

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007711.D
 Acq On : 27 Oct 2021 10:19
 Operator : CG/JU
 Sample : M4303-16
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P027-TW001-01

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-BP102521.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007711.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.364	7	13	20	rBV	49829	69300	1.24%	0.147%
2	3.787	82	85	94	rBV	229496	346205	6.19%	0.734%
3	4.022	120	125	129	rBV	30816	46631	0.83%	0.099%
4	4.117	137	141	145	rBV2	10297	13838	0.25%	0.029%
5	4.211	155	157	162	rVB3	12676	15843	0.28%	0.034%
6	5.034	294	297	301	rVB2	5669	7256	0.13%	0.015%
7	5.217	324	328	334	rVB5	11942	19687	0.35%	0.042%
8	6.022	462	465	473	rVB9	3435	6308	0.11%	0.013%
9	6.316	510	515	520	rBV4	11780	19021	0.34%	0.040%
10	7.099	641	648	658	rVB	181871	301263	5.39%	0.638%
11	7.263	669	676	686	rVB	616602	973474	17.42%	2.063%
12	7.463	702	710	717	rBV	638604	991930	17.75%	2.102%
13	7.534	717	722	730	rVV	148777	233937	4.19%	0.496%
14	7.934	781	790	798	rBV	645185	1001250	17.92%	2.122%
15	8.005	798	802	806	rVB6	6263	10542	0.19%	0.022%
16	8.181	827	832	836	rBV8	5035	8505	0.15%	0.018%
17	8.640	901	910	920	rBV	500491	798130	14.28%	1.691%
18	8.758	926	930	934	rVB5	5398	7806	0.14%	0.017%
19	9.093	978	987	997	rBV	696623	1162840	20.81%	2.464%
20	9.557	1062	1066	1071	rVB7	5891	9486	0.17%	0.020%
21	9.816	1101	1110	1122	rBV	526925	867728	15.53%	1.839%
22	10.357	1195	1202	1214	rBV	879725	1418201	25.38%	3.005%
23	10.440	1214	1216	1223	rVB8	4578	7023	0.13%	0.015%
24	10.522	1223	1230	1240	rBV	104992	201964	3.61%	0.428%
25	10.740	1258	1267	1278	rVV	837609	1425246	25.50%	3.020%
26	10.869	1281	1289	1301	rBV	780746	1351039	24.17%	2.863%
27	10.946	1301	1302	1311	rVB9	4782	7473	0.13%	0.016%
28	11.393	1372	1378	1383	rBV8	3897	7586	0.14%	0.016%
29	11.563	1400	1407	1411	rBV10	6146	10923	0.20%	0.023%
30	12.322	1529	1536	1543	rBV2	17740	31665	0.57%	0.067%
31	12.499	1560	1566	1567	rBV5	5967	7861	0.14%	0.017%
32	12.946	1637	1642	1647	rBV8	3646	8190	0.15%	0.017%
33	13.093	1659	1667	1671	rVB8	7459	15750	0.28%	0.033%
34	13.193	1678	1684	1688	rBV8	4688	6828	0.12%	0.014%
35	13.481	1729	1733	1737	rBV7	4087	6997	0.13%	0.015%
36	13.546	1740	1744	1750	rVB8	8383	12628	0.23%	0.027%

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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-BP102521.M
 Title : SVOA CALIBRATION

