

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006901.D
 Acq On : 09 Sep 2021 15:32
 Operator : CG/JU
 Sample : M3552-04 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AH3

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

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:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006901.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.264	27	30	39	rVB	20538	37732	0.18%	0.023%
2	3.370	44	48	51	rVV	59316	82253	0.40%	0.051%
3	3.411	51	55	66	rVB	589868	778417	3.80%	0.480%
4	3.723	104	108	112	rVB	25697	34227	0.17%	0.021%
5	3.787	116	119	125	rBV	419725	651374	3.18%	0.402%
6	3.840	125	128	137	rVB	113771	175318	0.86%	0.108%
7	4.028	155	160	167	rVB	38425	63155	0.31%	0.039%
8	4.117	169	175	180	rVB2	28793	48183	0.24%	0.030%
9	4.387	216	221	226	rBV	29681	37265	0.18%	0.023%
10	4.452	228	232	237	rVV	62050	77886	0.38%	0.048%
11	4.564	243	251	258	rBV	896229	1167452	5.70%	0.720%
12	4.628	258	262	269	rVB	80122	109830	0.54%	0.068%
13	5.022	323	329	339	rVB	530504	722010	3.52%	0.445%
14	5.887	472	476	481	rVB2	19452	26955	0.13%	0.017%
15	7.093	669	681	691	rBV	558415	902011	4.40%	0.556%
16	7.269	704	711	727	rBV	2276013	3602295	17.58%	2.222%
17	7.469	737	745	757	rBV	2007155	3158807	15.42%	1.948%
18	7.940	816	825	832	rBV	1719411	2681362	13.09%	1.654%
19	8.005	832	836	844	rVB	60556	90241	0.44%	0.056%
20	8.175	860	865	872	rVB5	15434	27083	0.13%	0.017%
21	8.640	936	944	951	rBV	1746792	2788753	13.61%	1.720%
22	8.693	951	953	959	rVV	29710	49797	0.24%	0.031%
23	8.763	959	965	971	rVB	64614	110615	0.54%	0.068%
24	9.099	1015	1022	1038	rBV	2596518	4255289	20.77%	2.625%
25	9.605	1100	1108	1110	rBV4	26658	51151	0.25%	0.032%
26	9.652	1114	1116	1124	rVB4	31119	48253	0.24%	0.030%
27	9.822	1136	1145	1153	rBV	1949857	3176670	15.50%	1.959%
28	10.046	1179	1183	1190	rVB6	13507	23857	0.12%	0.015%
29	10.358	1228	1236	1244	rBV	2991691	5037768	24.59%	3.107%
30	10.422	1244	1247	1252	rVB6	18522	32680	0.16%	0.020%
31	10.522	1259	1264	1275	rBV	147311	253335	1.24%	0.156%
32	10.657	1283	1287	1292	rVB5	12720	22153	0.11%	0.014%
33	10.740	1293	1301	1314	rVB	2543178	4247079	20.73%	2.620%
34	10.875	1316	1324	1335	rBV	2512664	4323282	21.10%	2.667%
35	11.287	1386	1394	1399	rBV2	31122	58786	0.29%	0.036%
36	11.346	1399	1404	1408	rVV2	37385	62889	0.31%	0.039%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 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:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
 Title : SVOA CALIBRATION

