

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU010921\
 Data File : VU041885.D
 Acq On : 07 Jan 2021 20:16
 Operator : SY/MD
 Sample : M1026-16
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BFXL7

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\SOMUTR121220WMA.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	1.478	37	43	51	rVB2	16720	17173	2.40%	0.315%
2	1.604	74	82	94	rVB	62873	67807	9.49%	1.242%
3	1.922	172	181	191	rBV	64391	78718	11.02%	1.442%
4	2.089	224	233	242	rBV	34357	46641	6.53%	0.854%
5	2.147	243	251	262	rVB	25056	35058	4.91%	0.642%
6	2.581	372	386	401	rBV2	122993	227682	31.86%	4.170%
7	3.060	522	535	555	rVB	63056	118160	16.54%	2.164%
8	3.199	567	578	596	rBV	4589	9786	1.37%	0.179%
9	3.379	617	634	651	rBV3	18788	44488	6.23%	0.815%
10	3.890	775	793	811	rBV	38708	92210	12.90%	1.689%
11	4.649	1011	1029	1078	rBV3	102999	330567	46.26%	6.055%
12	5.079	1146	1163	1190	rBV2	116857	290216	40.61%	5.316%
13	5.330	1223	1241	1257	rVB3	39408	90803	12.71%	1.663%
14	5.739	1347	1368	1384	rBV3	212159	550695	77.07%	10.087%
15	6.259	1513	1530	1552	rBV	174736	331101	46.34%	6.065%
16	6.552	1608	1621	1639	rBV	67269	126466	17.70%	2.316%
17	6.700	1653	1667	1687	rBV	131571	260274	36.42%	4.767%
18	7.571	1927	1938	1952	rBV2	60931	106785	14.94%	1.956%
19	7.902	2028	2041	2057	rBV	230449	396079	55.43%	7.255%
20	7.970	2057	2062	2077	rVB2	4582	8159	1.14%	0.149%
21	8.182	2116	2128	2146	rBV	40196	68945	9.65%	1.263%
22	8.558	2234	2245	2256	rBV2	35840	59973	8.39%	1.098%
23	8.639	2256	2270	2300	rBV	379867	714576	100.00%	13.089%
24	9.420	2500	2513	2532	rBV	224276	386153	54.04%	7.073%
25	10.754	2916	2928	2946	rVB	115083	188617	26.40%	3.455%
26	11.806	3240	3255	3271	rBV	206238	332898	46.59%	6.098%
27	12.053	3319	3332	3350	rBV	57835	113562	15.89%	2.080%
28	12.185	3361	3373	3388	rVB	227281	365969	51.21%	6.703%

