

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV101618\
 Data File : VV008297.D
 Acq On : 16 Oct 2018 14:29
 Operator : SY/MD
 Sample : J5406-09
 Misc : 25.0 mL/MSVOA V/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 C0AK3

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_V\METHOD\SOMVTR101518WMA.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	1.225	37	42	46	rBV	9493	8158	2.39%	0.312%
2	1.319	64	71	82	rVB	40835	44177	12.96%	1.689%
3	1.586	147	154	166	rVB	32202	36260	10.64%	1.387%
4	2.132	315	324	336	rBV	65550	93407	27.40%	3.572%
5	3.003	587	595	606	rVB2	13694	27009	7.92%	1.033%
6	3.984	882	900	928	rBV	35657	132519	38.88%	5.067%
7	4.405	1015	1031	1052	rBV	49727	118415	34.74%	4.528%
8	4.811	1148	1157	1174	rVB6	1615	4035	1.18%	0.154%
9	5.100	1230	1247	1266	rBV2	129375	296996	87.13%	11.357%
10	5.601	1394	1403	1405	rBV4	2671	3572	1.05%	0.137%
11	5.666	1411	1423	1440	rVB	105544	204694	60.05%	7.827%
12	5.952	1501	1512	1523	rVB4	7640	14640	4.30%	0.560%