37	12.322	1565	1570	1578	rVB	67177	106791	0.52%	0.066%
38	12.528	1599	1605	1613	rBV3	19730	46131	0.23%	0.028%
39	13.534	1771	1776	1784	rVB4	18667	34079	0.17%	0.021%
40	13.975	1841	1851	1857	rBV	6082318	8391687	40.95%	5.176%
41	14.016	1857	1858	1867	rVB3	38382	47813	0.23%	0.029%
42	14.257	1892	1899	1914	rVB	7040664	10460144	51.05%	6.452%
43	14.563	1940	1951	1962	rBV	3946213	5855563	28.58%	3.612%
44	14.746	1974	1982	1988	rBV	518687	800162	3.90%	0.494%
45	14.975	2017	2021	2026	rVB7	15637	24278	0.12%	0.015%
46	15.328	2077	2081	2087	rBV2	30890	48478	0.24%	0.030%
47	15.557	2113	2120	2132	rBV	9400399	13376982	65.28%	8.251%
48	15.669	2132	2139	2153	rVV	2847397	3996625	19.50%	2.465%
49	15.793	2156	2160	2166	rVB2	67228	105857	0.52%	0.065%
50	16.245	2231	2237	2242	rVB5	22224	41837	0.20%	0.026%
51	16.340	2248	2253	2261	rBV2	33981	55416	0.27%	0.034%
52	16.992	2359	2364	2368	rBV5	14459	23966	0.12%	0.015%
53	17.098	2372	2382	2386	rBV3	24234	55586	0.27%	0.034%
54	17.310	2412	2418	2428	rBV	4609058	6834134	33.35%	4.216%
55	17.416	2428	2436	2449	rBV	10705322	15599164	76.13%	9.622%
56	17.987	2529	2533	2537	rVB4	36915	49424	0.24%	0.030%
57	18.198	2564	2569	2578	rBV	324028	498244	2.43%	0.307%
58	19.057	2712	2715	2718	rVB5	19512	20579	0.10%	0.013%
59	19.222	2738	2743	2749	rBV2	135161	199846	0.98%	0.123%
60	19.286	2749	2754	2762	rBV	162122	250711	1.22%	0.155%
61	19.428	2774	2778	2787	rBV	239403	355230	1.73%	0.219%
62	19.645	2809	2815	2821	rBV	13950929	18815592	91.82%	11.606%
63	21.104	3060	3063	3070	rBV3	134574	196560	0.96%	0.121%
64	21.392	3107	3112	3125	rBV	5903492	7889507	38.50%	4.866%
65	23.686	3493	3502	3509	rBV2	9872238	20491048	100.00%	12.640%
66	23.839	3520	3528	3540	rVB2	3987556	8431283	41.15%	5.201%