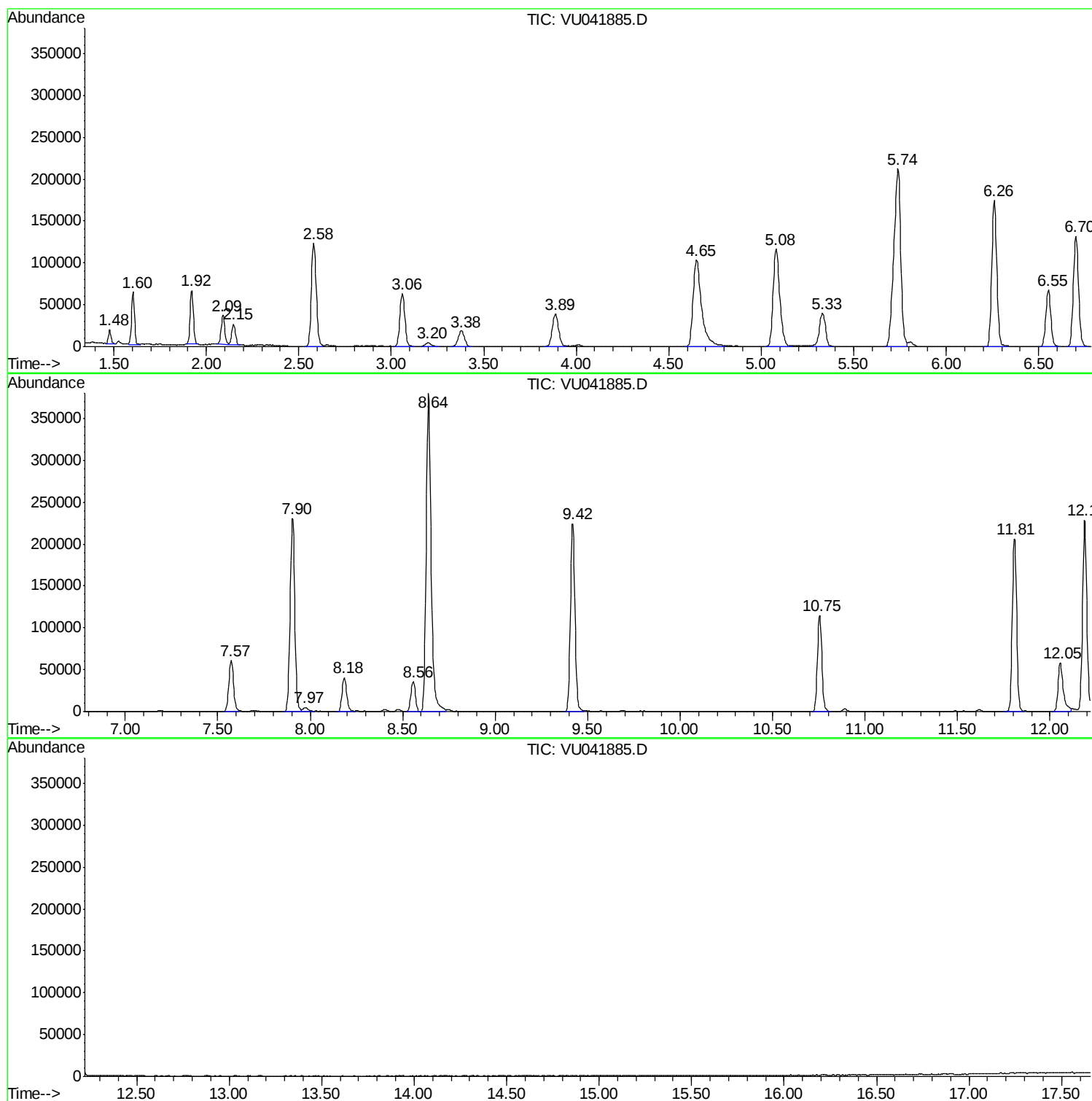
Sum of corrected areas: 5459561

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Instrument :
MSVOA_U
ClientSampleId :
BFXL7

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR121220WMA.M
Quant Title : TRACE VOA SOM01.0

