

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061622\
 Data File : VN072896.D
 Acq On : 16 Jun 2022 22:27
 Operator : JC\MD
 Sample : N3202-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-22

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\82N061622W.M

Title : SW846 8260

Signal : TIC: VN072896.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.292	564	579	604	rBV	455381	1664082	29.43%	6.292%
2	5.534	604	620	646	rVV	1593200	5653701	100.00%	21.375%
3	7.951	1020	1031	1036	rBV	565057	1360336	24.06%	5.143%
4	8.016	1036	1042	1053	rVB	788558	1820807	32.21%	6.884%
5	8.369	1092	1102	1112	rBV	606865	1395511	24.68%	5.276%
6	8.469	1112	1119	1127	rVV	222421	524164	9.27%	1.982%
7	8.904	1181	1193	1210	rBV	1135597	2393643	42.34%	9.050%
8	10.133	1393	1402	1406	rBV2	101932	205280	3.63%	0.776%
9	10.374	1434	1443	1456	rBV	1985442	3659003	64.72%	13.834%
10	10.486	1456	1462	1472	rVB	162032	296225	5.24%	1.120%
11	11.680	1655	1665	1679	rBV	1616680	2852548	50.45%	10.785%
12	12.668	1825	1833	1843	rBV	1431740	2465250	43.60%	9.321%
13	13.610	1985	1993	2007	rBV	1262217	2158971	38.19%	8.163%

