

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU121319\
 Data File : VU036218.D
 Acq On : 14 Dec 2019 00:07
 Operator : JC/MD
 Sample : K6240-22
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BFQS7

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

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_U\METHOD\SOMUTR121319WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.372	3	10	21	rBV2	56418	58776	5.88%	0.799%
2	1.610	76	84	94	rVB	148249	163759	16.39%	2.227%
3	1.932	175	184	195	rBV	126382	165117	16.53%	2.246%
4	2.591	377	389	403	rBV	168963	272812	27.31%	3.710%
5	2.684	409	418	433	rVB3	5970	14666	1.47%	0.199%
6	2.822	451	461	470	rBV	7198	12611	1.26%	0.172%
7	4.700	1025	1045	1084	rBV	116641	407533	40.80%	5.543%
8	5.112	1157	1173	1199	rBV	171080	372876	37.33%	5.071%
9	5.768	1357	1377	1403	rBV2	326182	845907	84.69%	11.505%
10	6.288	1520	1539	1561	rBV2	267812	518253	51.88%	7.049%
11	6.732	1663	1677	1697	rBV	205951	407765	40.82%	5.546%
12	7.603	1935	1948	1969	rBV2	93362	169515	16.97%	2.306%
13	7.931	2037	2050	2065	rBV	343246	609414	61.01%	8.289%
14	8.002	2065	2072	2081	rVB4	7465	13760	1.38%	0.187%
15	8.214	2125	2138	2155	rBV2	62840	111432	11.16%	1.516%
16	8.677	2266	2282	2320	rBV	452447	998856	100.00%	13.585%
17	8.822	2320	2327	2341	rVB7	6704	13056	1.31%	0.178%
18	9.446	2509	2521	2554	rBV	408879	723785	72.46%	9.844%
19	10.783	2924	2937	2953	rVB	197013	325465	32.58%	4.427%
20	11.835	3251	3264	3279	rBV	331265	555146	55.58%	7.550%
21	12.217	3370	3383	3398	rVB	353107	591982	59.27%	8.051%

