

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX110118\
 Data File : VX005724.D
 Acq On : 01 Nov 2018 12:57
 Operator : JC/MD
 Sample : J5199-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OK-01-103018

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

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_X\METHOD\82X103118W.M

Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.611	66	75	78	rBV4	9503	18040	2.48%	0.406%
2	2.446	206	212	220	rVB	6007	10930	1.50%	0.246%
3	2.861	274	280	288	rVB2	5287	10204	1.40%	0.229%
4	5.501	702	713	726	rBV	90707	253888	34.85%	5.710%
5	5.665	729	740	757	rVV2	151269	424067	58.22%	9.537%
6	6.074	796	807	823	rBV2	106259	283849	38.97%	6.384%
7	6.458	859	870	886	rBV	72748	188051	25.82%	4.229%
8	6.860	926	936	951	rBV	239766	579098	79.50%	13.023%
9	7.884	1097	1104	1119	rBV	89602	158612	21.77%	3.567%
10	8.713	1233	1240	1250	rBV	452315	728438	100.00%	16.382%
11	9.408	1350	1354	1358	rVB2	8497	10664	1.46%	0.240%
12	9.549	1370	1377	1383	rBV	44874	63054	8.66%	1.418%
13	9.646	1388	1393	1398	rVB	10694	15337	2.11%	0.345%
14	9.975	1441	1447	1458	rBV	49714	66657	9.15%	1.499%
15	10.116	1463	1470	1486	rBV	456449	628895	86.33%	14.143%
16	11.140	1632	1638	1652	rBV	353311	450529	61.85%	10.132%
17	12.079	1786	1792	1802	rBV	457451	556264	76.36%	12.510%

