

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082423\
 Data File : VV031901.D
 Acq On : 24 Aug 2023 11:22
 Operator : SY/MD
 Sample : 04115-02
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 C0AP5

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.1

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_V\Method\SFAMVTR082323WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV031901.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.217	49	55	61	rBV	8752	11797	2.58%	0.341%
2	1.536	147	154	166	rVB	36563	43603	9.52%	1.261%
3	2.063	310	318	331	rVB	48050	69986	15.28%	2.023%
4	3.815	851	863	904	rBV	47235	180824	39.48%	5.228%
5	4.256	981	1000	1029	rBV2	63577	180071	39.32%	5.206%
6	4.963	1203	1220	1245	rBV3	147531	382895	83.60%	11.069%
7	5.539	1387	1399	1427	rBV	132705	281421	61.45%	8.136%
8	5.834	1480	1491	1504	rBV3	5866	12534	2.74%	0.362%
9	5.995	1528	1541	1561	rBV	89266	193783	42.31%	5.602%
10	6.442	1668	1680	1695	rBV	16448	33460	7.31%	0.967%
11	6.924	1817	1830	1844	rBV2	24527	48650	10.62%	1.406%
12	7.249	1921	1931	1952	rBV	70817	132135	28.85%	3.820%
13	7.564	2020	2029	2047	rBV	14023	29183	6.37%	0.844%
14	7.879	2115	2127	2133	rBV3	6610	11711	2.56%	0.339%
15	8.030	2164	2174	2211	rBV	221570	458003	100.00%	13.241%
16	8.182	2211	2221	2243	rVB	29185	59775	13.05%	1.728%
17	8.789	2399	2410	2434	rBV	194812	337277	73.64%	9.750%
18	9.673	2675	2685	2697	rVB3	8429	14813	3.23%	0.428%
19	9.767	2704	2714	2721	rBV5	3263	5270	1.15%	0.152%
20	10.159	2824	2836	2849	rBV	62715	99313	21.68%	2.871%
21	10.320	2873	2886	2900	rBV3	31249	69160	15.10%	1.999%
22	10.435	2912	2922	2932	rBV2	10010	16649	3.64%	0.481%
23	11.098	3122	3128	3142	rBV7	3393	6841	1.49%	0.198%
24	11.188	3146	3156	3181	rVV	227714	361043	78.83%	10.438%
25	11.487	3238	3249	3262	rBV	42428	66846	14.60%	1.932%
26	11.564	3262	3273	3280	rBV	196213	318943	69.64%	9.220%
27	12.197	3454	3470	3478	rBV	9035	14565	3.18%	0.421%
28	12.297	3490	3501	3511	rBV8	4316	8459	1.85%	0.245%
29	13.387	3829	3840	3847	rBV9	2805	5067	1.11%	0.146%
30	13.992	4019	4028	4038	rVB	3135	5002	1.09%	0.145%