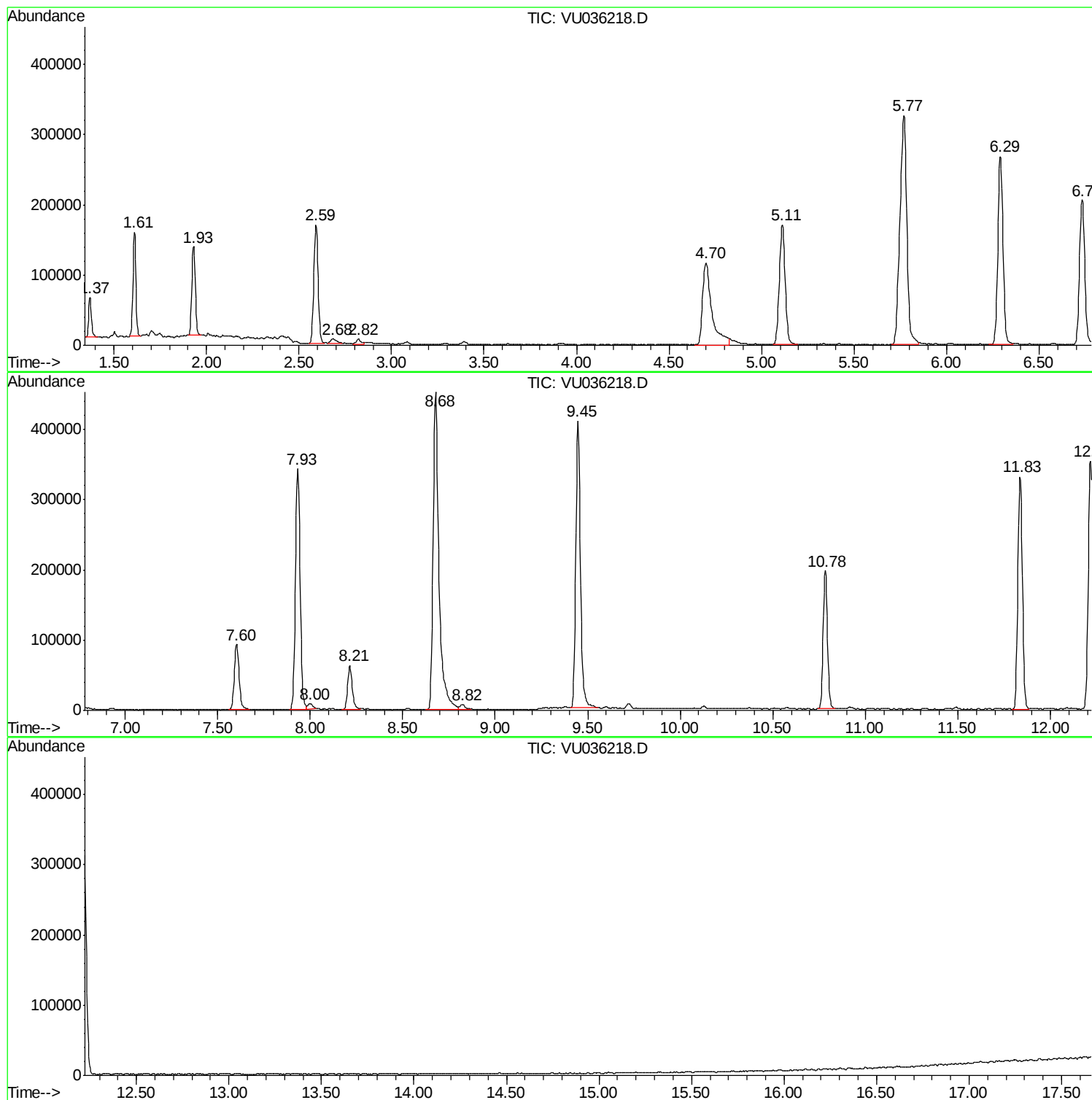
Sum of corrected areas: 7352486

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR121319WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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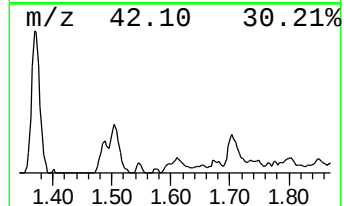
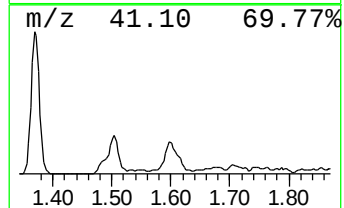
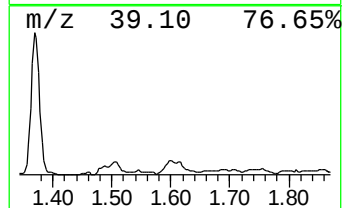
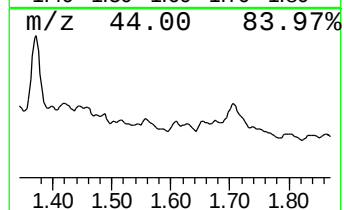
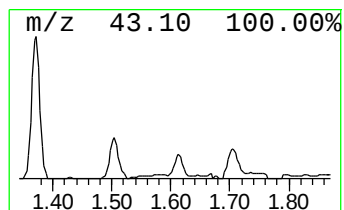
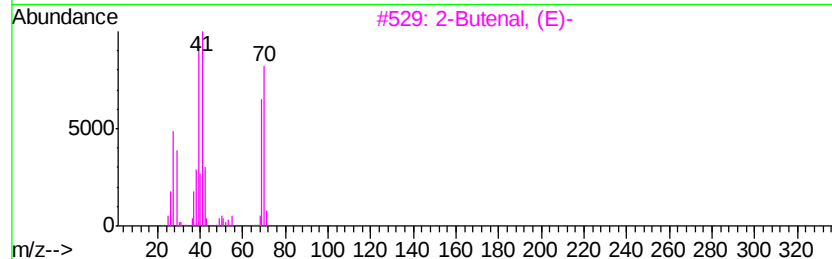
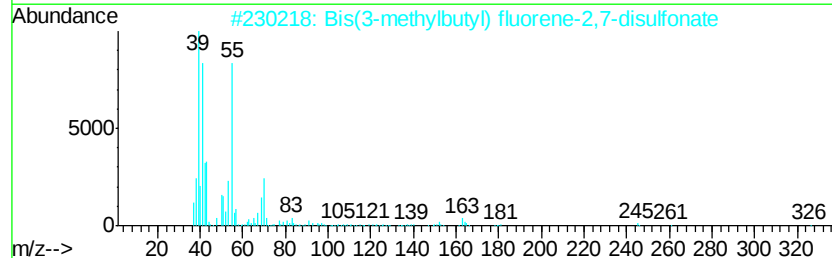
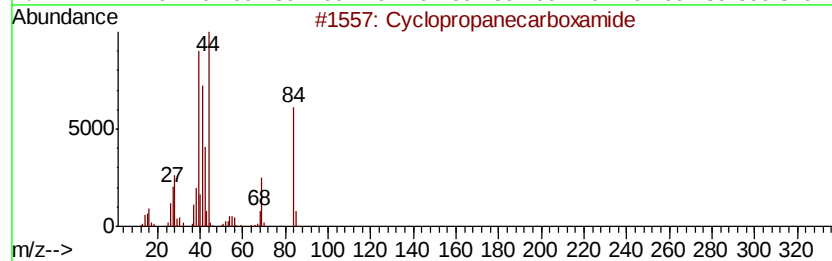
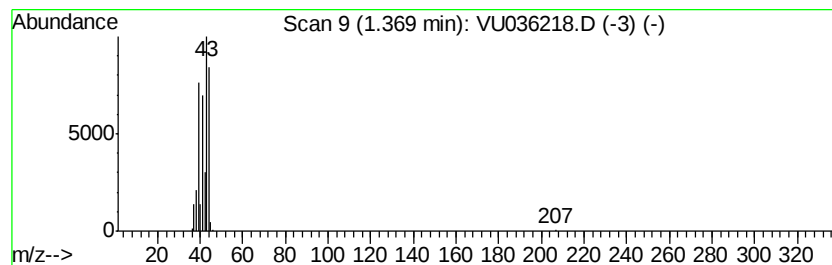
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR121319WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Cyclopropanecarboxamide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.37	0.57 ug/L	58776	1,4-Difluorobenzene	6.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropanecarboxamide	85	C4H7NO	006228-73-5	78
2		Bis(3-methylbutyl) fluorene-2,7-...	466	C23H30O6S2	253664-95-8	47
3		2-Butenal, (E)-	70	C4H6O	000123-73-9	47
4		2-Hexynoic acid	112	C6H8O2	000764-33-0	39
5		Imidazol-5(2H)-one, 4-amino-2-(2...	196	C7H8N4O3	318259-17-5	38



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclopropanecarbo...	1.37	0.6	ug/L	58776	1	6.29	518253	5.0