13	6.122	1550	1565	1585	rBV	72125	144978	42.53%	5.544%
14	6.634	1714	1724	1738	rVB8	3807	8288	2.43%	0.317%
15	7.035	1837	1849	1870	rBV2	36300	66013	19.37%	2.524%
16	7.199	1884	1900	1916	rVB2	26360	53425	15.67%	2.043%
17	7.363	1938	1951	1966	rBV	142472	251862	73.89%	9.631%
18	7.669	2033	2046	2059	rBV	21722	38232	11.22%	1.462%
19	8.148	2180	2195	2231	rBV5	127436	340858	100.00%	13.034%
20	8.293	2235	2240	2251	rVB5	2058	3947	1.16%	0.151%
21	8.900	2416	2429	2445	rBV	145688	243289	71.38%	9.303%
22	9.180	2506	2516	2524	rBV6	4422	8177	2.40%	0.313%
23	10.264	2843	2853	2864	rBV	39751	64507	18.92%	2.467%
24	11.299	3164	3175	3202	rVB	106872	198834	58.33%	7.603%
25	11.585	3251	3264	3279	rBV3	13827	29861	8.76%	1.142%
26	11.675	3282	3292	3311	rVB	101735	178983	52.51%	6.844%

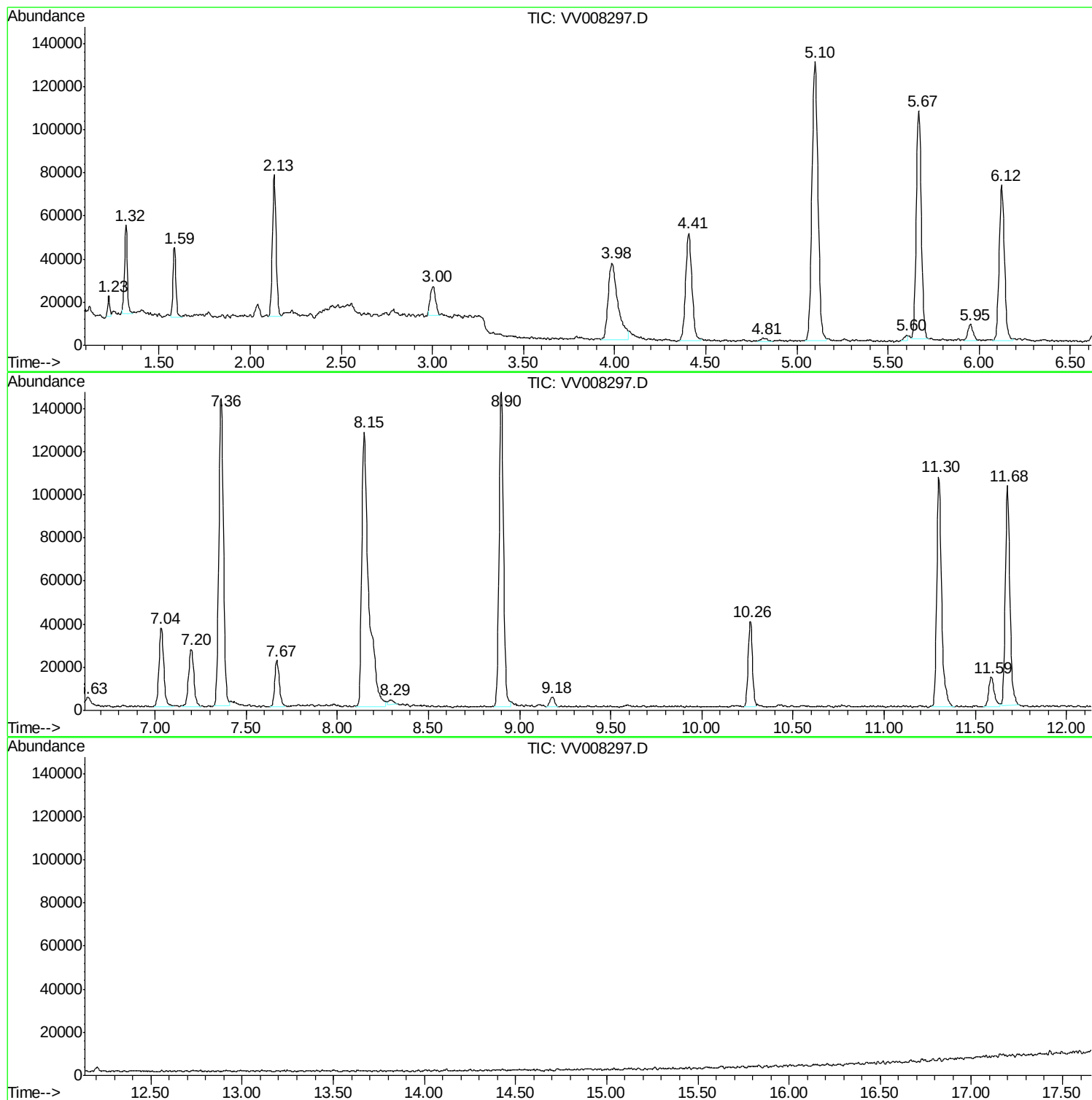
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

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