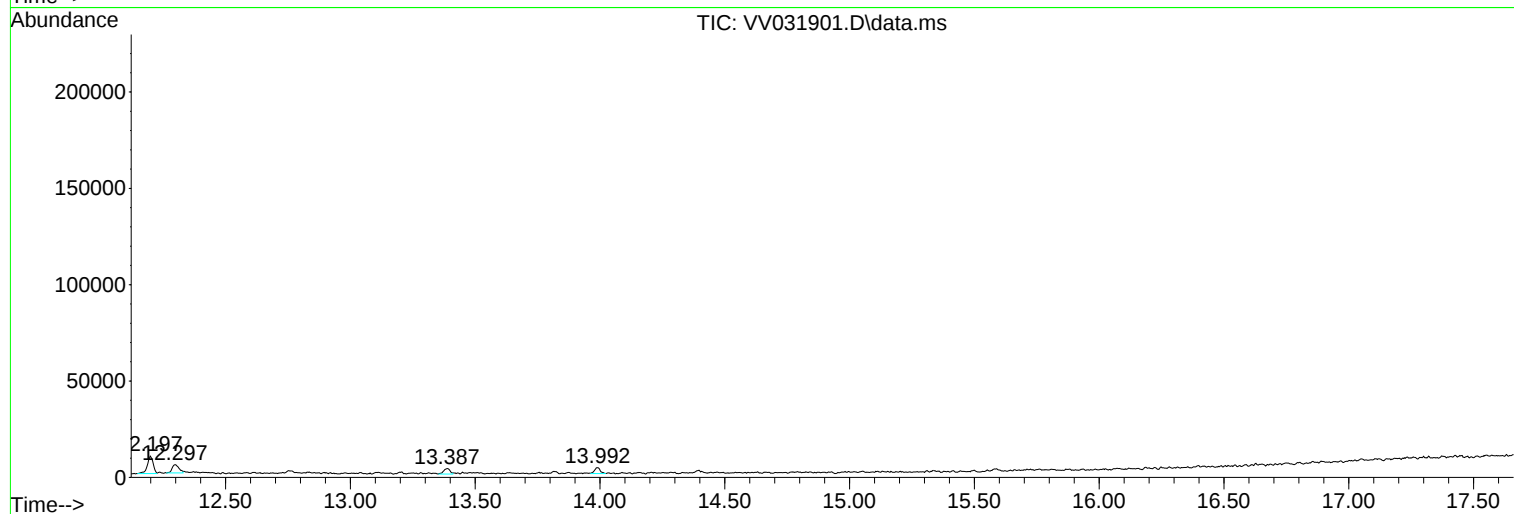
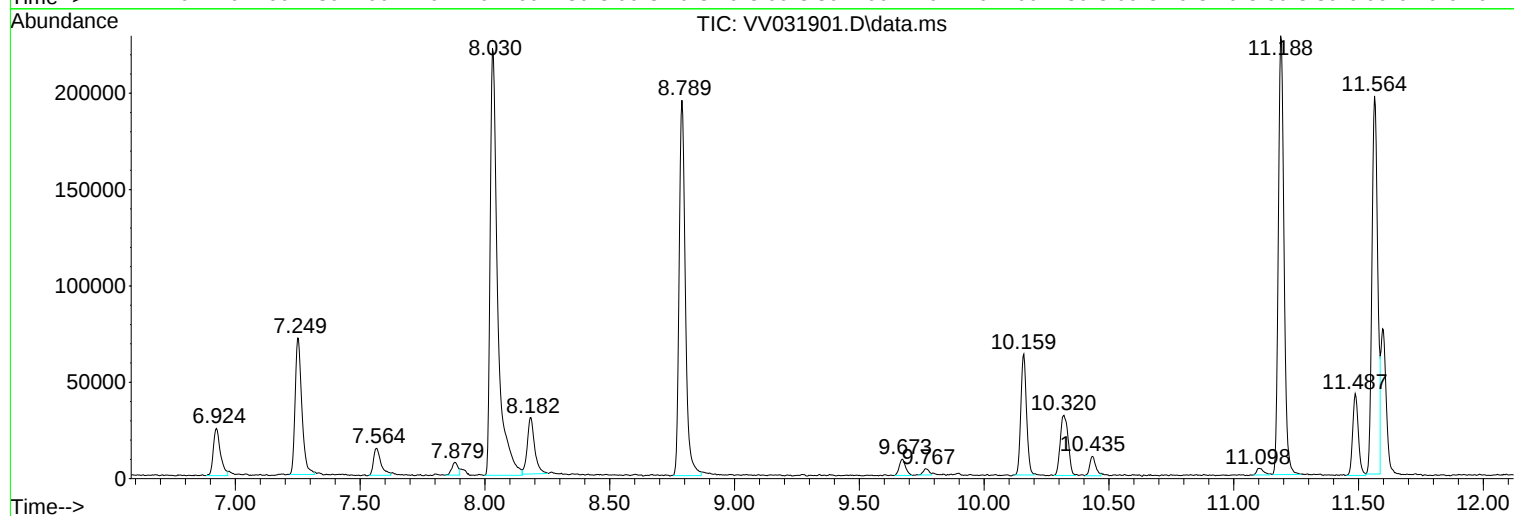
