

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007690.D
 Acq On : 26 Oct 2021 21:52
 Operator : CG/JU
 Sample : M4291-22
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P047-TW001-01

Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : 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: BP007690.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	10	13	20	rVB	36054	48635	1.00%	0.124%
2	3.646	58	61	66	rBV3	7329	12207	0.25%	0.031%
3	3.787	83	85	98	rBV	156607	237739	4.88%	0.608%
4	4.022	121	125	132	rVB2	25590	38987	0.80%	0.100%
5	4.117	137	141	145	rVB2	12373	15043	0.31%	0.038%
6	4.216	154	158	162	rVB4	9287	12865	0.26%	0.033%
7	5.040	294	298	303	rBV4	11768	21755	0.45%	0.056%
8	6.016	460	464	471	rBV4	6846	11916	0.24%	0.030%
9	6.316	511	515	522	rVB4	10633	17365	0.36%	0.044%
10	6.934	616	620	621	rBV4	6791	7192	0.15%	0.018%
11	7.099	642	648	657	rVB	129491	213031	4.38%	0.545%
12	7.263	669	676	687	rBV	438848	700871	14.40%	1.792%
13	7.463	702	710	720	rBV	457598	710058	14.59%	1.815%
14	7.563	725	727	732	rVB6	3516	5580	0.11%	0.014%
15	7.934	782	790	798	rBV	593090	935442	19.22%	2.392%
