

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072925\
Data File : VX047176.D
Acq On : 29 Jul 2025 14:42
Operator : JC/MD
Sample : Q2719-02 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
50765

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_X\Method\82X072125W.M

Title : SW846 8260

Signal : TIC: VX047176.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.385	693	705	721	rBV2	184852	589633	1.06%	0.897%
2	5.556	721	733	753	rVB	313905	985539	1.78%	1.499%
3	5.958	789	799	822	rBV2	195828	571277	1.03%	0.869%
4	6.763	921	931	942	rBV	502491	1291975	2.33%	1.966%
5	6.873	942	949	971	rVB	279327	740726	1.34%	1.127%
6	8.647	1233	1240	1258	rBV	903409	1473377	2.66%	2.242%
7	10.049	1465	1470	1487	rBV	1091687	1532269	2.77%	2.331%
8	11.079	1633	1639	1647	rBV2	1121691	1565874	2.83%	2.382%
9	12.018	1788	1793	1802	rVB	1283229	1580447	2.85%	2.405%
10	12.774	1907	1917	1928	rBV	40903070	55394876	100.00%	84.282%

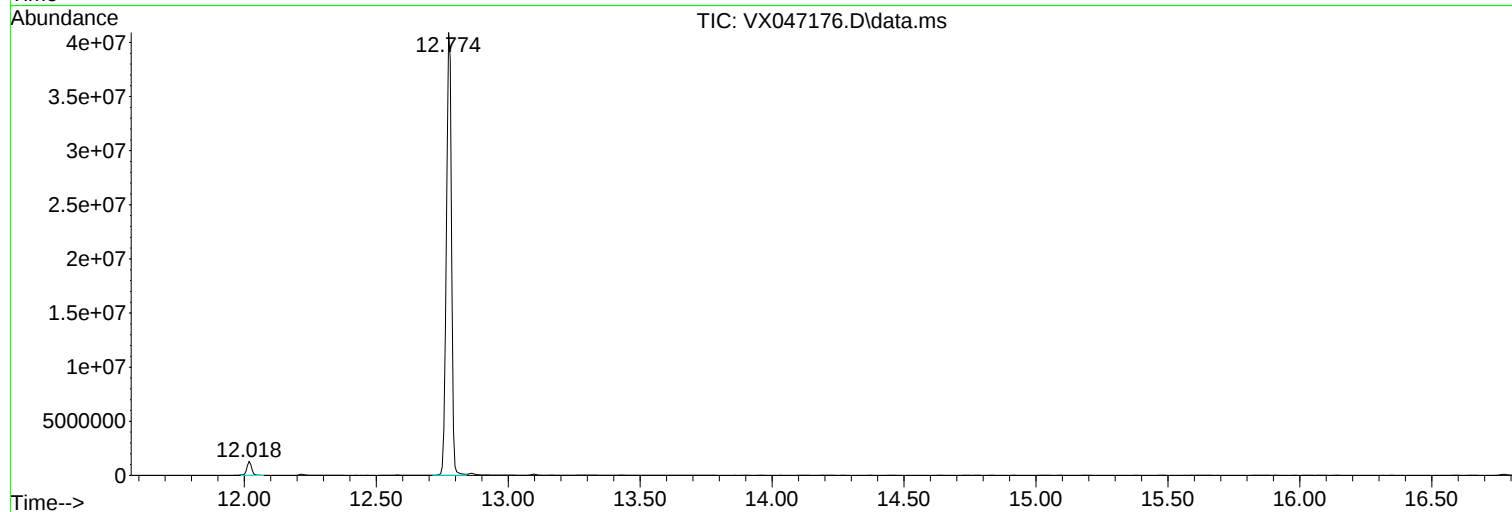
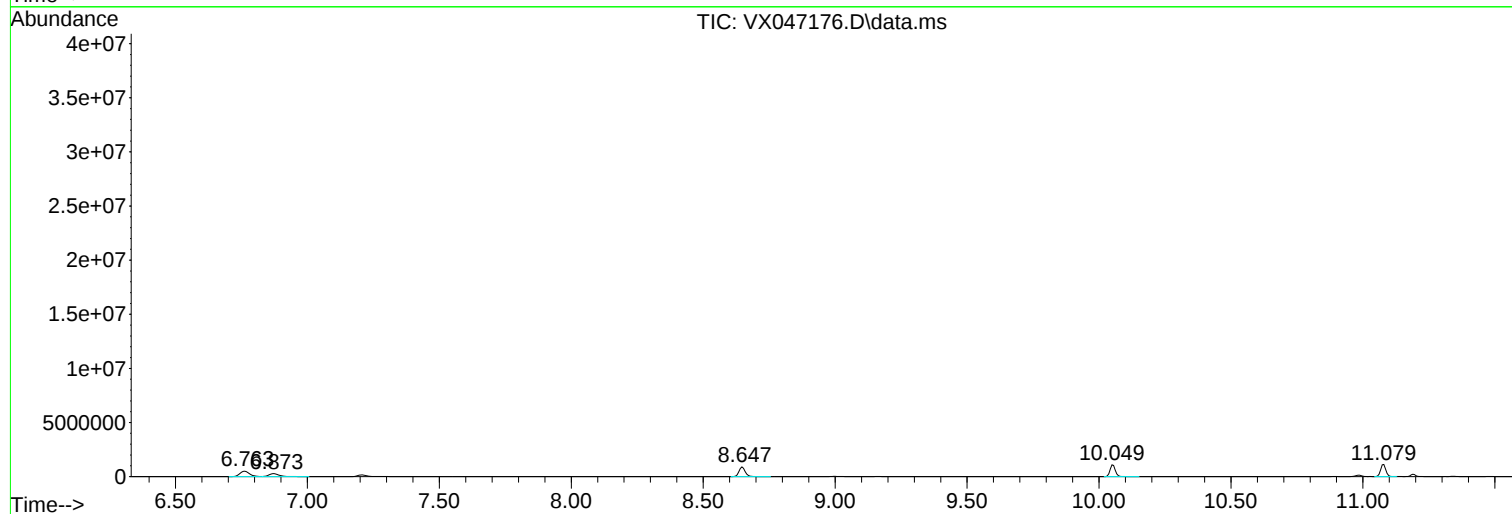
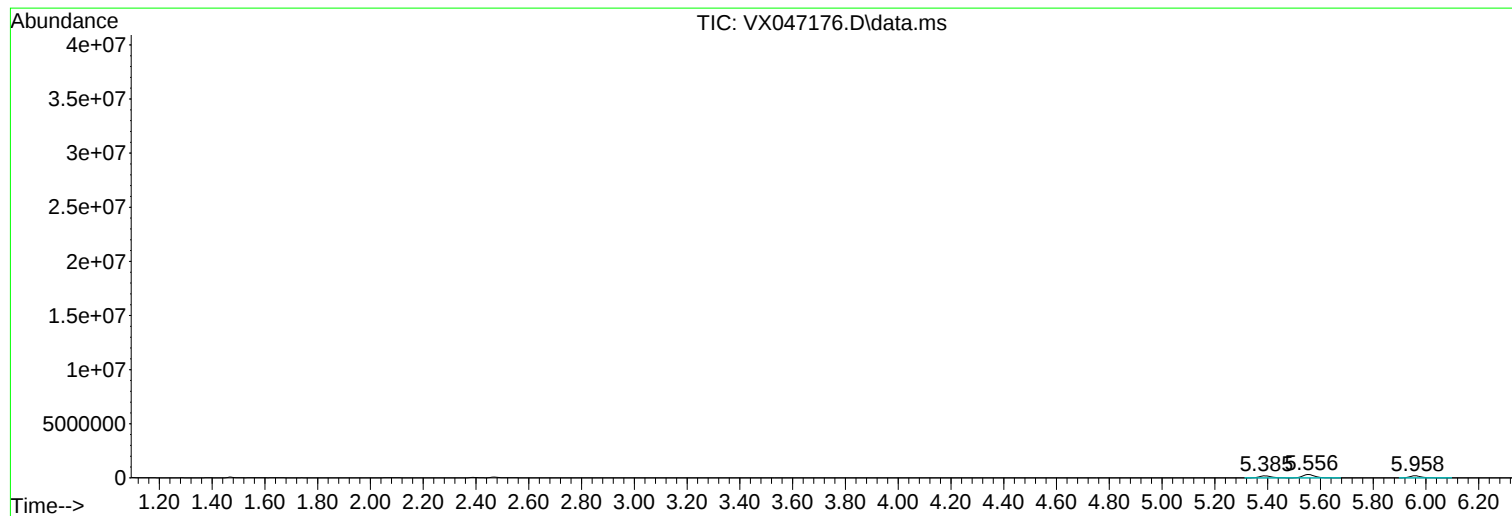
Sum of corrected areas: 65725993

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072125W.M
Quant Title : SW846 8260