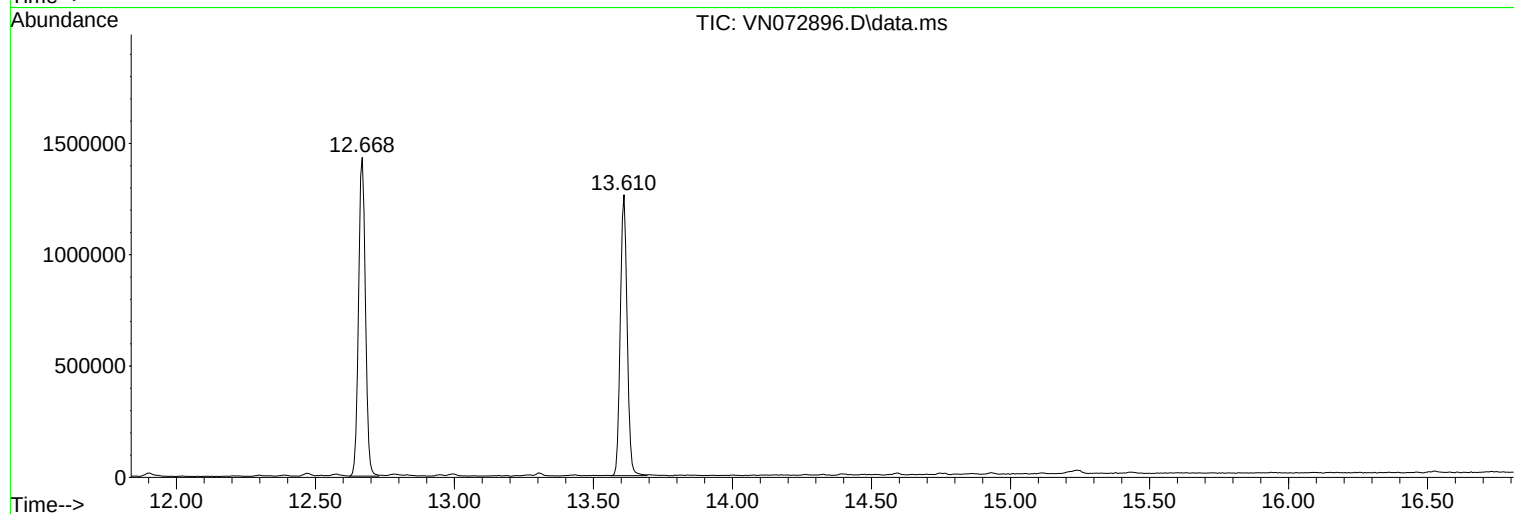
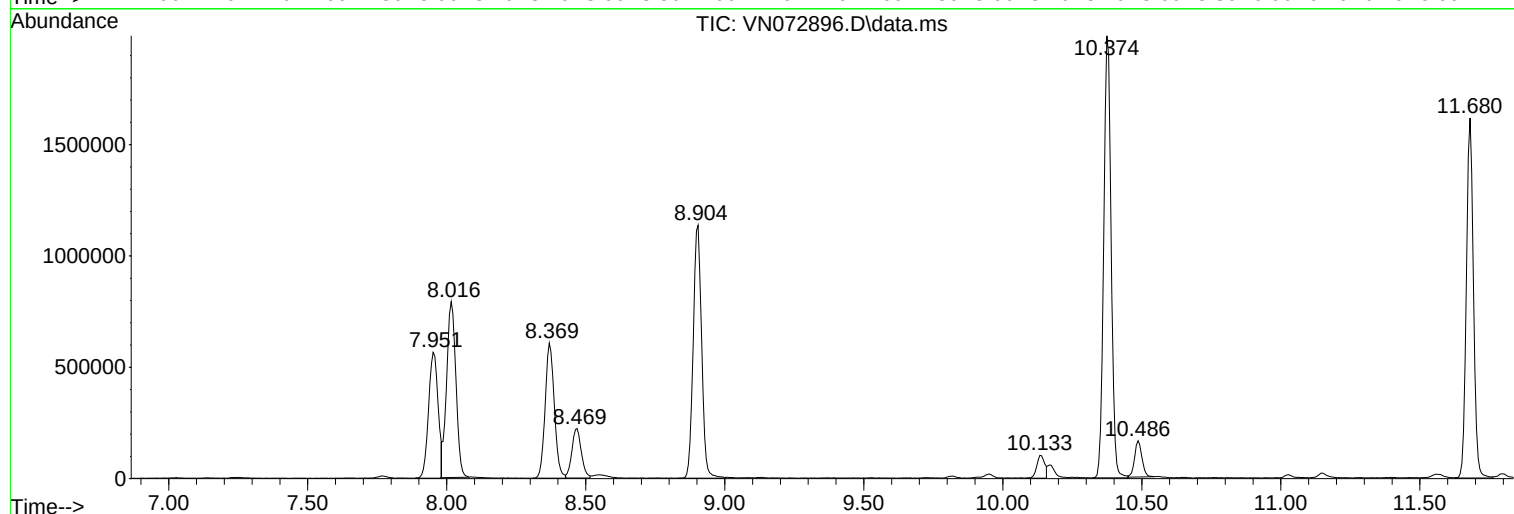
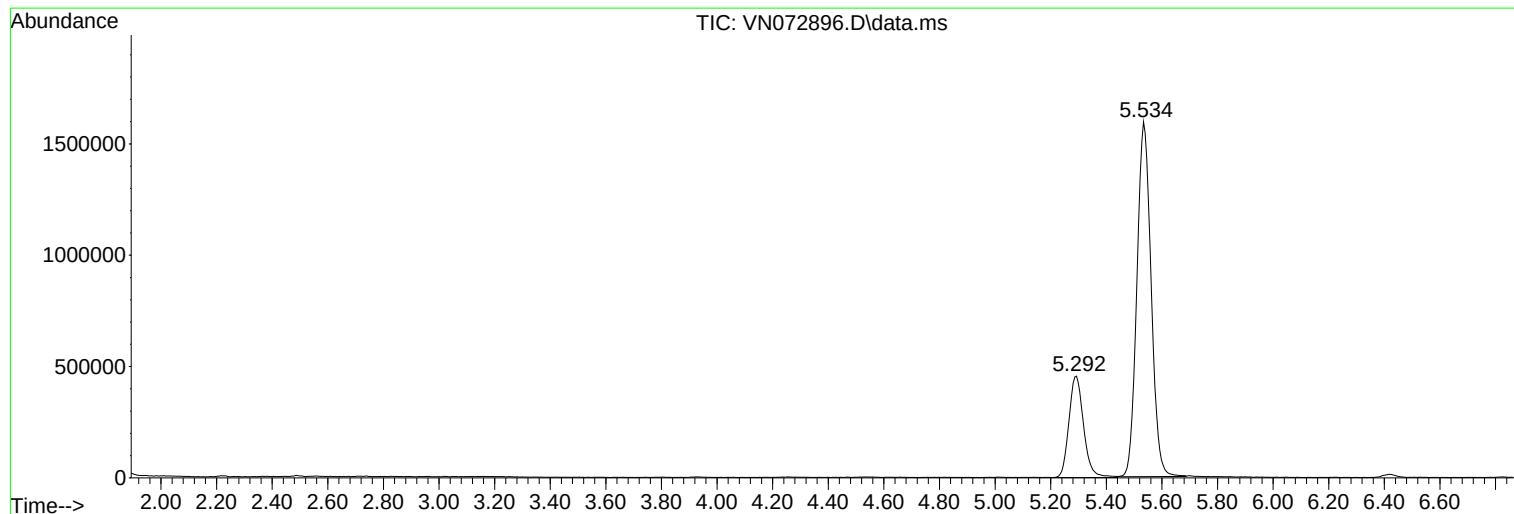
Sum of corrected areas: 26449521