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



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Instrument :
MSVOA_U
ClientSampleId :
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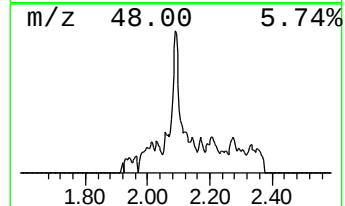
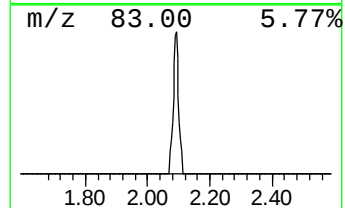
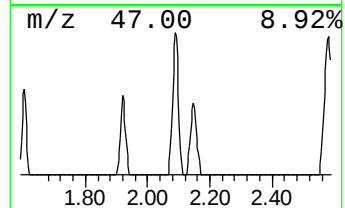
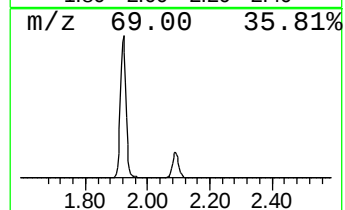
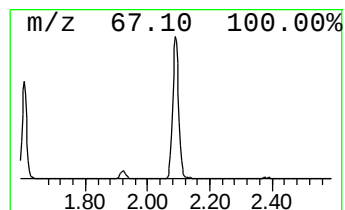
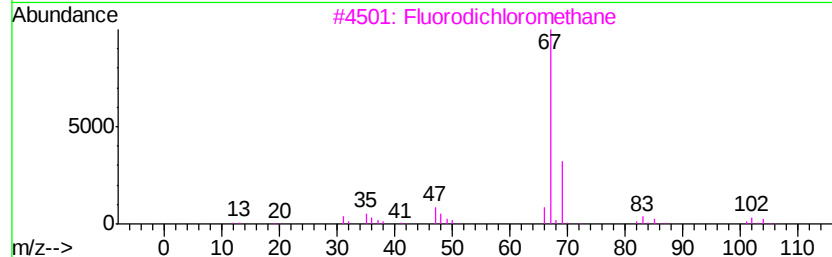
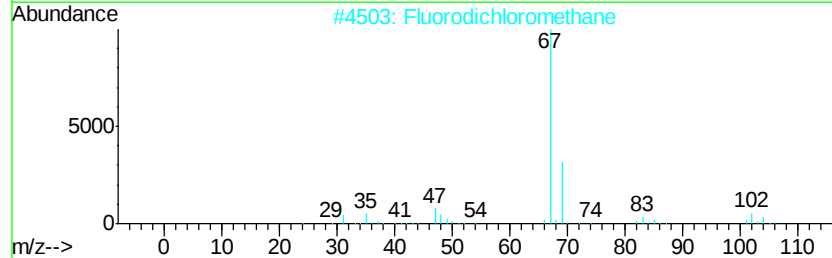
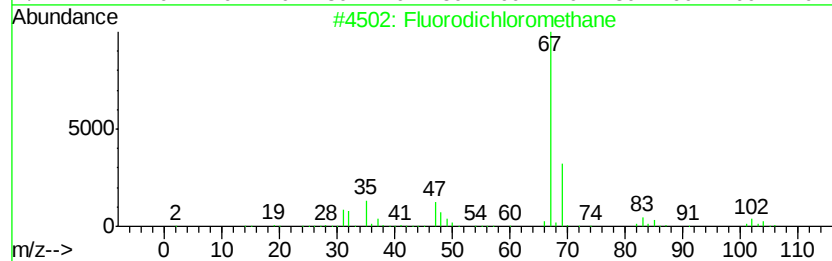
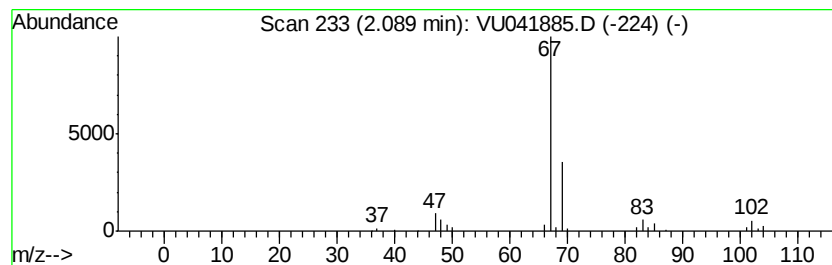
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR121220WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 Fluorodichloromethane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.09	0.70 ug/L	46641	1,4-Difluorobenzene	6.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluorodichloromethane	102	CHCl2F	000075-43-4	91
2			Fluorodichloromethane	102	CHCl2F	000075-43-4	91
3			Fluorodichloromethane	102	CHCl2F	000075-43-4	91
4			1-Butyne, 3,3-dimethyl-	82	C6H10	000917-92-0	7
5			Cyclopentene, 1-chloro-	102	C5H7Cl	000930-29-0	7