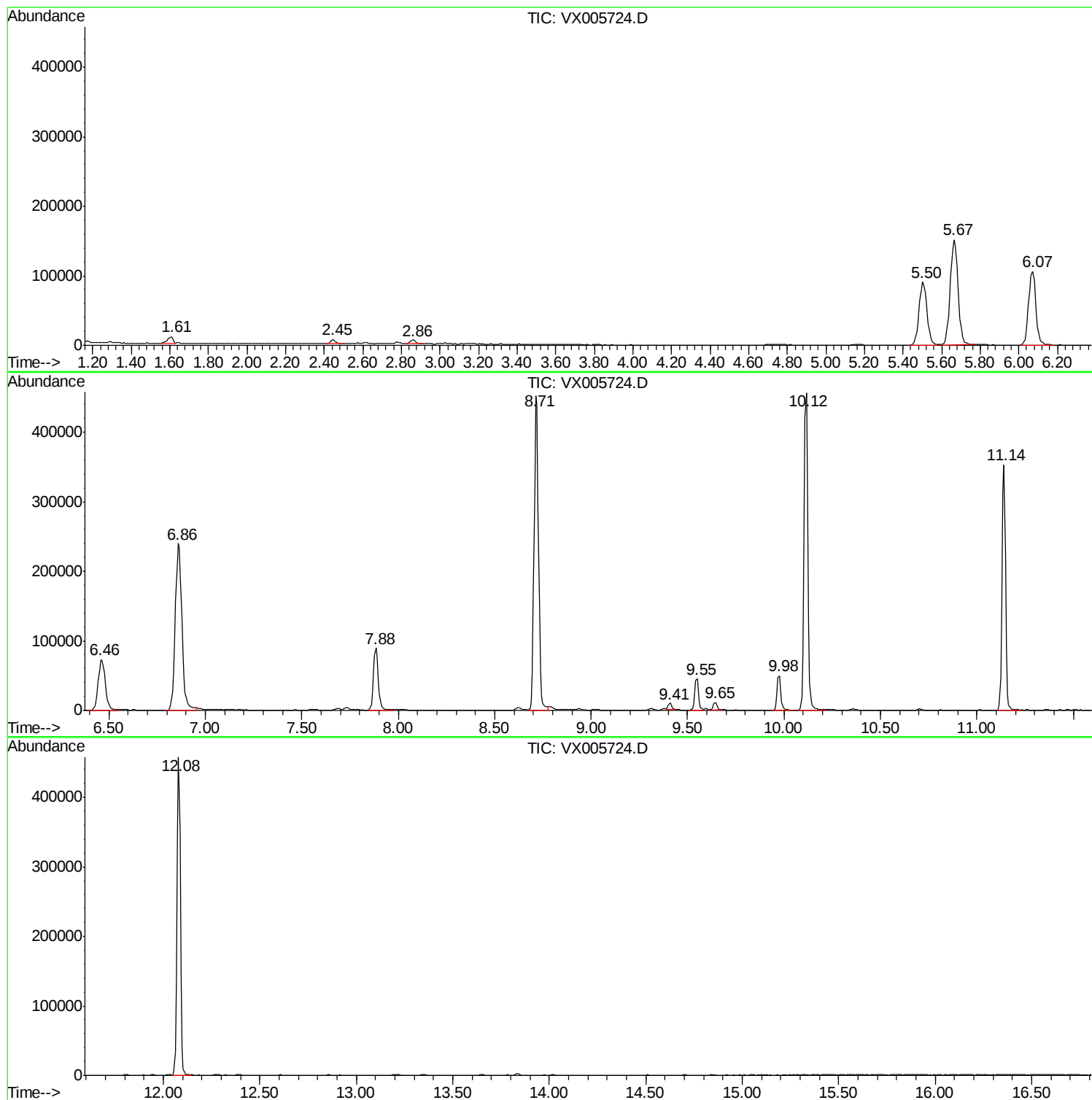
Sum of corrected areas: 4446577

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Instrument :
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ClientSampleId :
OK-01-103018

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X103118W.M
Quant Title : SW846 8260

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



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Instrument :
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ClientSampled :