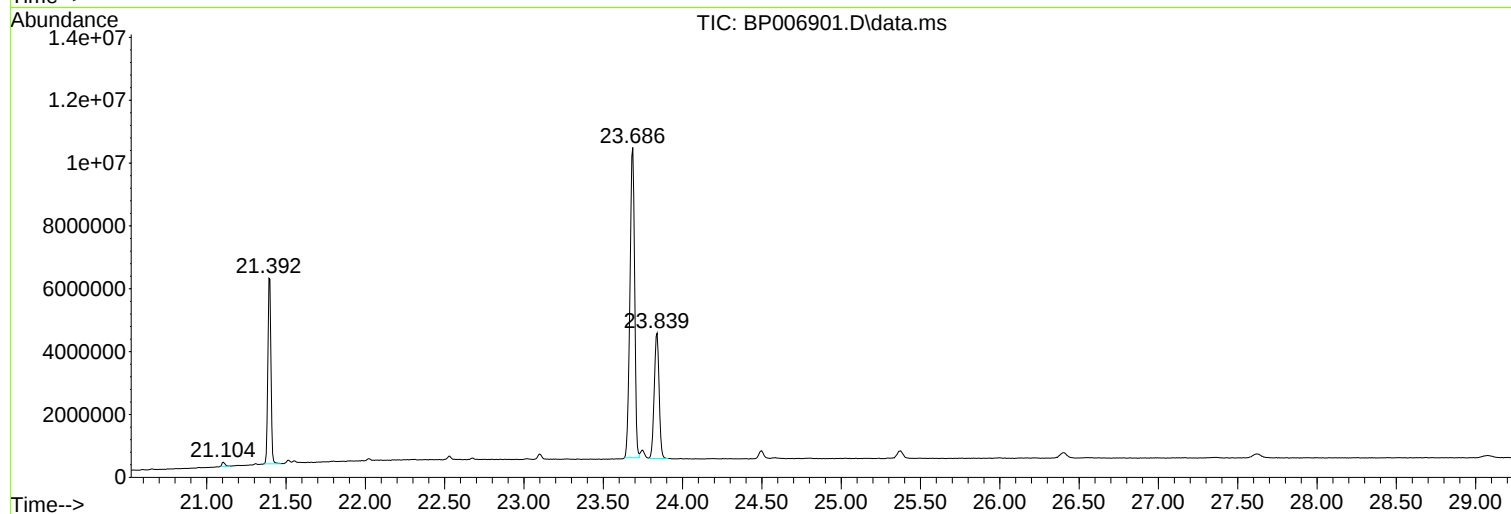
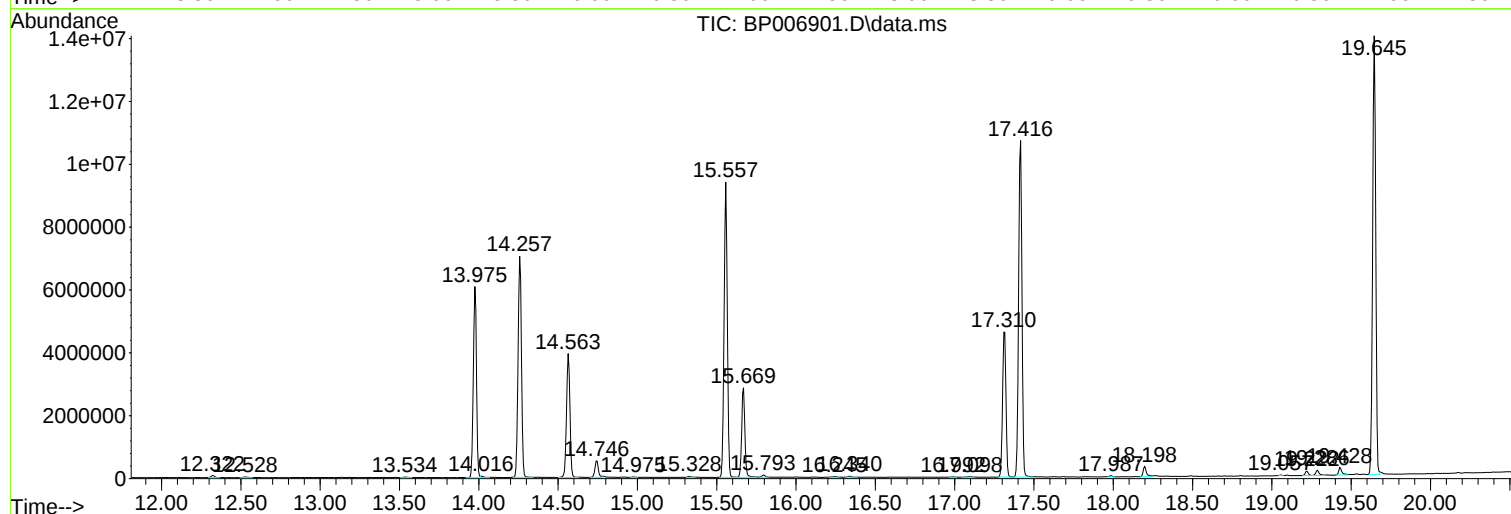
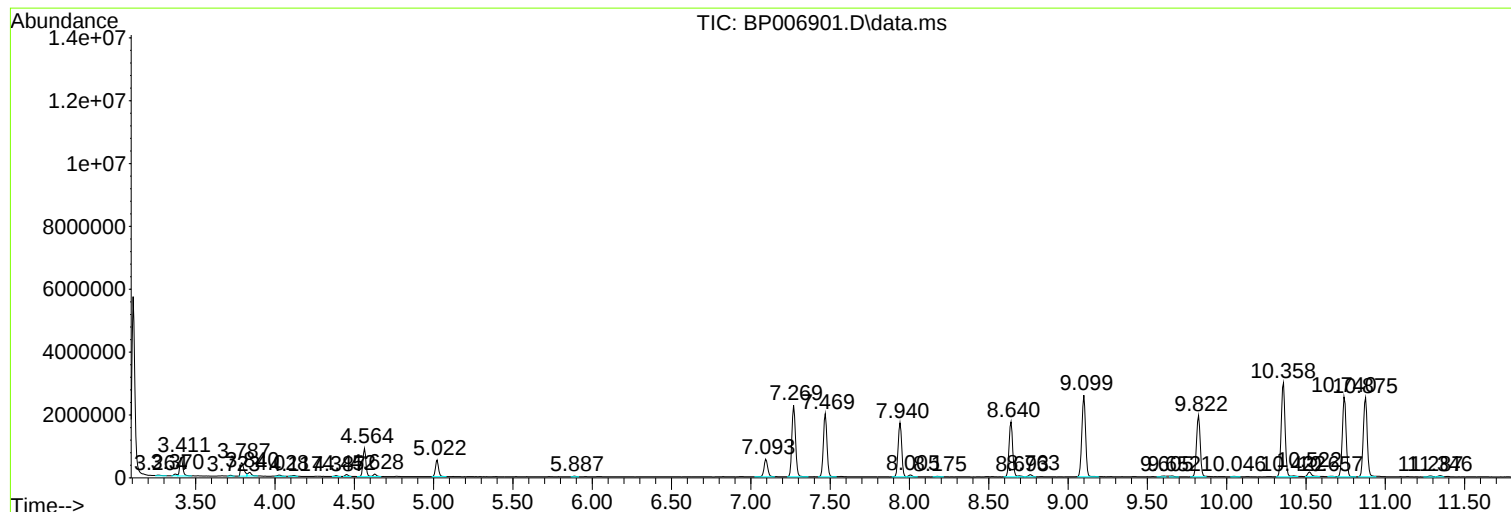
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ALS Vial : 11 Sample Multiplier: 1