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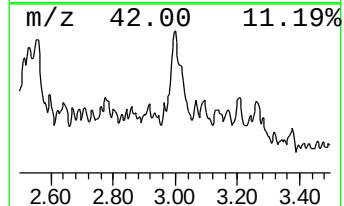
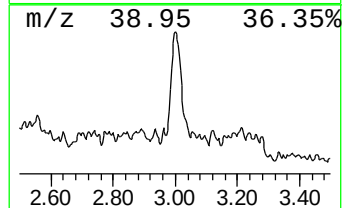
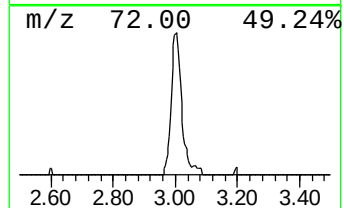
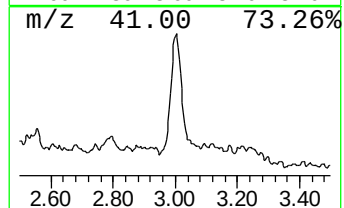
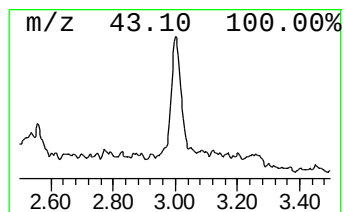
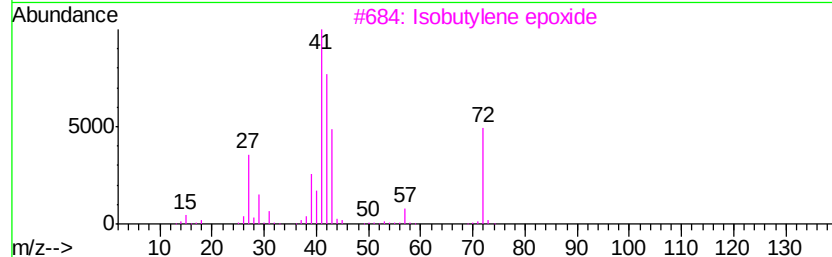
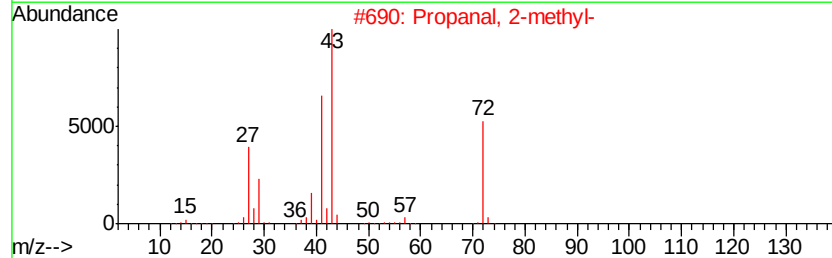
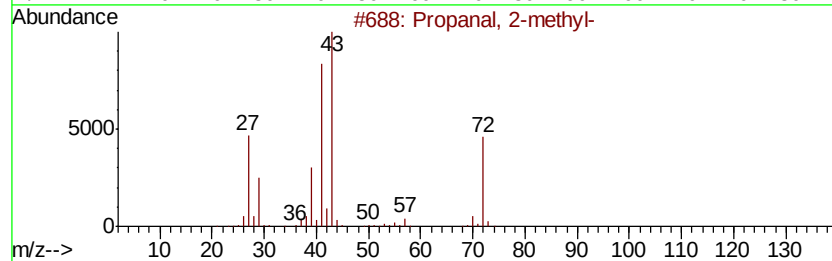
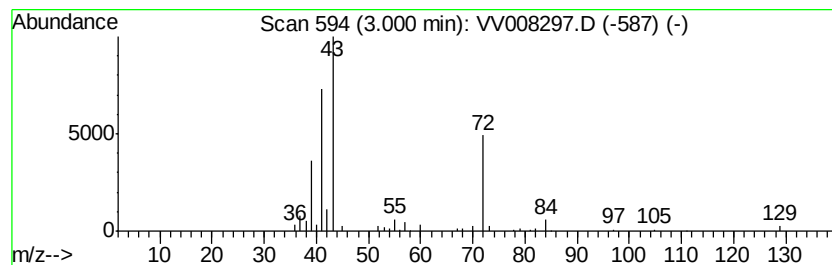
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 Propanal, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.00	0.66 ug/L	27009	1,4-Difluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanal, 2-methyl-	72	C4H8O	000078-84-2	72
2		Propanal, 2-methyl-	72	C4H8O	000078-84-2	72
3		Isobutylene epoxide	72	C4H8O	000558-30-5	50
4		2-Propenoic acid, 3-phenyl-, 2-m...	202	C13H14O2	054889-46-2	25
5		Isobutylene epoxide	72	C4H8O	000558-30-5	9



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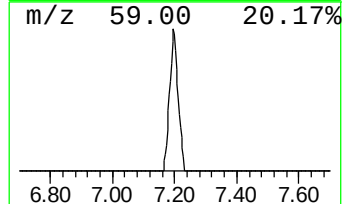
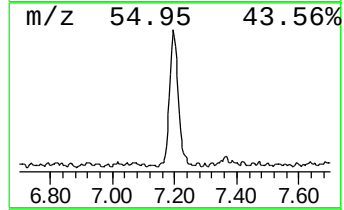