37	13.981	1811	1818	1831	rBV	1716358	2362663	42.28%	5.006%
38	14.263	1855	1866	1876	rBV	1842541	2711941	48.52%	5.747%
39	14.569	1911	1918	1931	rVB	1307258	1954107	34.96%	4.141%
40	14.763	1943	1951	1964	rBV	138635	240932	4.31%	0.511%
41	14.987	1984	1989	1996	rBV	7159	13351	0.24%	0.028%
42	15.392	2052	2058	2062	rBV3	11570	19292	0.35%	0.041%
43	15.563	2080	2087	2101	rBV	2507929	3667780	65.63%	7.772%
44	15.681	2101	2107	2120	rBV	661556	897165	16.05%	1.901%
45	15.804	2122	2128	2133	rVV3	16836	26439	0.47%	0.056%
46	16.263	2196	2206	2211	rBV	144305	222976	3.99%	0.472%
47	16.351	2218	2221	2227	rVB6	6138	10191	0.18%	0.022%
48	17.010	2327	2333	2334	rBV6	5046	7611	0.14%	0.016%
49	17.116	2346	2351	2356	rBV2	7467	13907	0.25%	0.029%
50	17.263	2371	2376	2379	rBV6	4387	6444	0.12%	0.014%
51	17.328	2379	2387	2397	rVV	1616915	2354633	42.13%	4.989%
52	17.428	2397	2404	2416	rVV	2954633	4235553	75.79%	8.975%
53	17.522	2416	2420	2425	rVB7	16394	25651	0.46%	0.054%
54	17.622	2434	2437	2440	rBV3	7522	8566	0.15%	0.018%
55	18.010	2498	2503	2508	rVB4	25498	38837	0.69%	0.082%
56	18.222	2534	2539	2543	rBV6	29006	46882	0.84%	0.099%
57	18.292	2547	2551	2556	rVB	8448	14931	0.27%	0.032%
58	19.239	2708	2712	2719	rBV3	43568	60492	1.08%	0.128%
59	19.304	2719	2723	2728	rBV	44366	61426	1.10%	0.130%
60	19.457	2746	2749	2753	rBV5	12981	16344	0.29%	0.035%
61	19.604	2772	2774	2777	rVB3	14357	11918	0.21%	0.025%
62	19.663	2778	2784	2792	rBV	3983429	5243236	93.82%	11.110%
63	21.133	3031	3034	3038	rBV2	59081	76812	1.37%	0.163%
64	21.416	3077	3082	3088	rBV	2195540	2841881	50.85%	6.022%
65	23.704	3464	3471	3482	rBV2	2759313	5588753	100.00%	11.843%
66	23.863	3491	3498	3507	rVB2	1452692	2980924	53.34%	6.317%

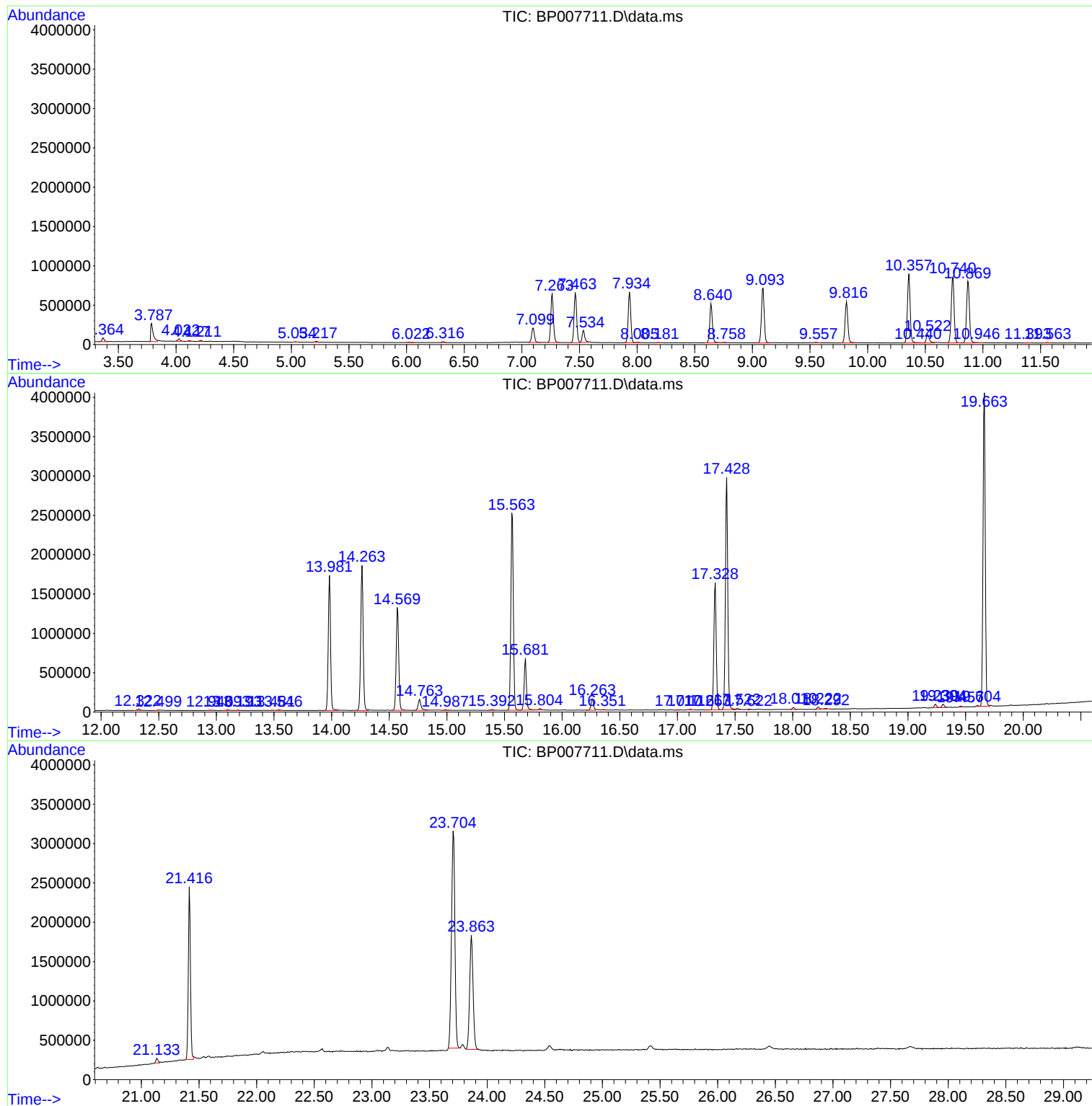
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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



