

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV072122\
 Data File : VV026876.D
 Acq On : 21 Jul 2022 15:56
 Operator : SY/MD
 Sample : N3809-06
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VHBLK001

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\SFAMVTR071122WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV026876.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.301	72	81	102	rVB	157633	188548	20.56%	2.561%
2	1.561	155	162	179	rVB	132208	154026	16.79%	2.092%
3	2.101	320	330	346	rVV	257732	370466	40.39%	5.033%
4	2.179	348	354	368	rVB3	8677	18077	1.97%	0.246%
5	2.503	444	455	469	rBV	42785	71512	7.80%	0.971%
6	2.671	491	507	514	rBV4	8654	24040	2.62%	0.327%
7	3.034	606	620	635	rBV3	12452	25313	2.76%	0.344%
8	3.892	876	887	938	rBV	86050	309735	33.77%	4.208%
9	4.343	1011	1027	1056	rBV	168721	436814	47.62%	5.934%
10	5.043	1229	1245	1275	rBV2	367024	917250	100.00%	12.460%
11	5.613	1408	1422	1461	rBV	243311	520826	56.78%	7.075%
12	5.908	1503	1514	1530	rBV9	8459	19710	2.15%	0.268%
13	6.066	1549	1563	1594	rBV	212183	455792	49.69%	6.192%
14	6.577	1712	1722	1742	rVB2	26469	54283	5.92%	0.737%
15	6.985	1836	1849	1869	rBV	78256	154878	16.89%	2.104%
16	7.313	1939	1951	1971	rBV	386250	699926	76.31%	9.508%
17	7.622	2035	2047	2077	rBV	46250	91789	10.01%	1.247%
18	7.937	2133	2145	2152	rBV4	9870	18207	1.98%	0.247%
19	8.088	2182	2192	2235	rBV	343066	697225	76.01%	9.472%
20	8.850	2417	2429	2454	rBV	388682	658689	71.81%	8.948%
21	10.214	2841	2853	2870	rBV	146272	239058	26.06%	3.248%
22	10.381	2893	2905	2916	rBV3	21542	34816	3.80%	0.473%
23	11.249	3164	3175	3195	rBV	363182	581071	63.35%	7.894%
24	11.622	3279	3291	3311	rBV	348134	570998	62.25%	7.757%
25	12.252	3475	3487	3496	rVB3	6588	11146	1.22%	0.151%
26	13.265	3794	3802	3811	rVB6	6040	9582	1.04%	0.130%
27	13.506	3868	3877	3887	rBV	8528	14737	1.61%	0.200%
28	13.744	3944	3951	3961	rVB6	8157	12769	1.39%	0.173%

