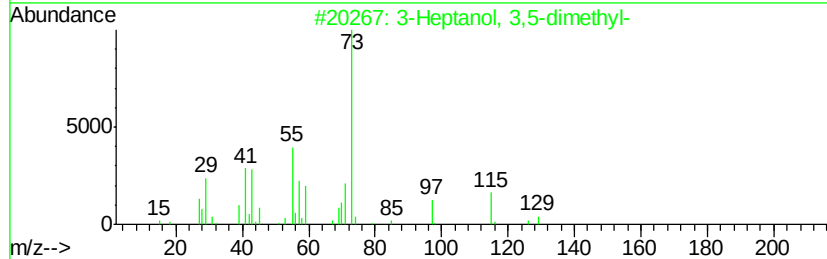
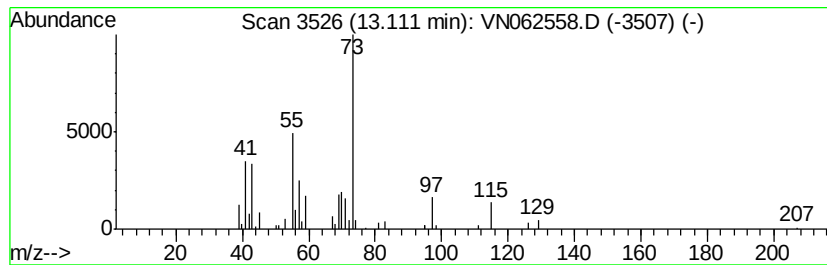
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Heptanol, 3,5-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.11	5.44 ug/l	105890	Chlorobenzene-d5	11.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanol, 3,5-dimethyl-	144	C9H20O	019549-74-7	78
2	3-Heptanol, 3,6-dimethyl-	144	C9H20O	001573-28-0	59
3	3-Heptanol, 3,6-dimethyl-	144	C9H20O	001573-28-0	53
4	Sodium N-methyl dithiocarbamate	129	C2H4NNaS2	1000292-95-8	25
5	1,1,1,4-Tetramethyl-4-chloro-4-v...	206	C8H19ClSi2	102105-49-7	25



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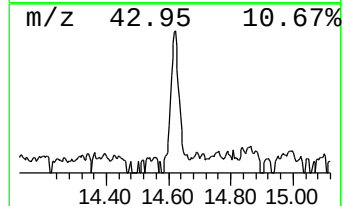
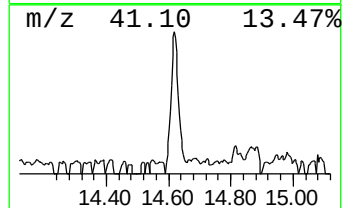
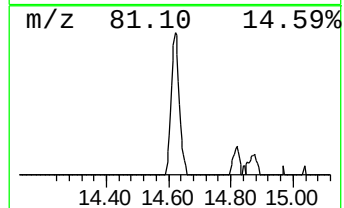
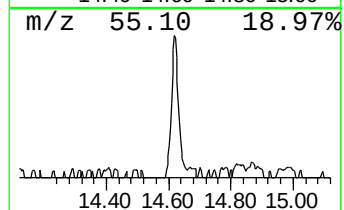
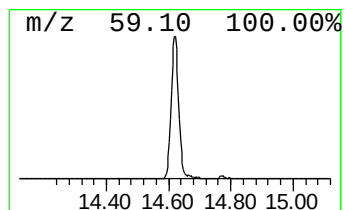
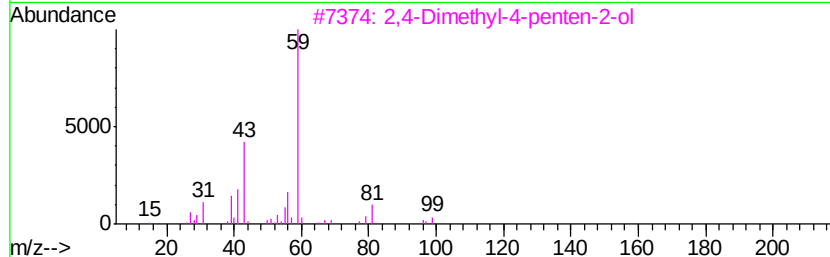
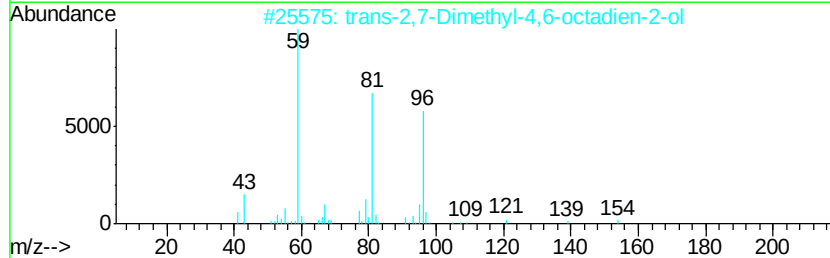
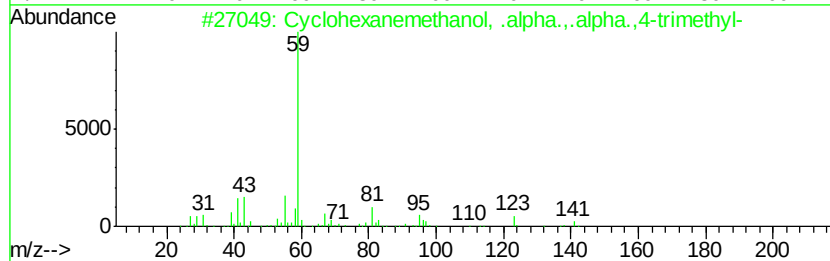
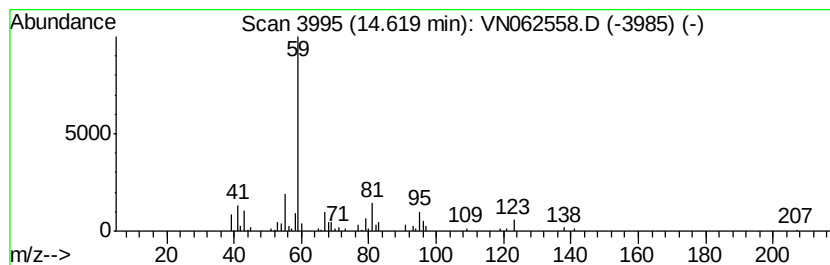
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexanemethanol, .alpha... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.62	4.14 ug/l	80588	Chlorobenzene-d5	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanemethanol, .alpha...al...	156	C10H20O	000498-81-7	83
2		trans-2,7-Dimethyl-4,6-octadien-...	154	C10H18O	1000281-69-5	59
3		2,4-Dimethyl-4-penten-2-ol	114	C7H14O	019781-53-4	53
4		Cyclohexanemethanol, .alpha...al...	156	C10H20O	000498-81-7	50
5		Cyclohexanemethanol, .alpha...al...	156	C10H20O	005114-00-1	50



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Heptanol, 3,5-d...	13.11	5.4	ug/l	105890	3	11.38	583671	30.0
Cyclohexanemethan...	14.62	4.1	ug/l	80588	3	11.38	583671	30.0