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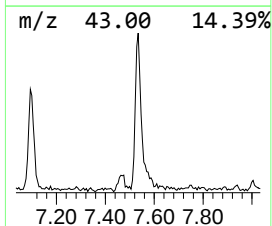
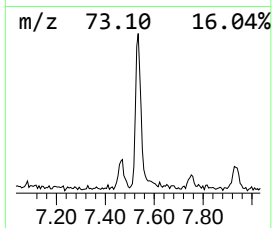
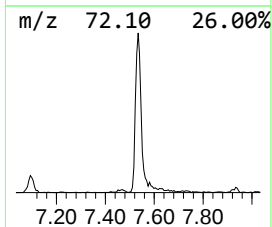
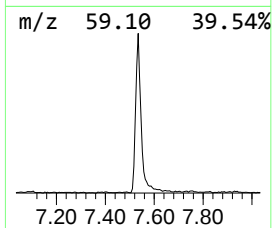
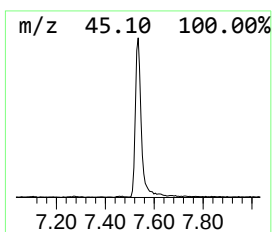
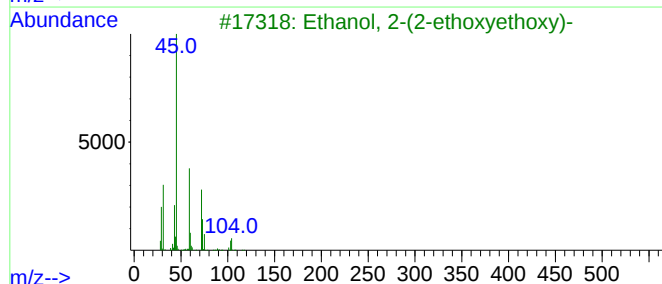
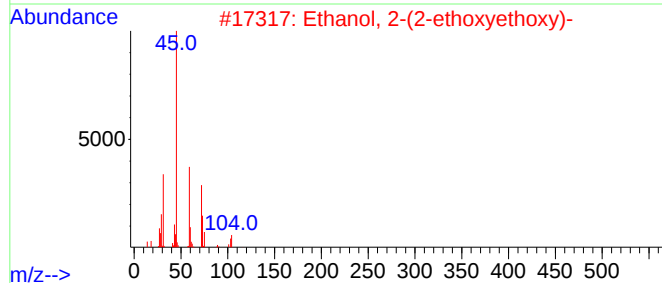
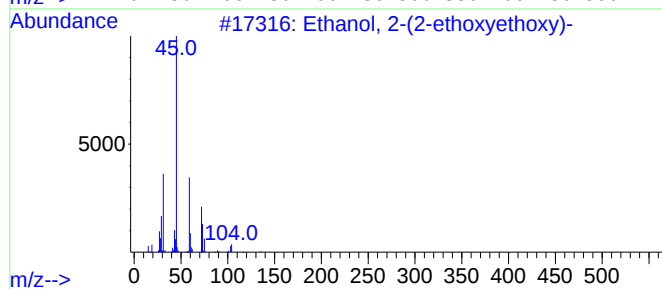
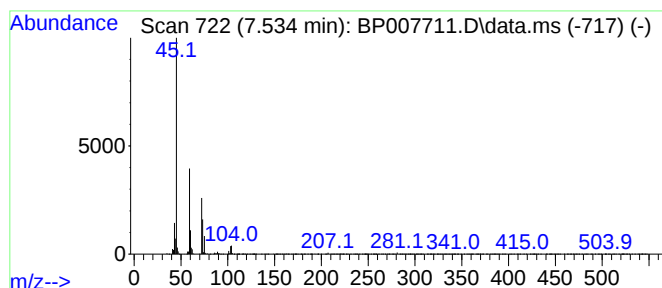
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.534	4.67 ng/ul	233937	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	86
3			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
4			Diethyl carbitol	162	C8H18O3	000112-36-7	72
5			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	72