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



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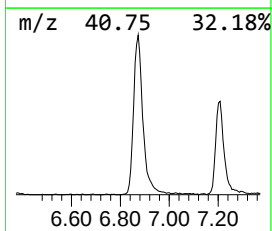
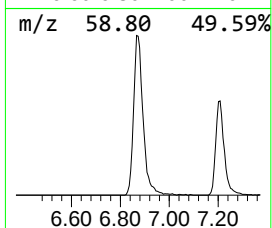
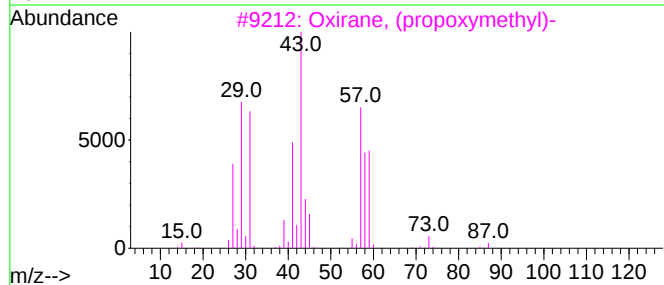
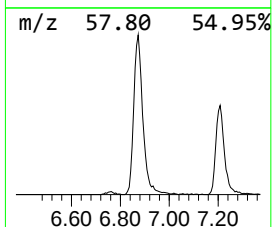
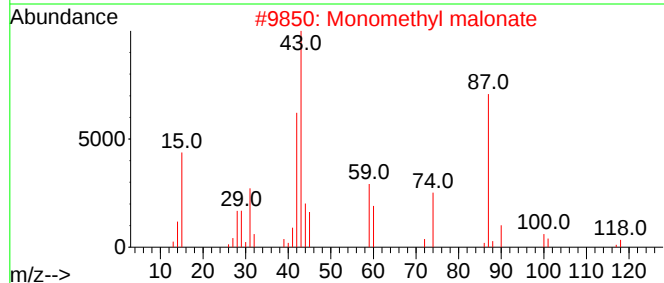
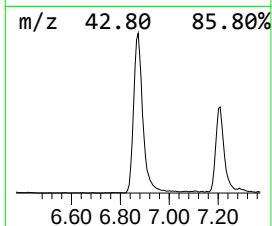
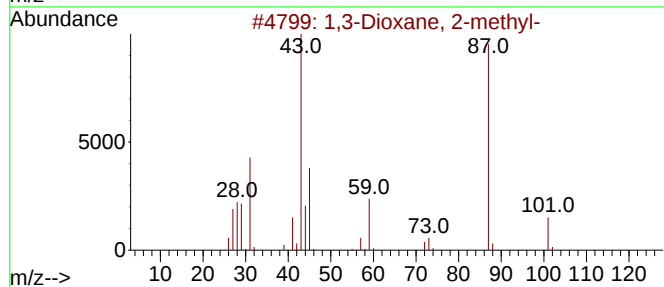
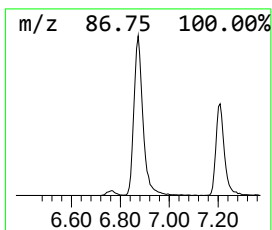
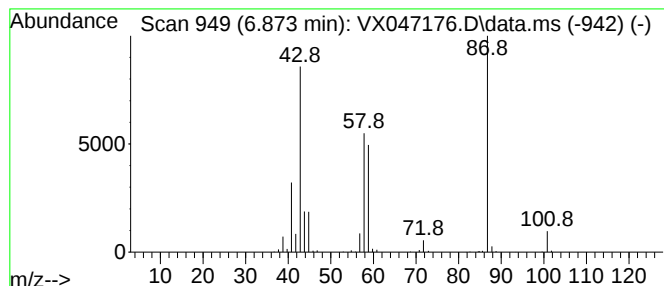
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072125W.M
Quant Title : SW846 8260

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

Peak Number 1 1,3-Dioxane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.873	28.67 ug/l	740726	1,4-Difluorobenzene	6.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxane, 2-methyl-			102	C5H10O2	000626-68-6	50
2	Monomethyl malonate			118	C4H6O4	016695-14-0	42
3	Oxirane, (propoxymethyl)-			116	C6H12O2	003126-95-2	37
4	1,3-Dioxolane, 4-methyl-2-propyl-			130	C7H14O2	004352-99-2	25
5	N,N-Dimethylacetamide			87	C4H9NO	000127-19-5	9