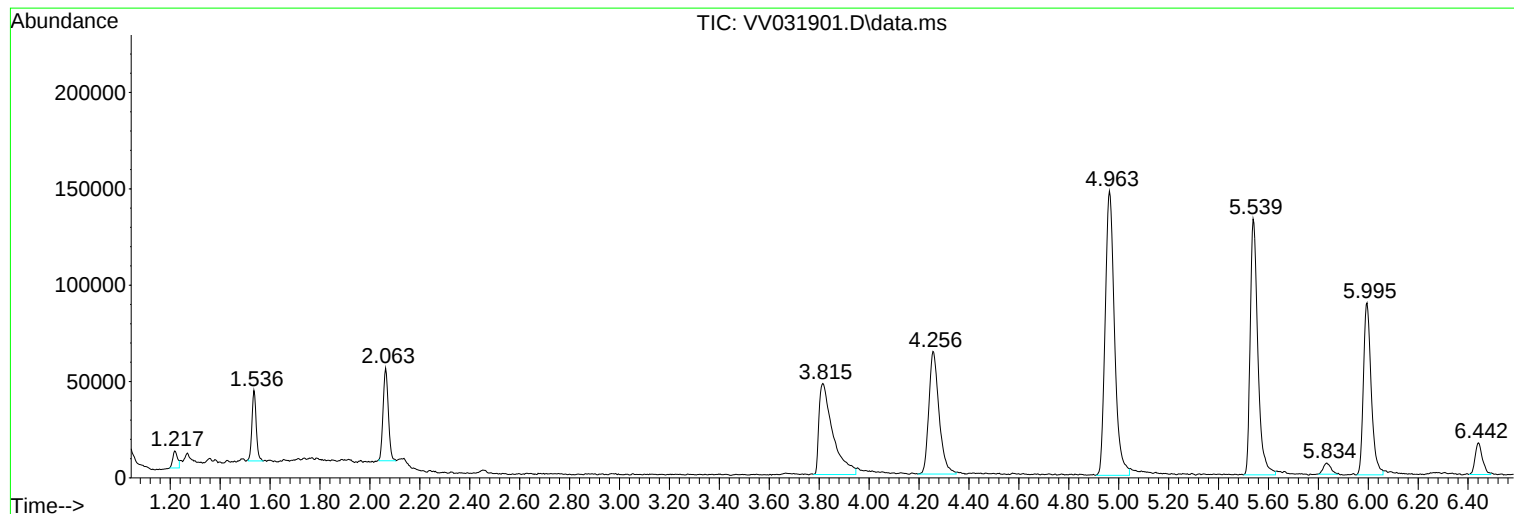
Sum of corrected areas: 3459079