OK-01-103018

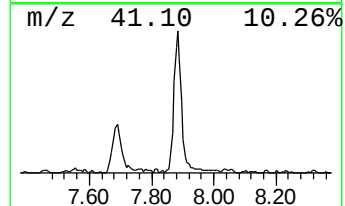
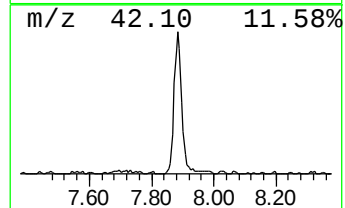
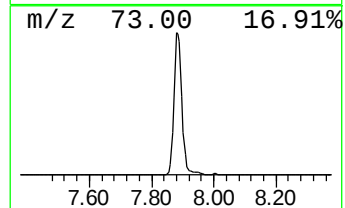
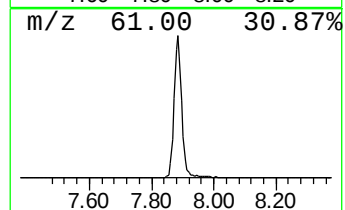
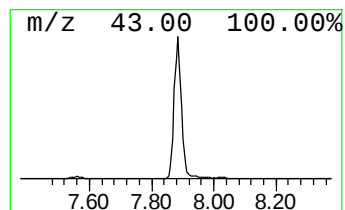
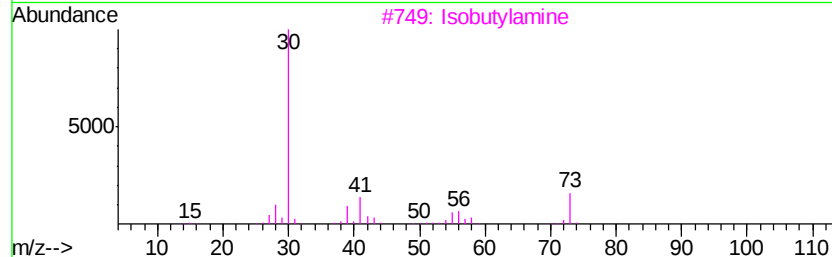
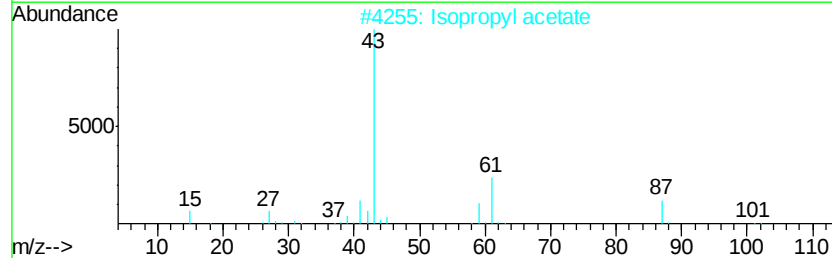
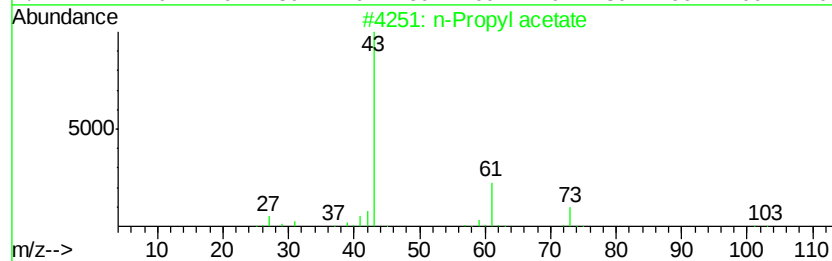
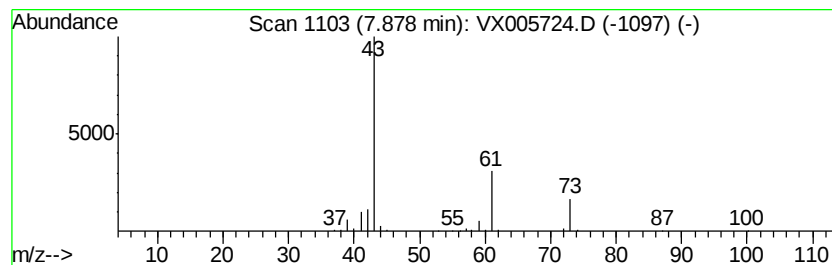
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X103118W.M
Quant Title : SW846 8260