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP006901.M
Quant Title : SVOA CALIBRATION

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



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Sample : M3552-04 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AH3

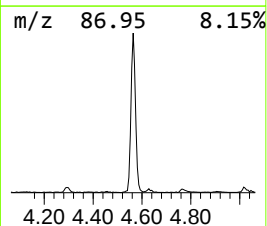
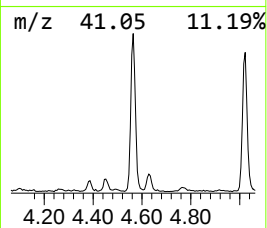
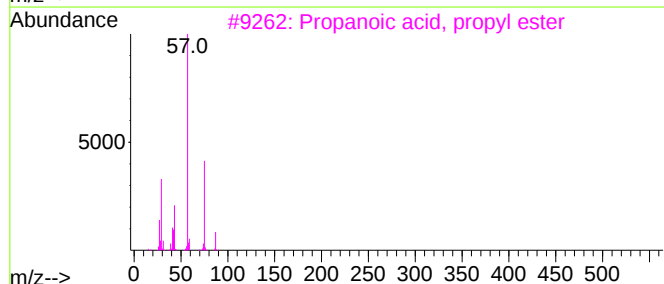
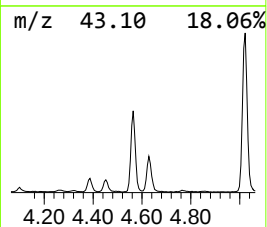
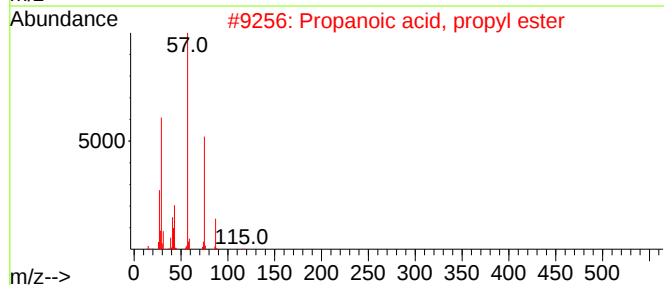
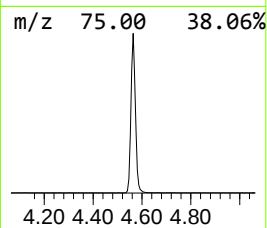
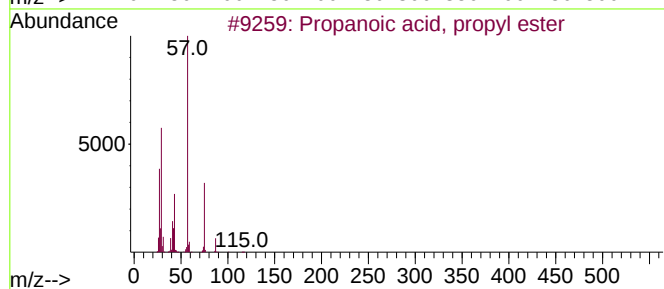
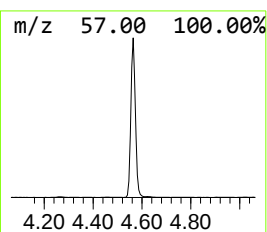
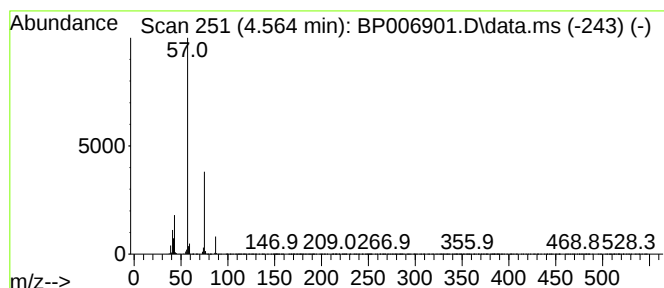
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propanoic acid, propyl ester Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.564	8.71 ng/ul	1167450	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
2			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	74
3			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	72
4			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	64
5			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	47