16	8.005	800	802	806	rVV5	4866	6493	0.13%	0.017%
17	8.075	810	814	817	rBV6	3513	5042	0.10%	0.013%
18	8.175	825	831	838	rBV7	9978	20645	0.42%	0.053%
19	8.487	881	884	888	rVB6	5355	7458	0.15%	0.019%
20	8.640	899	910	920	rBV	343004	571917	11.75%	1.462%
21	9.087	977	986	997	rBV	482418	810291	16.65%	2.072%
22	9.257	1011	1015	1016	rBV3	4206	5349	0.11%	0.014%
23	9.552	1057	1065	1071	rVB10	6484	14416	0.30%	0.037%
24	9.816	1103	1110	1120	rBV	365779	600363	12.33%	1.535%
25	10.357	1194	1202	1213	rBV	615747	1019116	20.94%	2.605%
26	10.522	1224	1230	1241	rBV	101774	180830	3.71%	0.462%
27	10.740	1259	1267	1275	rBV	753681	1288391	26.47%	3.294%
28	10.869	1281	1289	1299	rBV	537228	928193	19.07%	2.373%
29	12.263	1519	1526	1533	rBV	157596	261157	5.36%	0.668%
30	12.328	1533	1537	1543	rVB2	16104	28666	0.59%	0.073%
31	13.193	1677	1684	1688	rVB9	3762	8546	0.18%	0.022%
32	13.481	1728	1733	1737	rVV8	3422	5106	0.10%	0.013%
33	13.545	1741	1744	1750	rVB8	3550	5995	0.12%	0.015%
34	13.981	1811	1818	1834	rBV	1221648	1665207	34.21%	4.257%
35	14.263	1856	1866	1875	rBV	1290409	1919520	39.43%	4.907%
36	14.569	1910	1918	1930	rBV	1191342	1772366	36.41%	4.531%