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

Peak Number 1 n-Propyl acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.88	13.69 ug/l	158612	1,4-Difluorobenzene	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Propyl acetate	102	C5H10O2	000109-60-4	83
2		Isopropyl acetate	102	C5H10O2	000108-21-4	36
3		Isobutylamine	73	C4H11N	000078-81-9	9
4		Isobutane	58	C4H10	000075-28-5	7
5		1-Butanamine	73	C4H11N	000109-73-9	5



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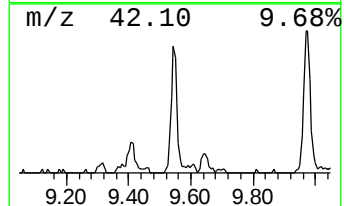
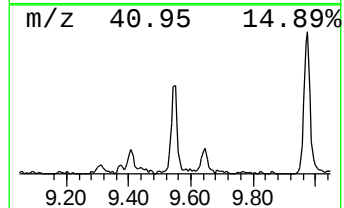
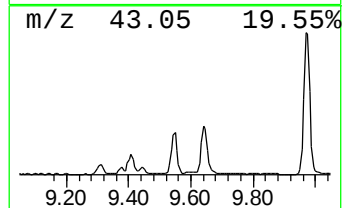
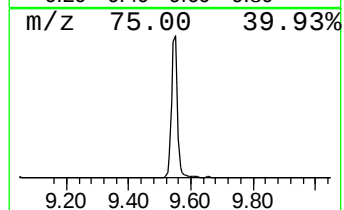
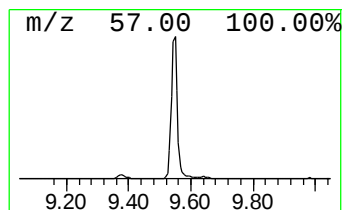
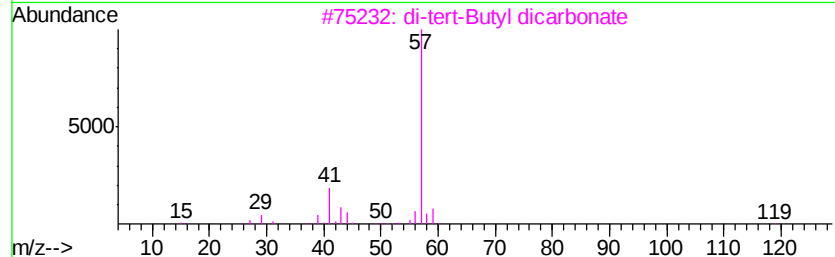
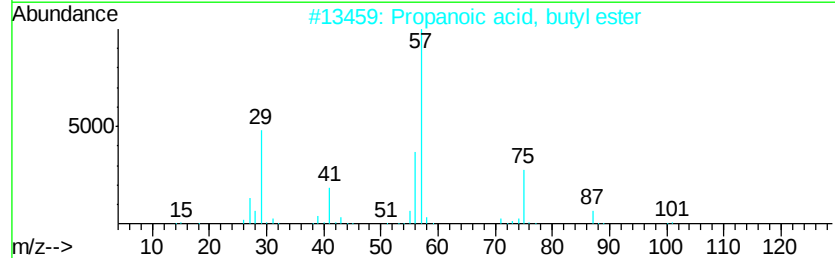
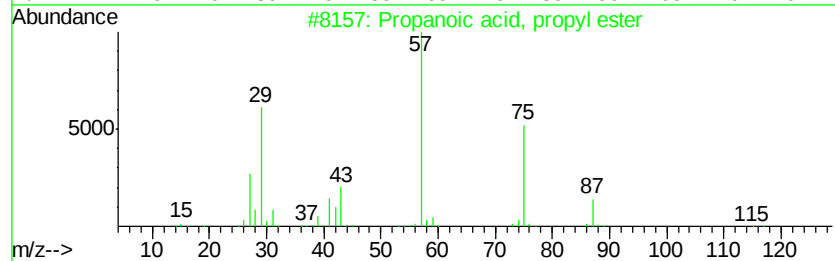
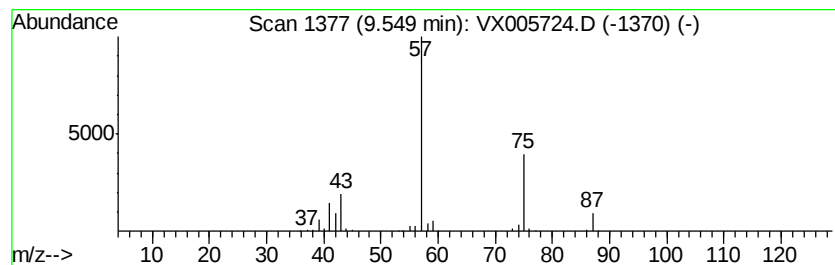
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 Peak Number 2 Propanoic acid, propyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.55	5.01 ug/l	63054	Chlorobenzene-d5	10.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
2		Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	50
3		di-tert-Butyl dicarbonate	218	C10H18O5	024424-99-5	10
4		S-(-)-2-Methylbutylamine	87	C5H13N	020626-52-2	9
5		1-Pentanamine	87	C5H13N	000110-58-7	9