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Instrument :
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ClientSampleId :
C0AP5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR082323WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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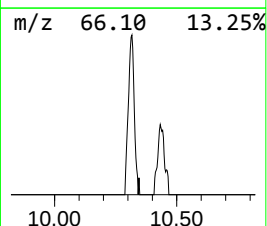
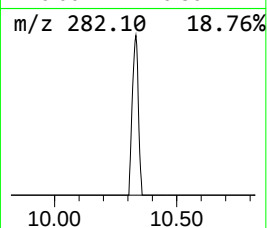
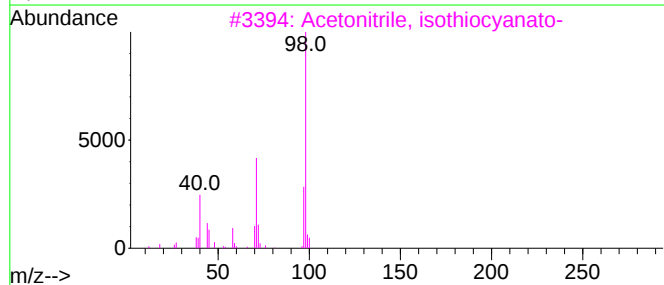
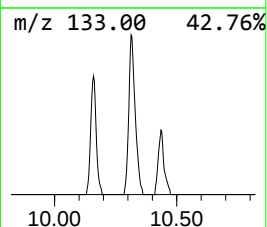
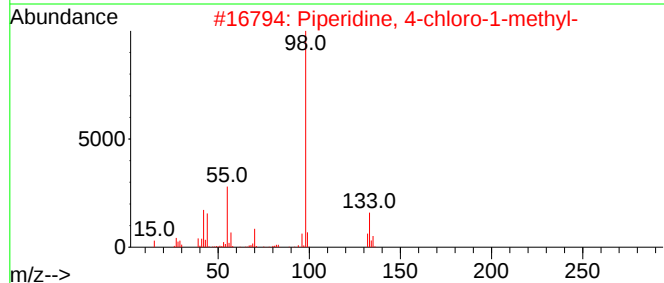
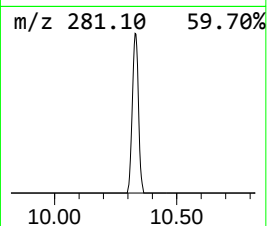
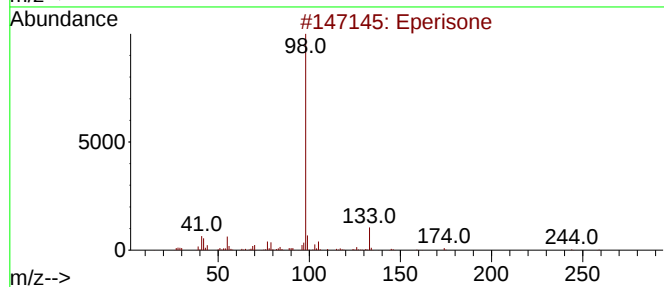
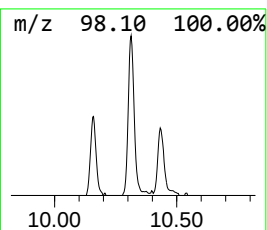
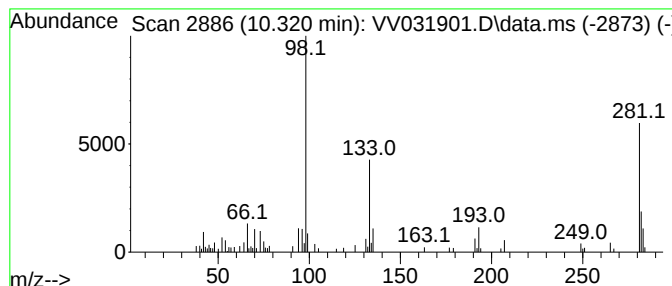
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR082323WMA.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.320	0.96 ug/L	69160	1,4-Dichlorobenzene-d4	11.188