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061622W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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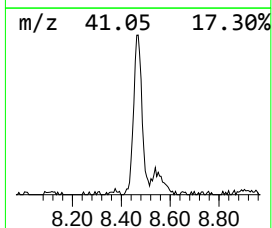
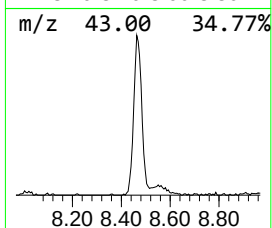
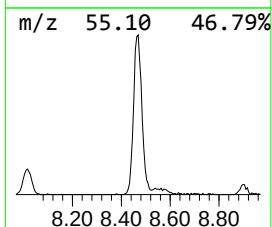
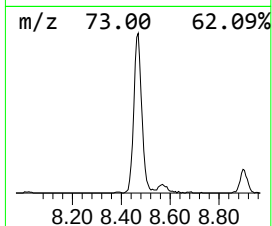
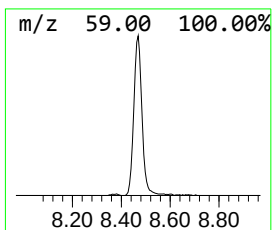
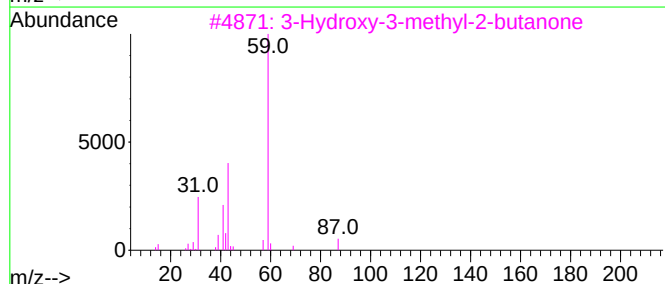
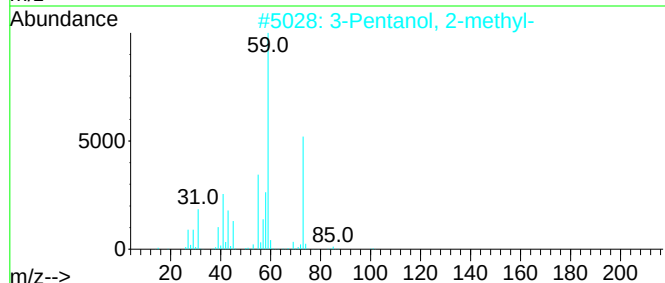
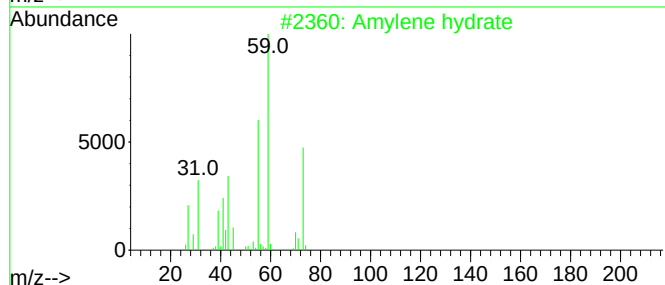
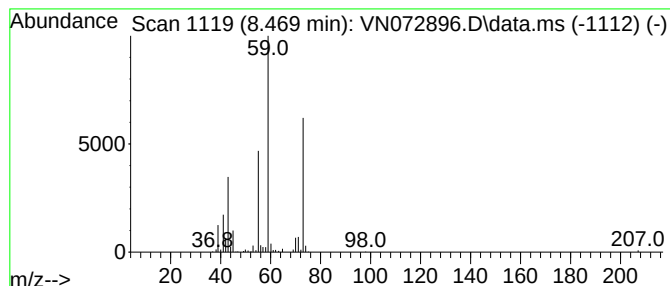
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Amylene hydrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.469	10.95 ug/l	524164	1,4-Difluorobenzene	8.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Amylene hydrate	88	C5H12O	000075-85-4	83
2			3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	50
3			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	40
4			3-Hexanol	102	C6H14O	000623-37-0	39
5			3-Heptanol, 4-methyl-	130	C8H18O	014979-39-6	39