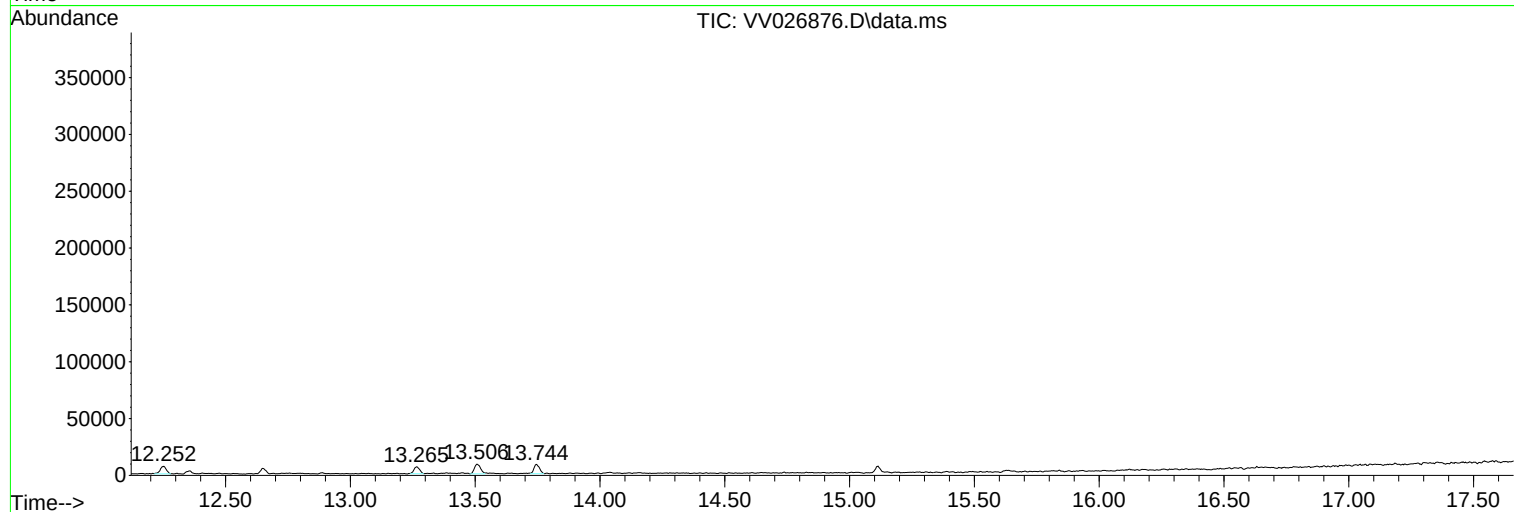
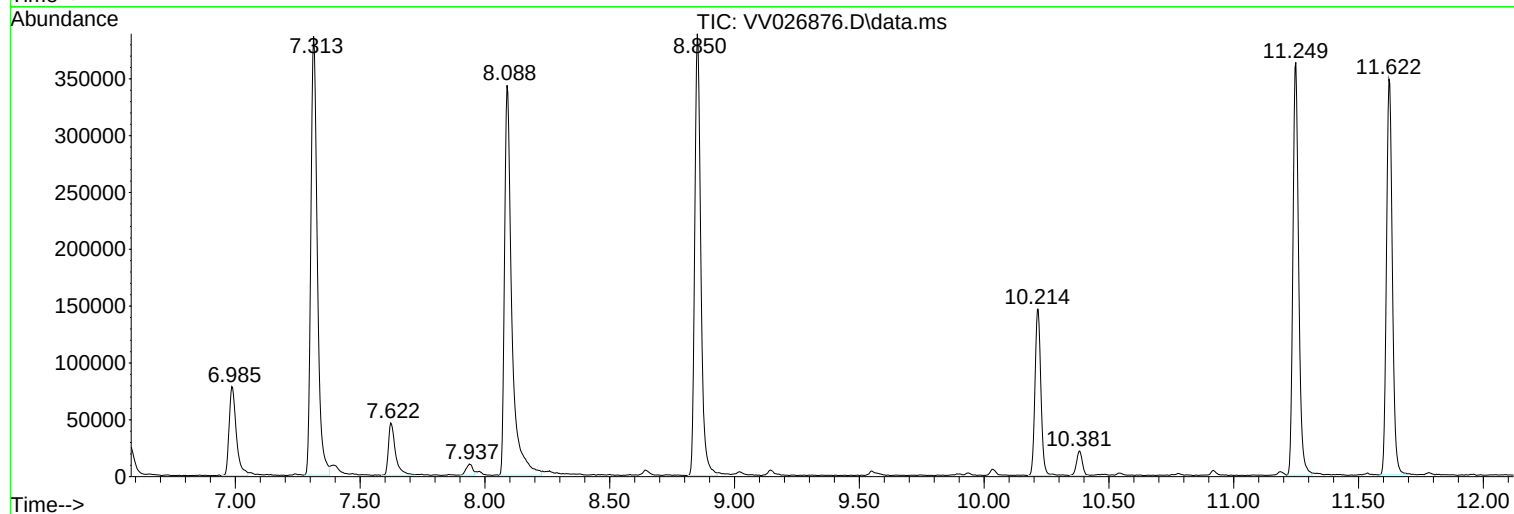
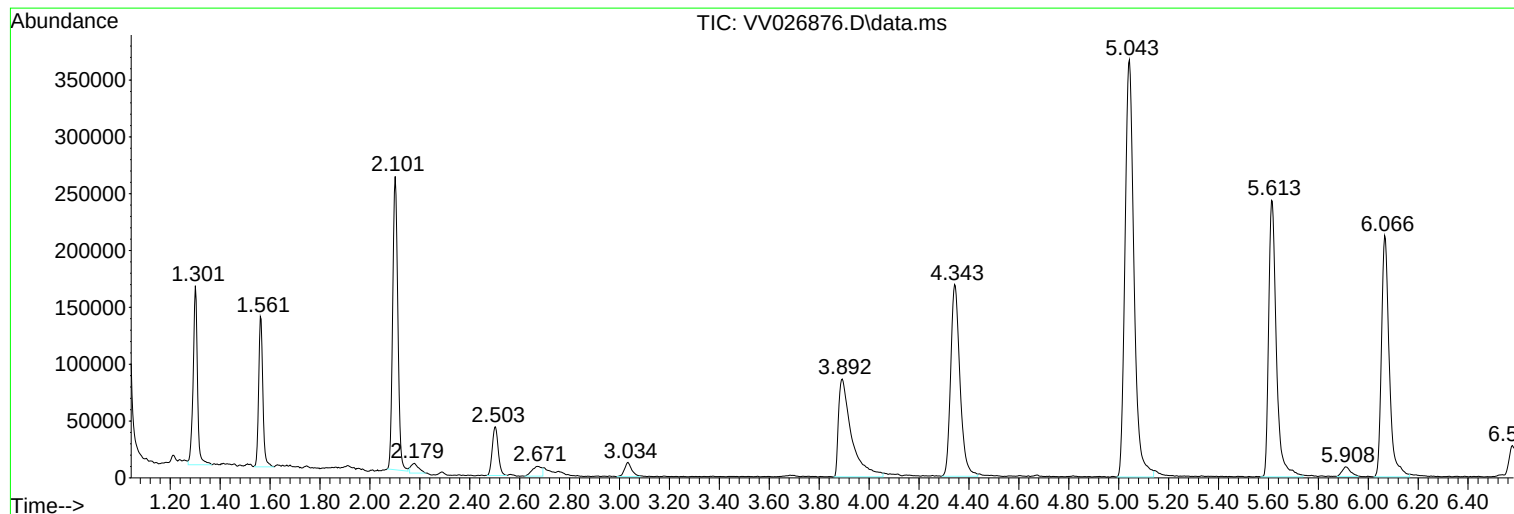
Sum of corrected areas: 7361283