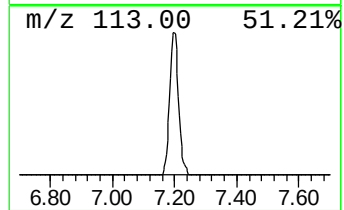
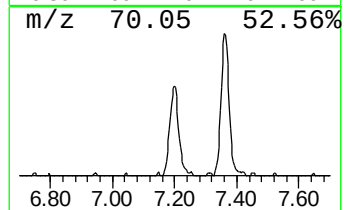
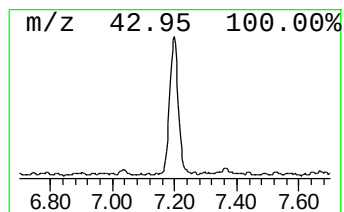
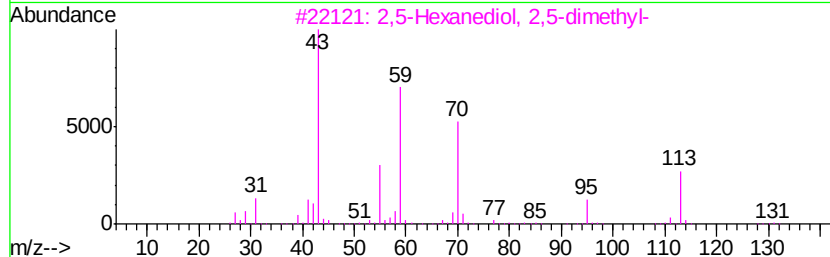
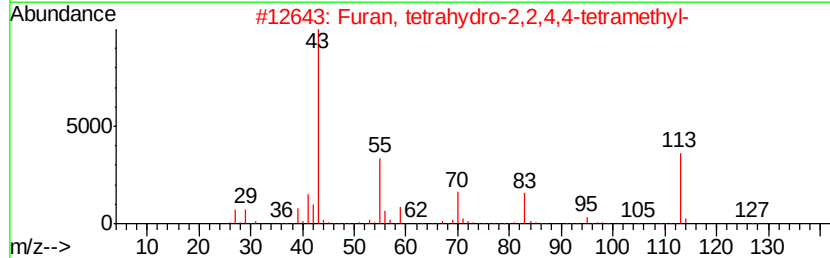
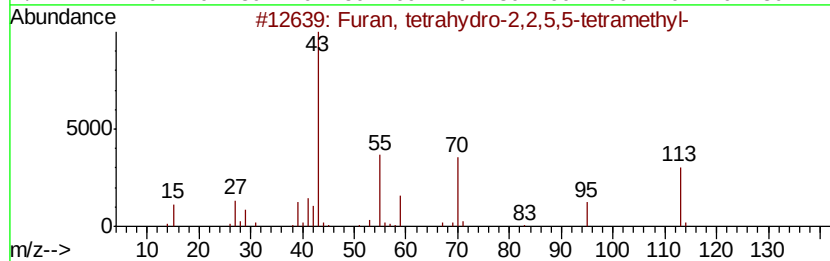
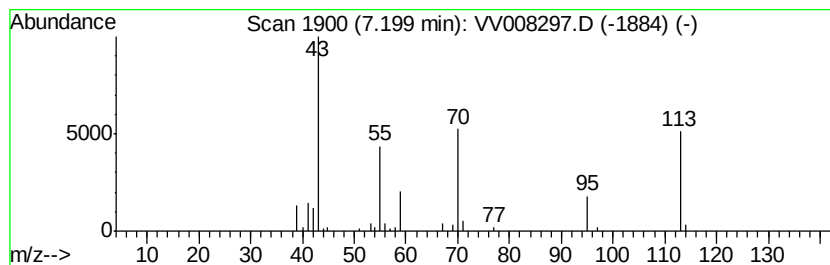
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Peak Number 3 Furan, tetrahydro-2,2,5,5-t... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	1.31 ug/L	53425	1,4-Difluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Furan, tetrahydro-2,2,5,5-tetram...	128	C8H16O	015045-43-9	83
2		Furan, tetrahydro-2,2,4,4-tetram...	128	C8H16O	003358-28-9	59
3		2,5-Hexanediol, 2,5-dimethyl-	146	C8H18O2	000110-03-2	53
4		Furan, tetrahydro-2,2,4,4-tetram...	128	C8H16O	003358-28-9	45
5		Furan, tetrahydro-2,2,5,5-tetram...	128	C8H16O	015045-43-9	40



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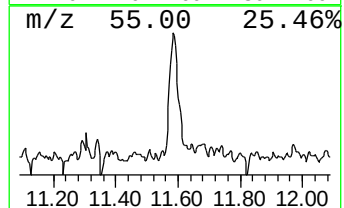
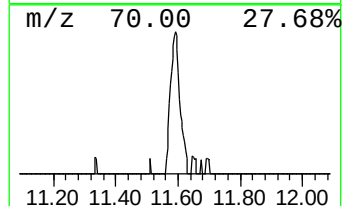