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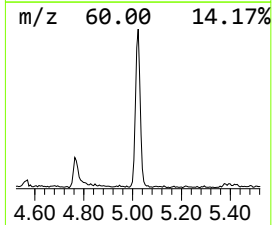
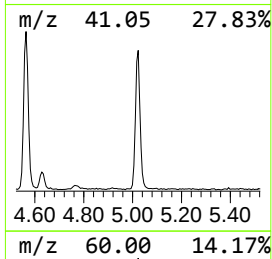
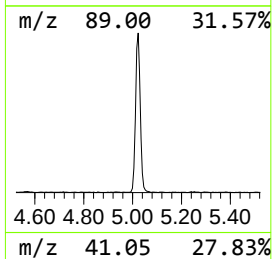
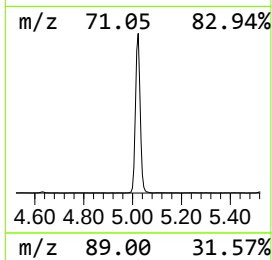
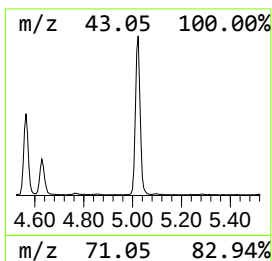
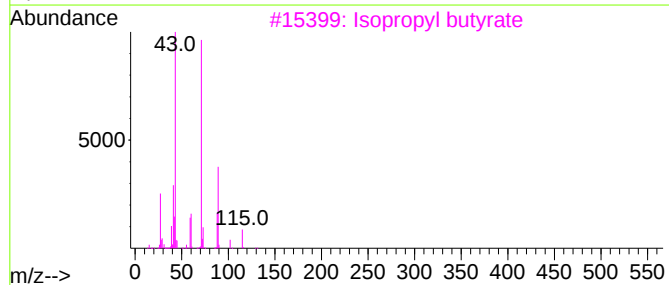
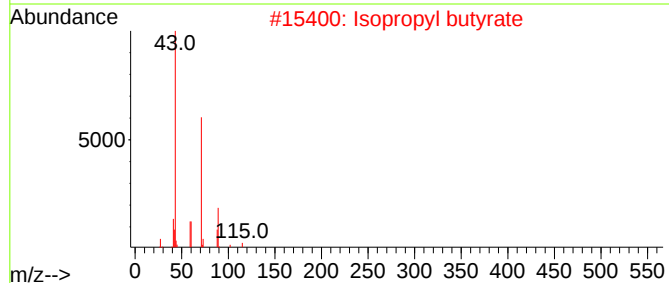
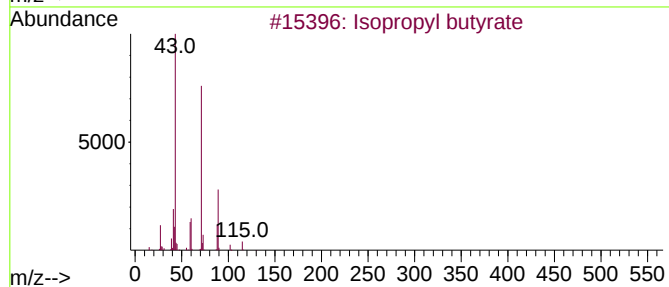
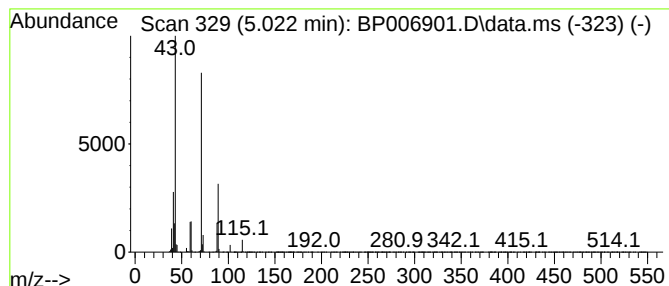
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Isopropyl butyrate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.022	5.39 ng/ul	722010	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl butyrate	130	C7H14O2	000638-11-9	94
2			Isopropyl butyrate	130	C7H14O2	000638-11-9	90
3			Isopropyl butyrate	130	C7H14O2	000638-11-9	90
4			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	78
5			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propanoic acid,...	4.564	8.7	ng/ul	1167450	1	7.940	2681360	20.0
Isopropyl butyrate	5.022	5.4	ng/ul	722010	1	7.940	2681360	20.0