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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	14.763	1940	1951	1963	rBV	98839	183615	3.77%	0.469%
38	14.987	1985	1989	1996	rBV9	10394	15298	0.31%	0.039%
39	15.563	2080	2087	2098	rBV	1775166	2675621	54.96%	6.840%
40	15.681	2100	2107	2123	rBV	503813	722026	14.83%	1.846%
41	15.804	2124	2128	2133	rVB	11650	18731	0.38%	0.048%
42	16.257	2199	2205	2210	rBV4	11212	24346	0.50%	0.062%
43	16.357	2216	2222	2227	rVB10	5718	9956	0.20%	0.025%
44	17.104	2345	2349	2354	rBV5	8602	15619	0.32%	0.040%
45	17.328	2380	2387	2397	rBV	1513296	2208565	45.37%	5.646%
46	17.428	2397	2404	2416	rVV	2333620	3482431	71.54%	8.903%
47	17.516	2417	2419	2427	rVB9	11990	22114	0.45%	0.057%
48	17.886	2476	2482	2484	rBV6	4042	6357	0.13%	0.016%
49	18.010	2498	2503	2508	rVB	20885	26583	0.55%	0.068%
50	18.222	2534	2539	2545	rBV2	42650	72459	1.49%	0.185%
51	19.239	2707	2712	2716	rBV3	37009	54291	1.12%	0.139%
52	19.304	2719	2723	2731	rVB	35841	56663	1.16%	0.145%
53	19.451	2745	2748	2754	rBV6	16755	29899	0.61%	0.076%
54	19.598	2771	2773	2777	rBV4	15388	17159	0.35%	0.044%