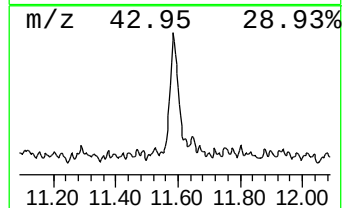
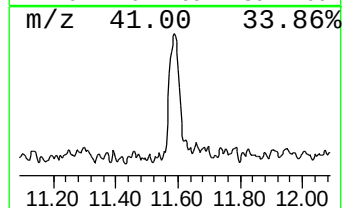
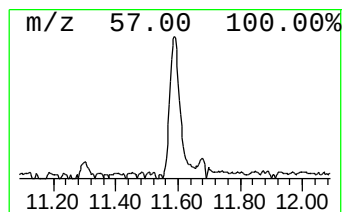
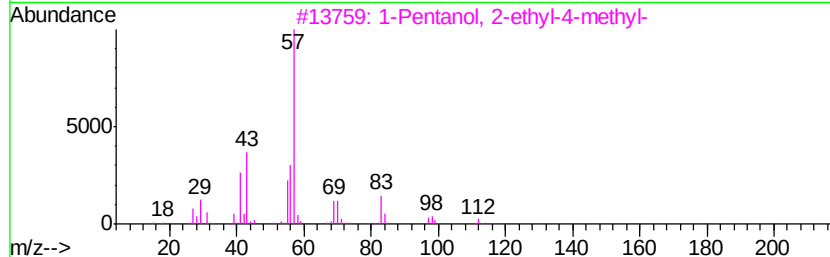
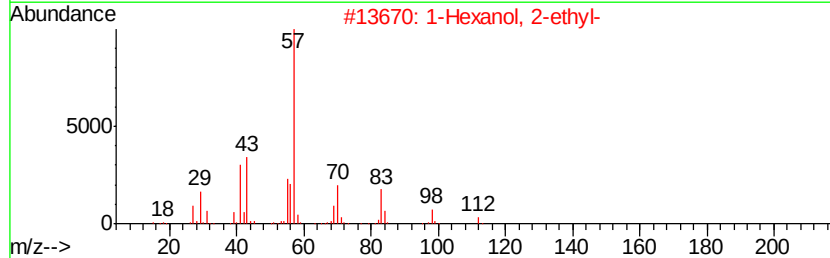
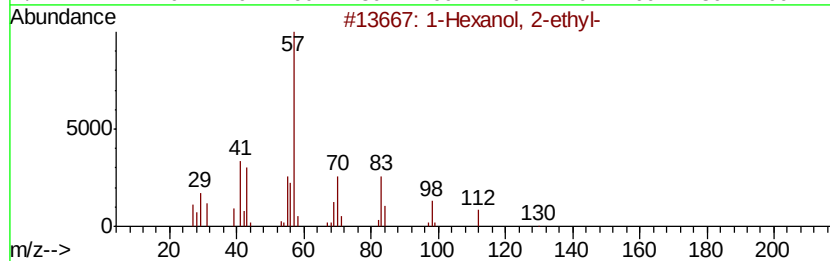
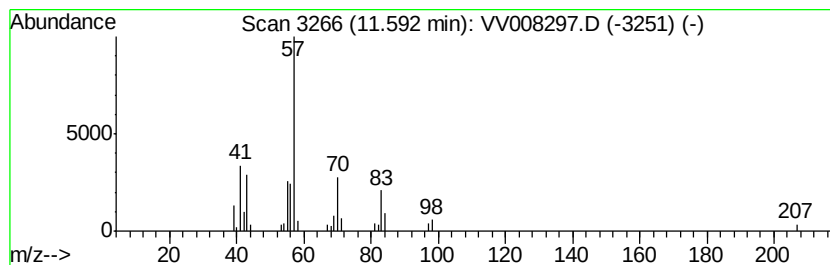
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Peak Number 4 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.59	0.75 ug/L	29861	1,4-Dichlorobenzene-d4	11.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Hexanol, 2-ethyl-	130 C8H18O	000104-76-7	59
2	1-Hexanol, 2-ethyl-	130 C8H18O	000104-76-7	59
3	1-Pentanol, 2-ethyl-4-methyl-	130 C8H18O	000106-67-2	53
4	1-Hexanol, 2-ethyl-	130 C8H18O	000104-76-7	50
5	2-Propyl-1-pentanol	130 C8H18O	058175-57-8	50



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propanal, 2-methyl-	3.00	0.7	ug/L	27009	1	5.67	204694	5.0
Furan, tetrahydro...	7.20	1.3	ug/L	53425	1	5.67	204694	5.0
1-Hexanol, 2-ethyl-	11.59	0.8	ug/L	29861	3	11.30	198834	5.0