

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009898.D
 Acq On : 16 Apr 2022 12:55
 Operator : CG/JU
 Sample : N2421-17
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J77

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-BP040722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP009898.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.117	3	5	18	rVB	356692	498375	7.92%	0.959%
2	3.381	45	50	54	rBV	29265	41677	0.66%	0.080%
3	3.664	94	98	99	rBV3	6108	6341	0.10%	0.012%
4	3.799	118	121	134	rBV	144224	246178	3.91%	0.474%
5	4.128	174	177	184	rVB2	9801	14404	0.23%	0.028%
6	5.264	366	370	377	rVB5	6669	11649	0.19%	0.022%
7	7.111	677	684	693	rBV	187352	307350	4.88%	0.591%
8	7.281	705	713	722	rBV	534994	883344	14.04%	1.700%
9	7.352	724	725	731	rVB	6724	6676	0.11%	0.013%
10	7.481	740	747	756	rBV	568548	921178	14.64%	1.772%
11	7.952	819	827	837	rBV	512127	861549	13.69%	1.658%
12	8.022	837	839	847	rVB5	9090	13371	0.21%	0.026%
13	8.305	884	887	894	rVB9	4091	7727	0.12%	0.015%
14	8.652	938	946	955	rBV	533135	883994	14.05%	1.701%
15	9.110	1016	1024	1039	rBV	756914	1272360	20.22%	2.448%
16	9.281	1047	1053	1058	rVB9	5454	10150	0.16%	0.020%
17	9.569	1096	1102	1108	rBV2	24867	41103	0.65%	0.079%
18	9.840	1137	1148	1159	rBV	614017	1038272	16.50%	1.998%
19	10.375	1231	1239	1252	rBV	915456	1531054	24.33%	2.946%
20	10.552	1263	1269	1275	rBV6	7255	12797	0.20%	0.025%
21	10.763	1297	1305	1314	rVV	787454	1362521	21.65%	2.622%
22	10.893	1317	1327	1340	rVV	814660	1465937	23.30%	2.821%
23	12.063	1520	1526	1533	rBV4	9792	17429	0.28%	0.034%
24	12.199	1545	1549	1554	rVB7	4260	6438	0.10%	0.012%
25	12.346	1569	1574	1580	rVB2	20974	35050	0.56%	0.067%
26	13.322	1735	1740	1745	rVB8	3616	7901	0.13%	0.015%
27	13.422	1753	1757	1763	rVV3	14672	21373	0.34%	0.041%
28	13.516	1768	1773	1777	rVV6	7900	14558	0.23%	0.028%
29	13.563	1777	1781	1788	rVB7	5648	12014	0.19%	0.023%
30	13.993	1847	1854	1861	rBV	2128491	2872292	45.65%	5.527%
31	14.281	1892	1903	1917	rBV	2188232	3208985	51.00%	6.174%
32	14.493	1936	1939	1944	rVB6	5975	7836	0.12%	0.015%
33	14.587	1947	1955	1965	rVV	1347296	2069161	32.88%	3.981%
34	14.763	1978	1985	1991	rBV	226698	350121	5.56%	0.674%
35	14.816	1991	1994	2002	rVB	26993	36579	0.58%	0.070%
36	14.998	2021	2025	2031	rBV8	7605	11144	0.18%	0.021%

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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-BP040722.M
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37	15.228	2059	2064	2071	rVB9	4077	8774	0.14%	0.017%
38	15.393	2086	2092	2097	rBV2	19233	29925	0.48%	0.058%
39	15.581	2115	2124	2136	rBV	2892266	4278818	68.00%	8.233%
40	15.687	2136	2142	2155	rVV	1111059	1507702	23.96%	2.901%
41	15.822	2159	2165	2173	rVB	33594	53736	0.85%	0.103%
42	16.363	2254	2257	2262	rVB5	11866	16226	0.26%	0.031%
43	16.734	2315	2320	2328	rBV10	6367	11320	0.18%	0.022%
44	17.016	2363	2368	2372	rBV7	5977	10014	0.16%	0.019%
45	17.104	2378	2383	2384	rBV5	7023	7431	0.12%	0.014%
46	17.339	2416	2423	2430	rVV	1802549	2607790	41.44%	5.018%
47	17.439	2432	2440	2450	rVV	3605419	5262140	83.62%	10.125%
48	17.528	2451	2455	2458	rVV4	15549	25376	0.40%	0.049%
49	17.557	2458	2460	2466	rVB7	8212	14016	0.22%	0.027%
50	18.010	2533	2537	2540	rBV4	9985	10439	0.17%	0.020%
51	18.098	2548	2552	2557	rBV8	4281	8503	0.14%	0.016%
52	18.216	2568	2572	2579	rBV	105394	150780	2.40%	0.290%
53	18.516	2621	2623	2632	rVB9	6563	9410	0.15%	0.018%
54	19.245	2742	2747	2752	rBV	51638	73417	1.17%	0.141%
55	19.310	2754	2758	2766	rBV	61312	94143	1.50%	0.181%
56	19.451	2779	2782	2787	rVB5	23236	27148	0.43%	0.052%
57	19.598	2804	2807	2811	rVB6	15469	16378	0.26%	0.032%
58	19.669	2812	2819	2831	rVV	4679930	6292671	100.00%	12.108%
59	21.128	3063	3067	3077	rBV	330818	453769	7.21%	0.873%
60	21.416	3111	3116	3126	rBV	2072509	2817092	44.77%	5.420%
61	22.710	3333	3336	3341	rVB	98112	126356	2.01%	0.243%
62	23.721	3500	3508	3516	rBV	2665984	5469687	86.92%	10.524%
63	23.880	3528	3535	3547	rVB2	1176292	2480790	39.42%	4.773%