55	19.663	2778	2784	2797	rVV	3250911	4457323	91.57%	11.395%
56	20.563	2934	2937	2943	rVB2	60638	68106	1.40%	0.174%
57	21.133	3031	3034	3042	rBV5	93418	138898	2.85%	0.355%
58	21.416	3077	3082	3093	rBV	2158002	2844530	58.43%	7.272%
59	23.704	3464	3471	3481	rVB2	2478059	4867915	100.00%	12.445%
60	23.863	3491	3498	3508	rVB2	1446100	2982871	61.28%	7.626%

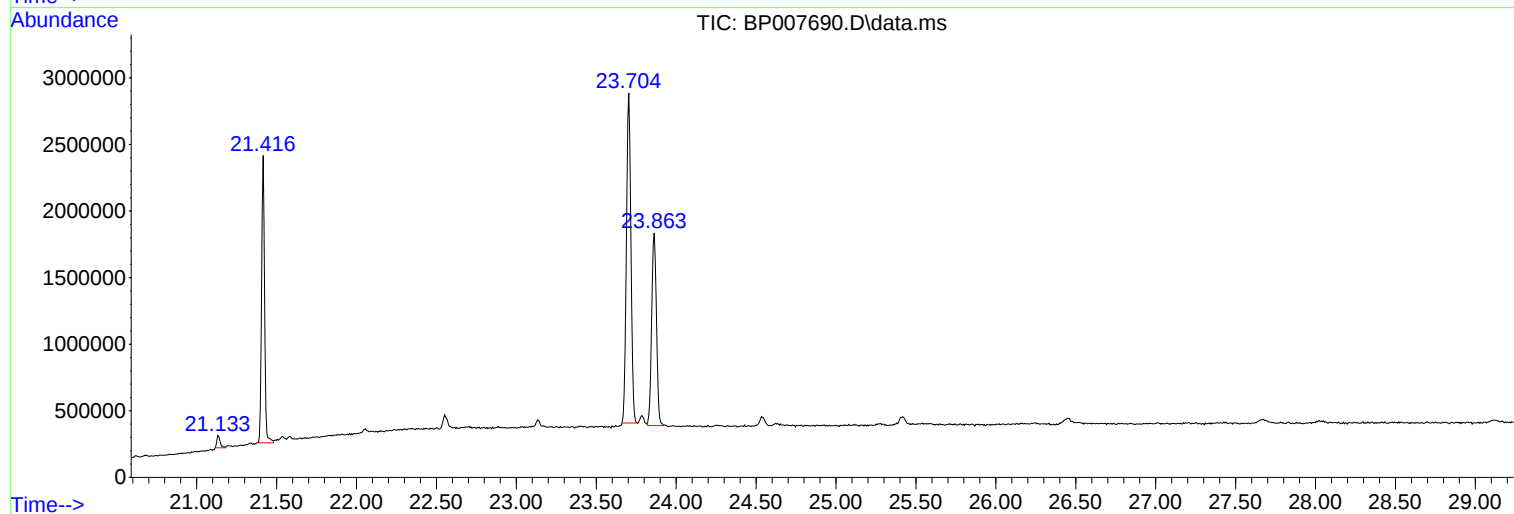
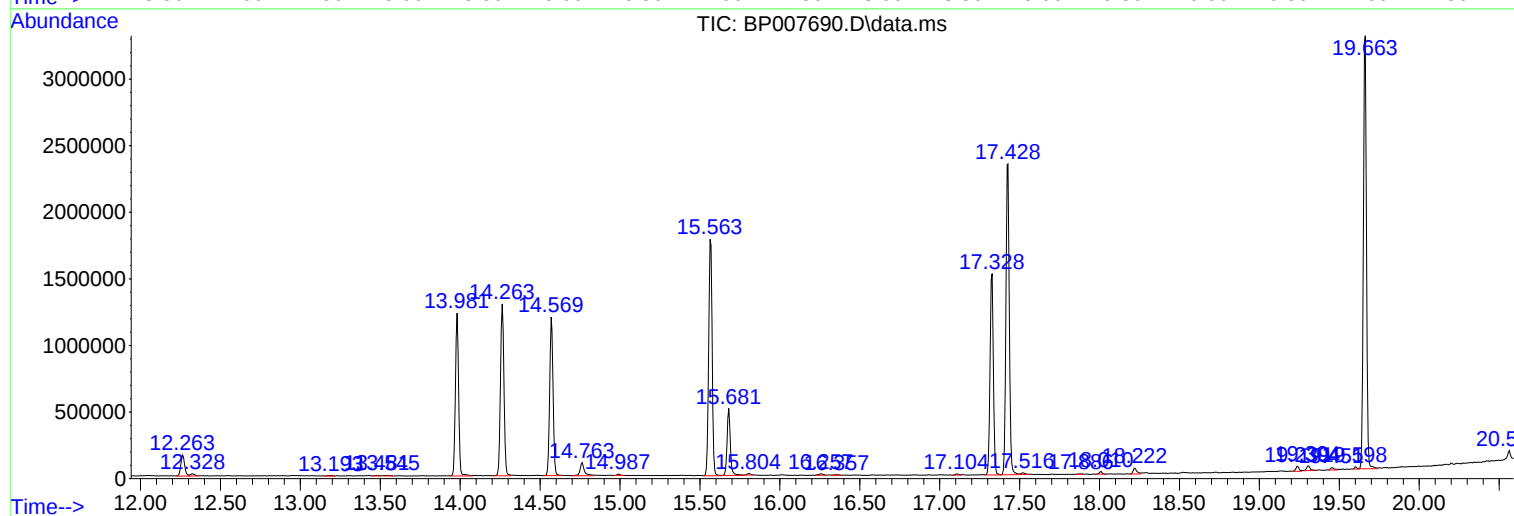
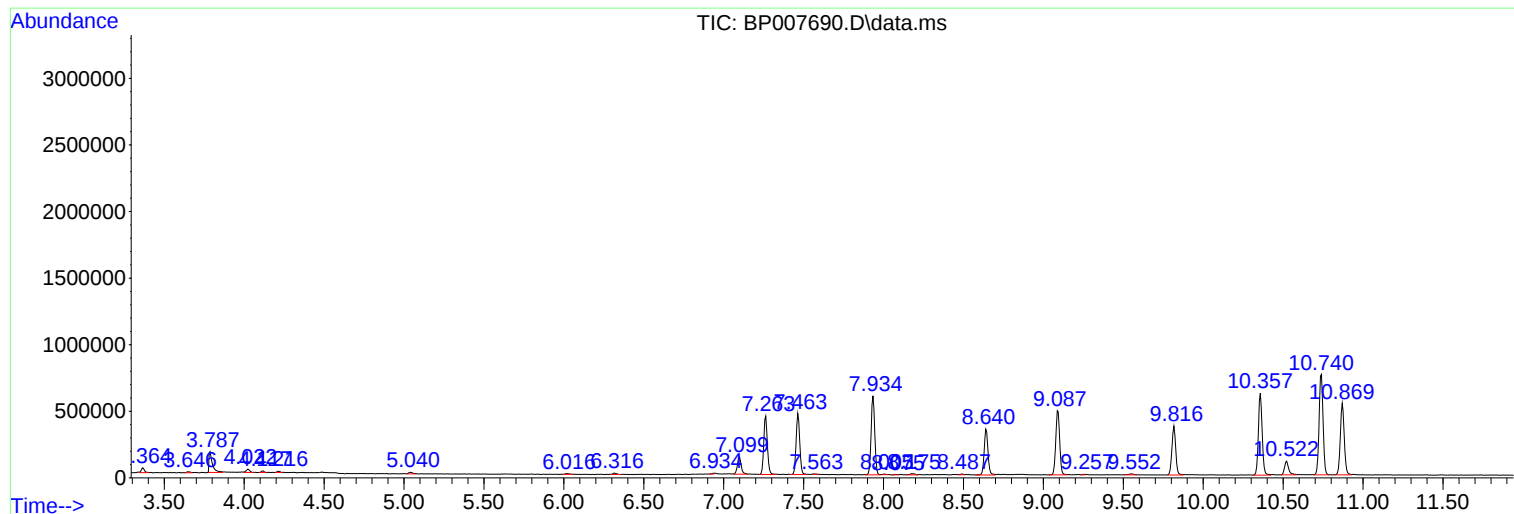
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 48 Sample Multiplier: 1

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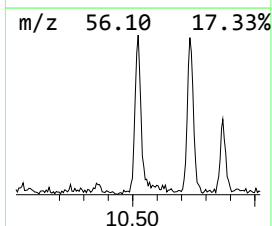
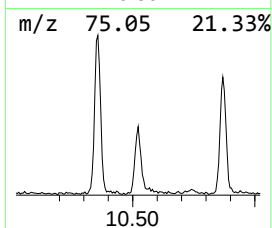
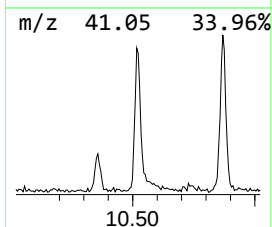
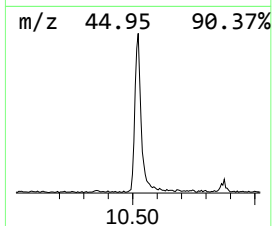
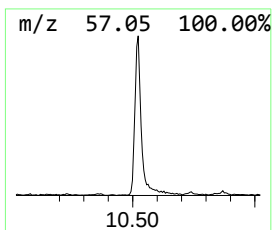
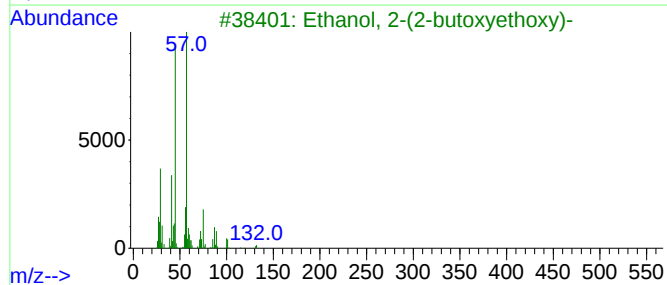
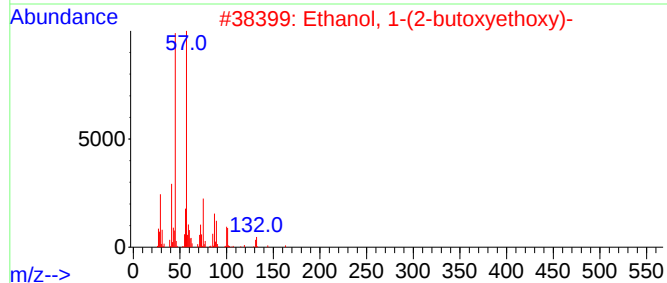
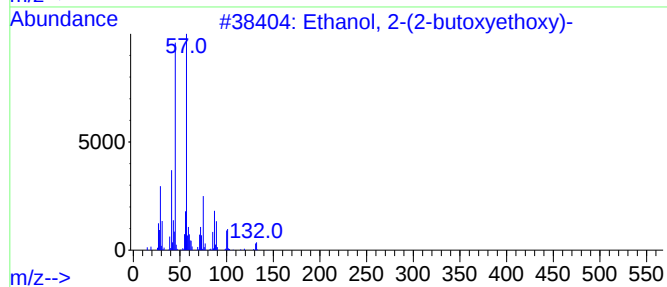
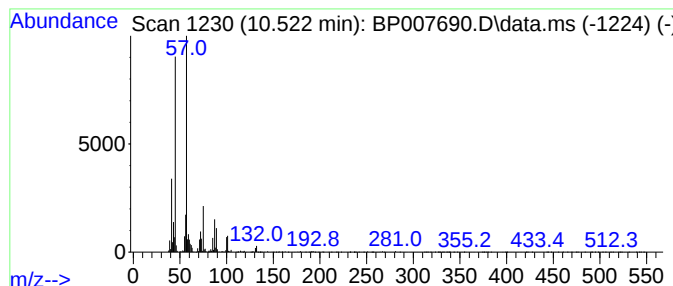
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.522	2.81 ng/ul	180830	Naphthalene-d8	10.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	91