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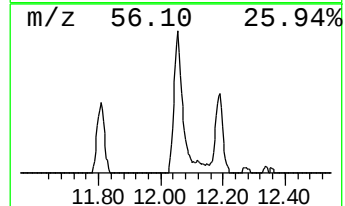
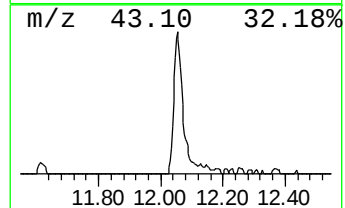
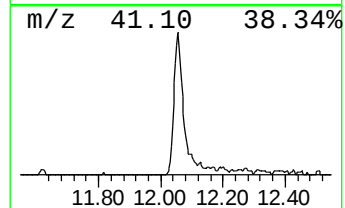
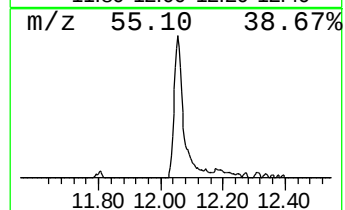
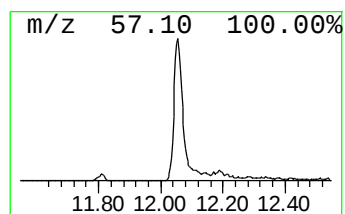
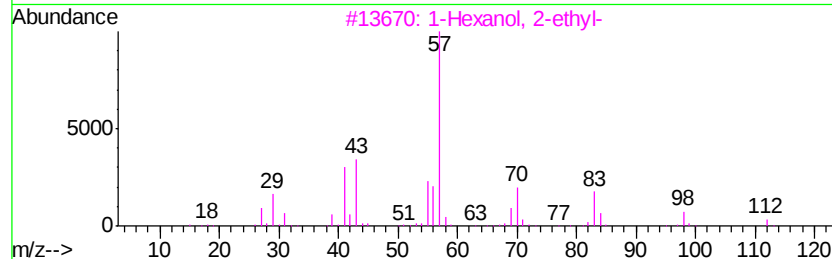
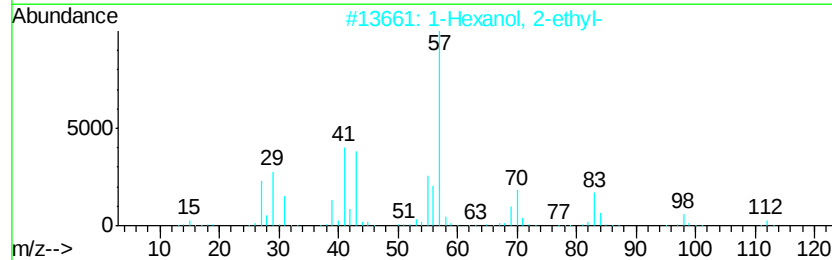
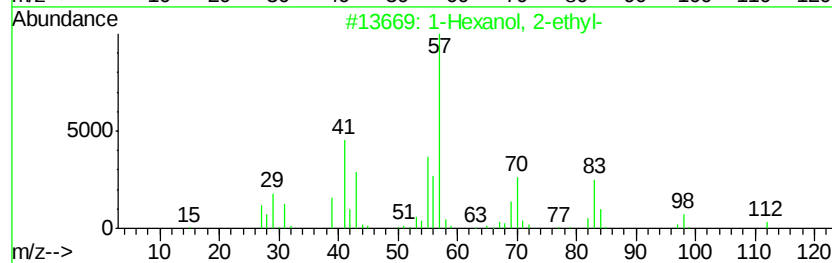
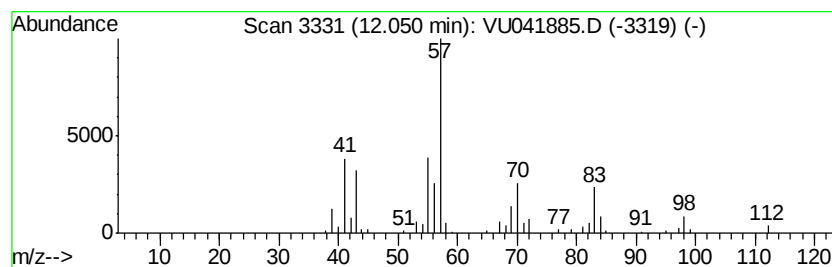
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Quant Title : TRACE VOA SOM01.0

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TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.05	1.71 ug/L	113562	1,4-Dichlorobenzene-d4	11.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	86
2			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
3			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
4			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	59
5			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Fluorodichloromet...	2.09	0.7	ug/L	46641	1	6.26	331101	5.0
1-Hexanol, 2-ethyl-	12.05	1.7	ug/L	113562	3	11.81	332898	5.0