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eperisone	259	C17H25NO	064840-90-0	37
2			Piperidine, 4-chloro-1-methyl-	133	C6H12ClN	005570-77-4	37
3			Acetonitrile, isothiocyanato-	98	C3H2N2S	032772-75-1	27
4			Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	27
5			Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	22



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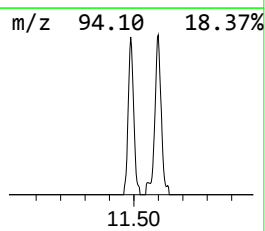
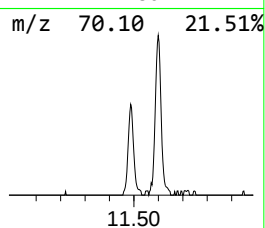
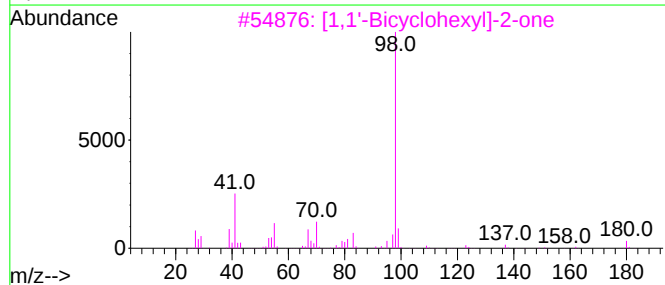
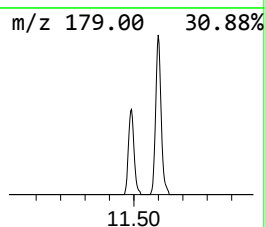
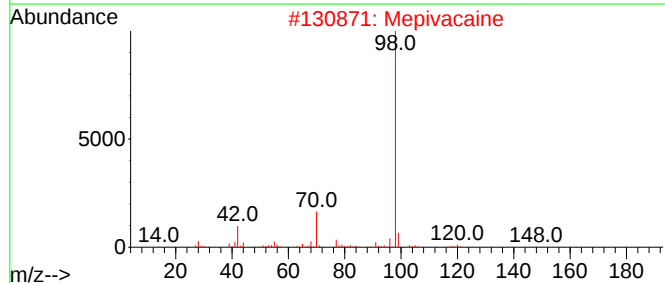
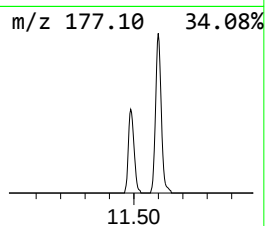
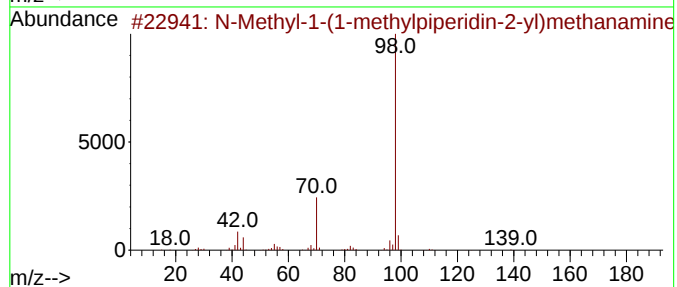
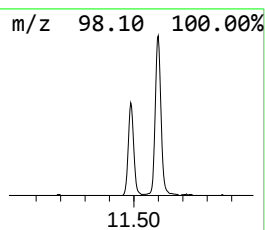
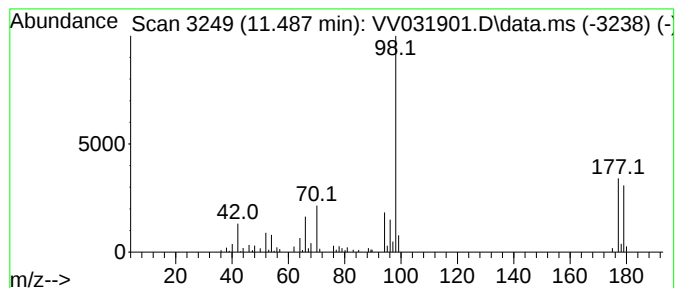
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Quant Title : TRACE VOA SFAM1.0

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TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-02 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.487	0.93 ug/L	66846	1,4-Dichlorobenzene-d4	11.188

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Methyl-1-(1-methylpiperidin-2-yl)methanamine	142	C8H18N2	1000484-42-5	47
2			Mepivacaine	246	C15H22N2O	000096-88-8	38
3			[1,1'-Bicyclohexyl]-2-one	180	C12H20O	000090-42-6	38
4			Benzen-d5-amine	98	C6H2D5N	004165-61-1	38
5			(R)-1-Ethyl-2-pyrrolidinecarboxamide	142	C7H14N2O	1000237-69-9	37



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	10.320	1.0	ug/L	69160	3	11.188	361043	5.0
unknown-02	11.487	0.9	ug/L	66846	3	11.188	361043	5.0