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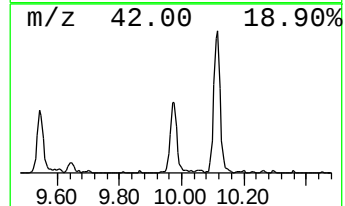
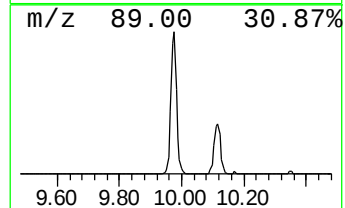
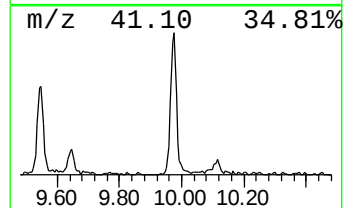
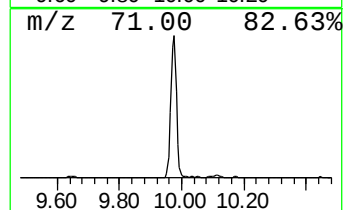
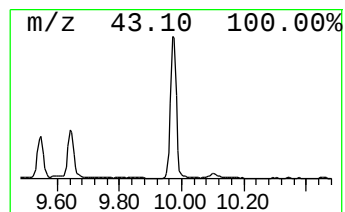
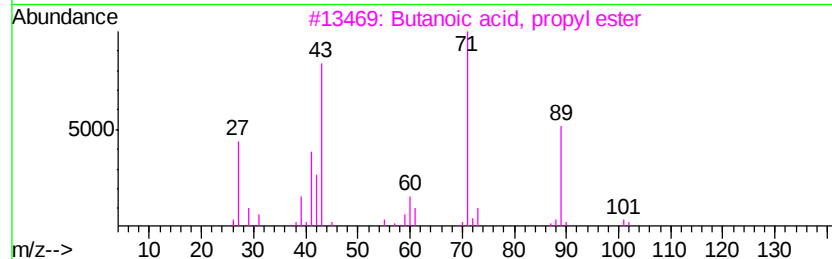
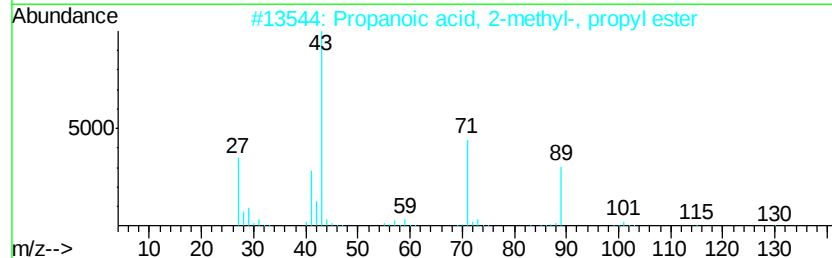
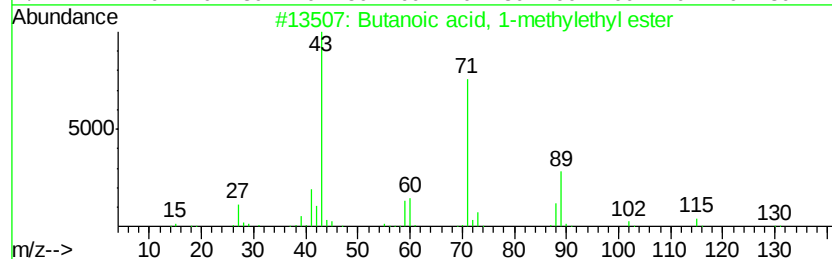
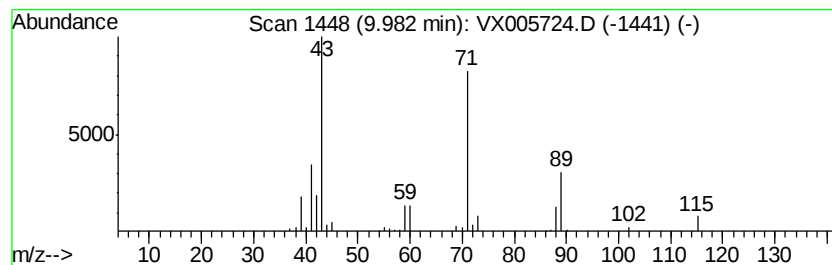
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Peak Number 3 Butanoic acid, 1-methylethy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.98	5.30 ug/l	66657	Chlorobenzene-d5	10.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
2		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72
3		Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	72
4		Butanoic acid, pentyl ester	158	C9H18O2	000540-18-1	64
5		Propanoic acid, 2-methyl-, penty...	158	C9H18O2	002445-72-9	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Propyl acetate	7.88	13.7	ug/l	158612	2	6.87	579098	50.0
Propanoic acid, p...	9.55	5.0	ug/l	63054	3	10.12	628895	50.0
Butanoic acid, 1-...	9.98	5.3	ug/l	66657	3	10.12	628895	50.0