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90



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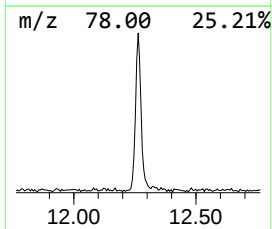
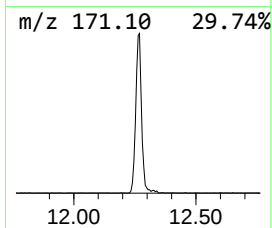
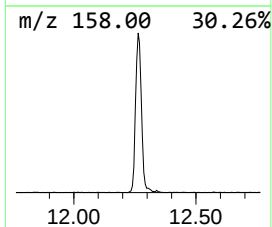
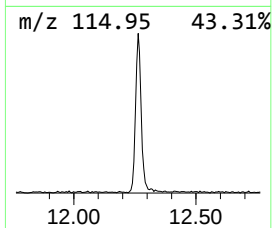
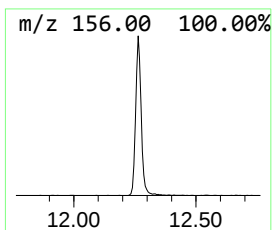
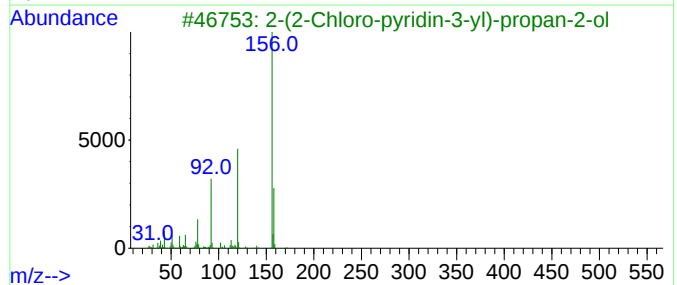
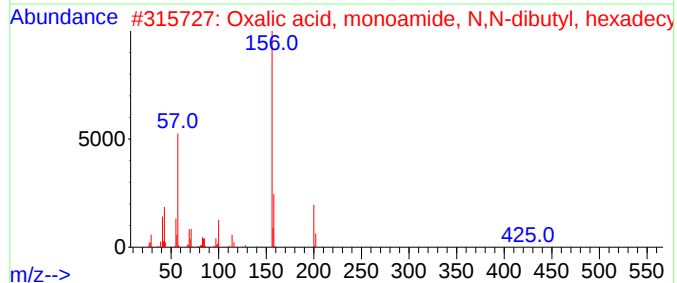
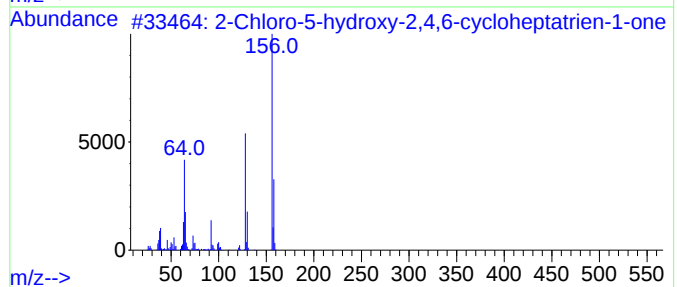
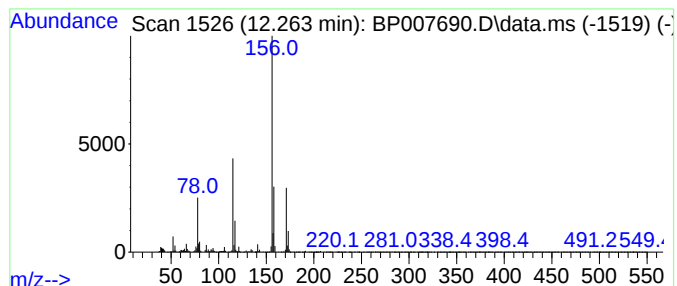
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.263	4.05 ng/ul	261157	Naphthalene-d8	10.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	43
2			Oxalic acid, monoamide, N,N-dibu...	425	C26H51NO3	1000309-46-9	32
3			2-(2-Chloro-pyridin-3-yl)-propan...	171	C8H10ClNO	267003-35-0	25
4			6-Isopropylquinoline	171	C12H13N	000135-79-5	23
5			Tebuthiuron	228	C9H16N4OS	034014-18-1	23



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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-b...	10.522	2.8	ng/u1	180830	2	10.740	1288390	20.0
unknown-01	12.263	4.0	ng/u1	261157	2	10.740	1288390	20.0