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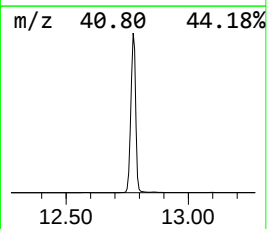
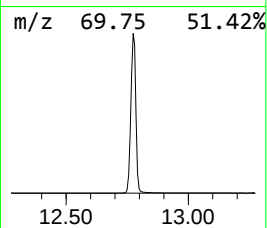
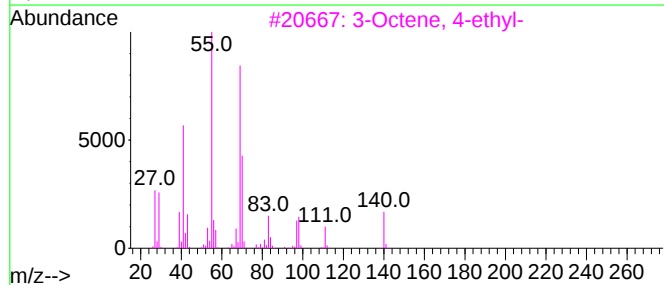
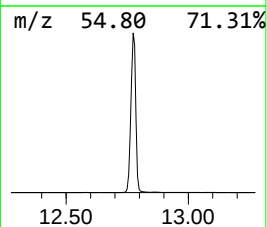
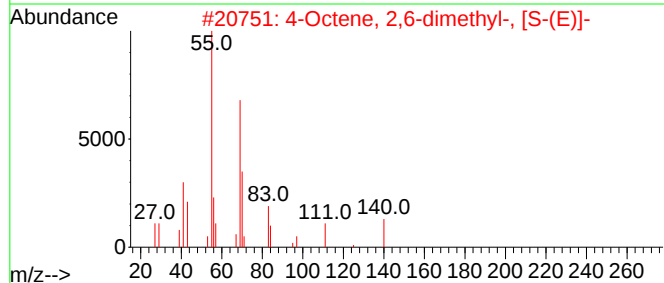
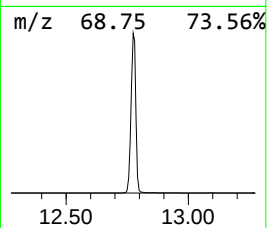
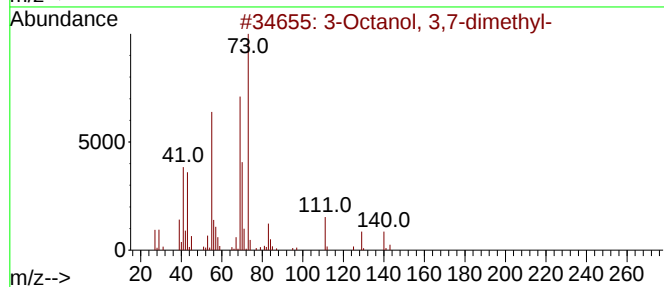
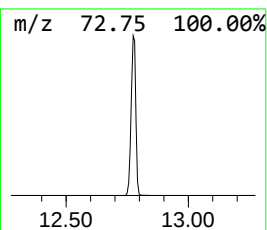
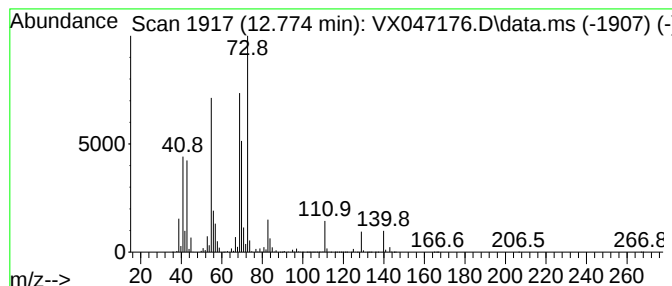
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TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Octanol, 3,7-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.774	1752.51 ug/l	55394900	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	91
2			4-Octene, 2,6-dimethyl-, [S-(E)]-	140	C10H20	062960-76-3	68
3			3-Octene, 4-ethyl-	140	C10H20	053966-51-1	52
4			Cyanamide, dimethyl-	70	C3H6N2	001467-79-4	46
5			1-Heptene, 5-methyl-	112	C8H16	013151-04-7	42



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,3-Dioxane, 2-...	6.873	28.7	ug/l	740726	2	6.763	1291980	50.0
3-Octanol, 3,7-...	12.774	1752.5	ug/l	55394900	4	12.018	1580450	50.0