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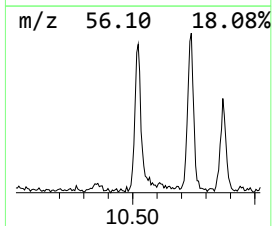
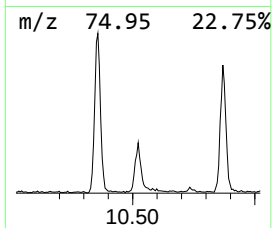
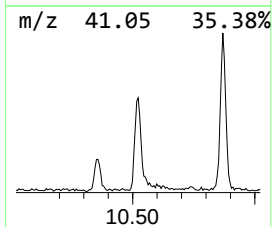
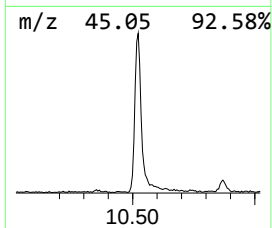
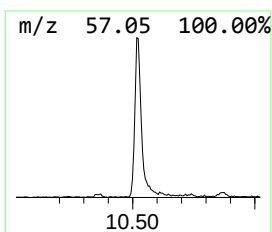
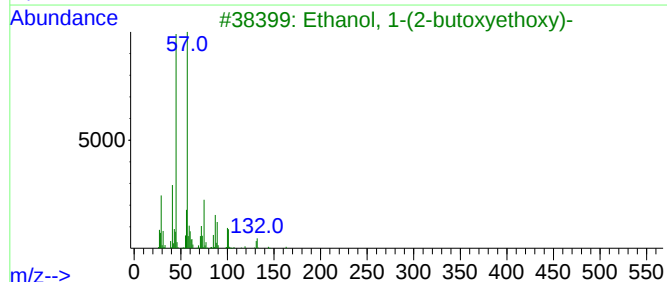
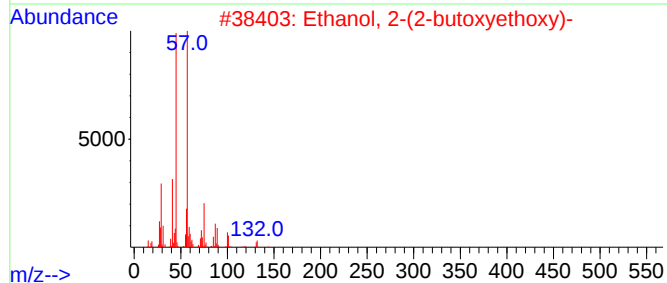
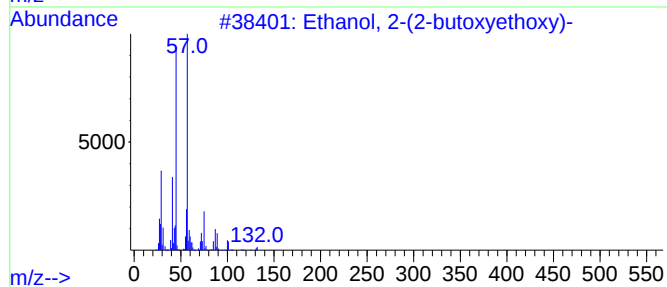
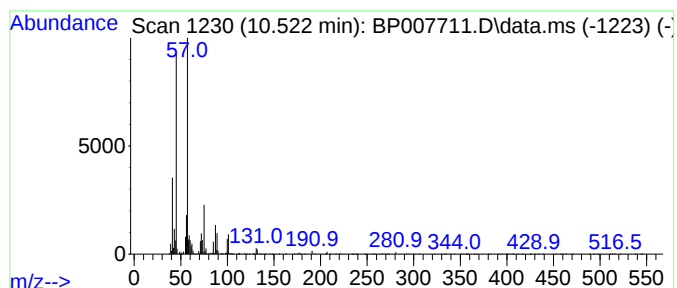
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.522	2.83 ng/ul	201964	Naphthalene-d8	10.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-e...	7.534	4.7	ng/u1	233937	1	7.934	1001250	20.0
Ethanol, 2-(2-b...	10.522	2.8	ng/u1	201964	2	10.740	1425250	20.0