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR071122WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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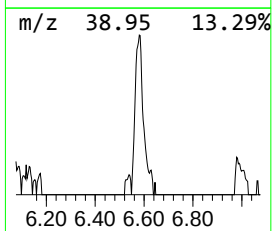
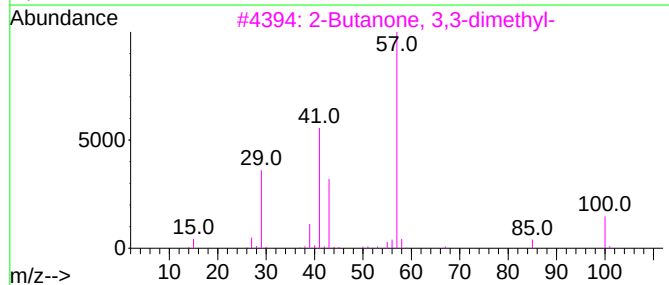
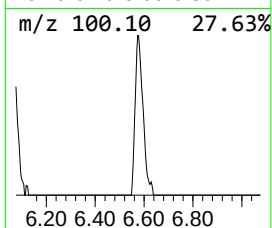
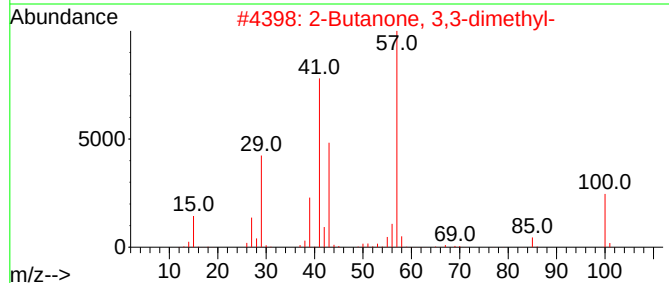
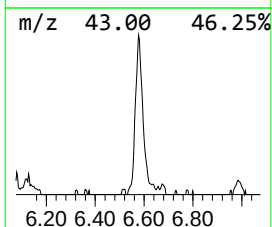
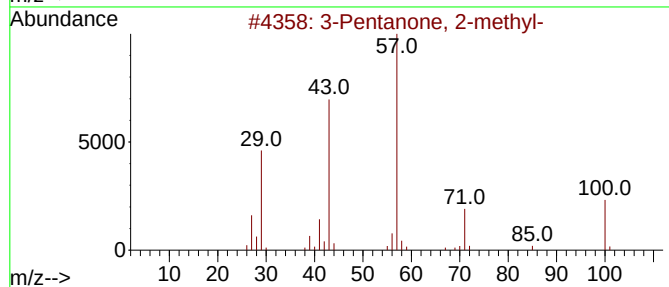
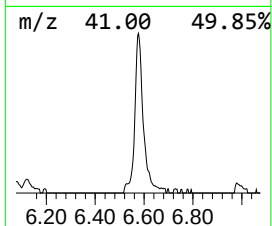
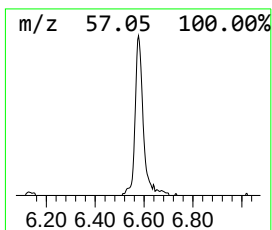
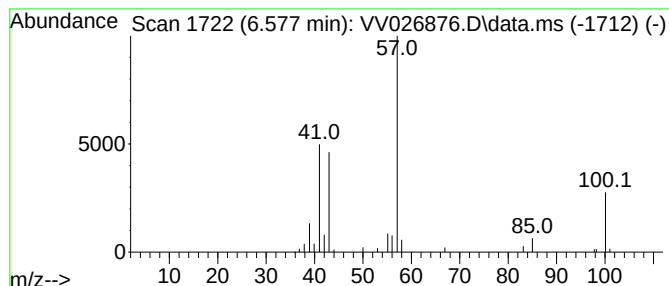
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR071122WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.577	0.52 ug/L	54283	1,4-Difluorobenzene	5.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Pentanone, 2-methyl-	100	C6H12O	000565-69-5	80
2			2-Butanone, 3,3-dimethyl-	100	C6H12O	000075-97-8	72
3			2-Butanone, 3,3-dimethyl-	100	C6H12O	000075-97-8	64
4			2-Butanone, 3,3-dimethyl-	100	C6H12O	000075-97-8	53
5			3-Pentanone, 2-methyl-	100	C6H12O	000565-69-5	42



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Pentanone, 2-...	6.577	0.5	ug/L	54283	1	5.616	520826	5.0