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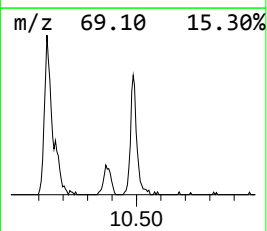
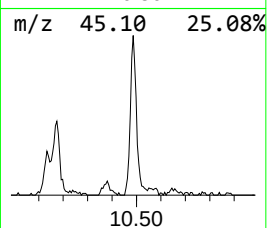
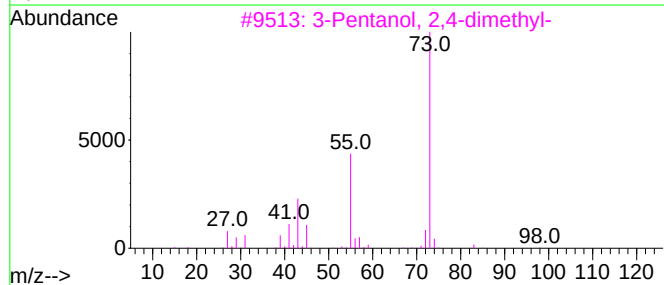
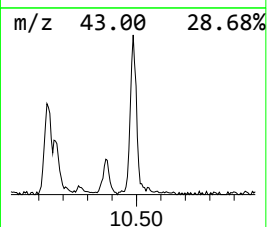
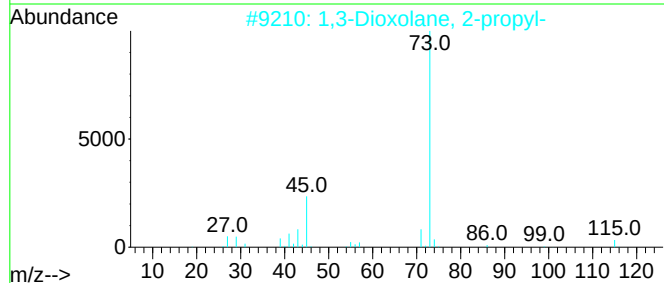
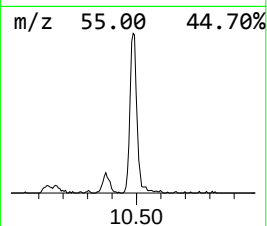
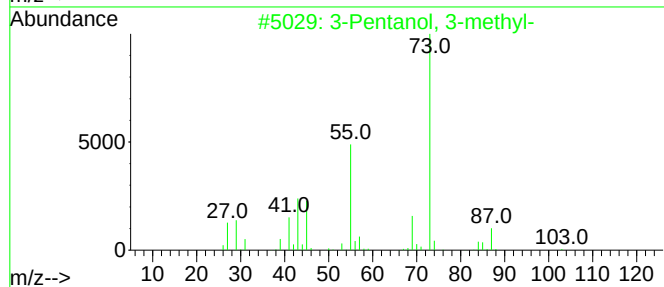
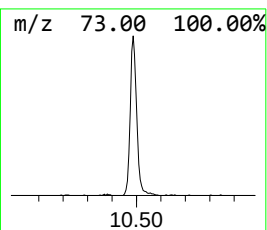
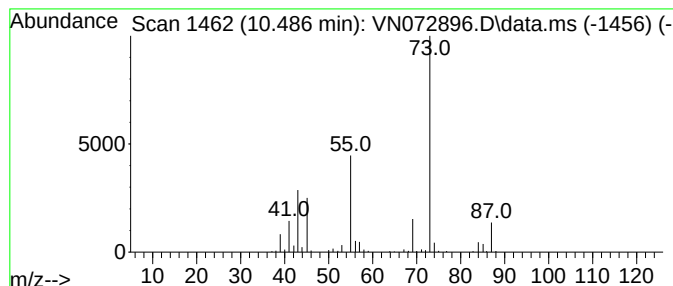
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Peak Number 2 3-Pentanol, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.486	5.19 ug/l	296225	Chlorobenzene-d5	11.680

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	83
2			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	43
3			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4			1,3-Dioxolane, 2-hexyl-	158	C9H18O2	001708-34-5	32
5			3-Methylpentan-3-yl propyl carbo...	188	C10H20O3	1000372-82-3	17



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Amylene hydrate	8.469	10.9	ug/l	524164	2	8.904	2393640	50.0
3-Pentanol, 3-m...	10.486	5.2	ug/l	296225	3	11.680	2852550	50.0