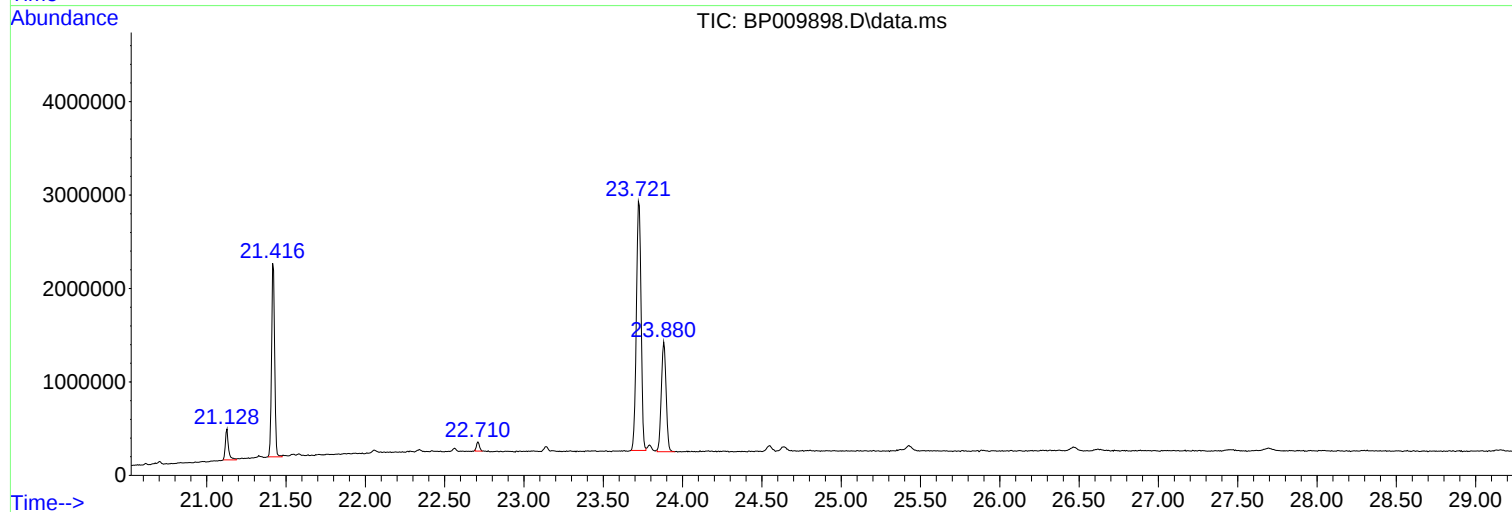
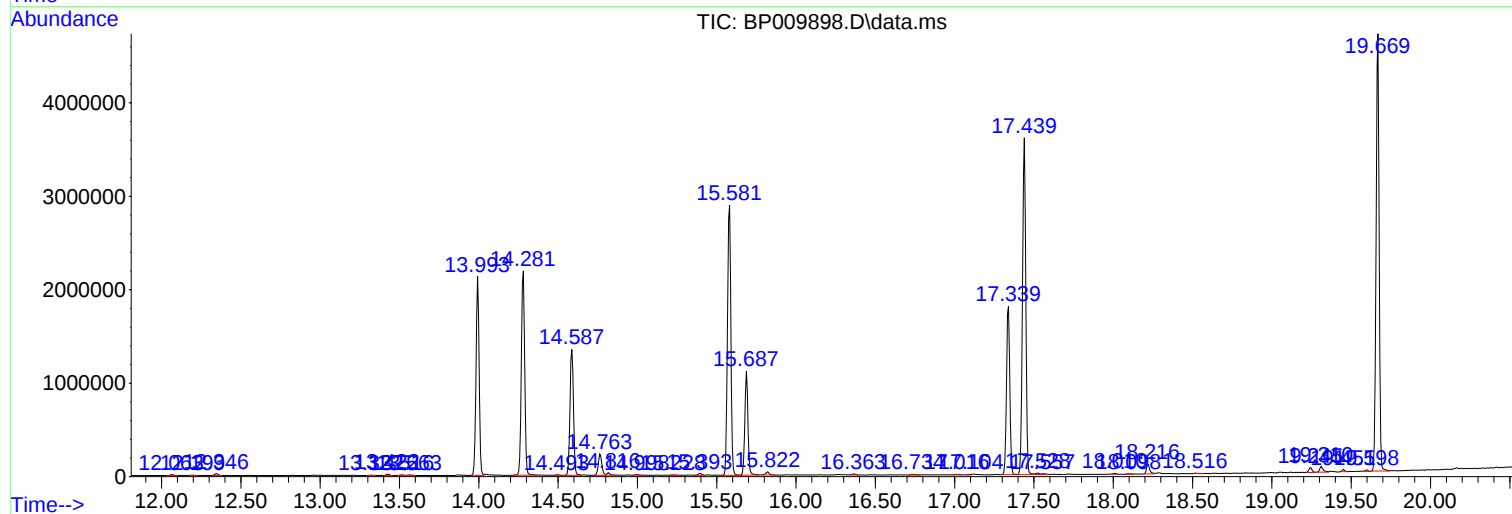
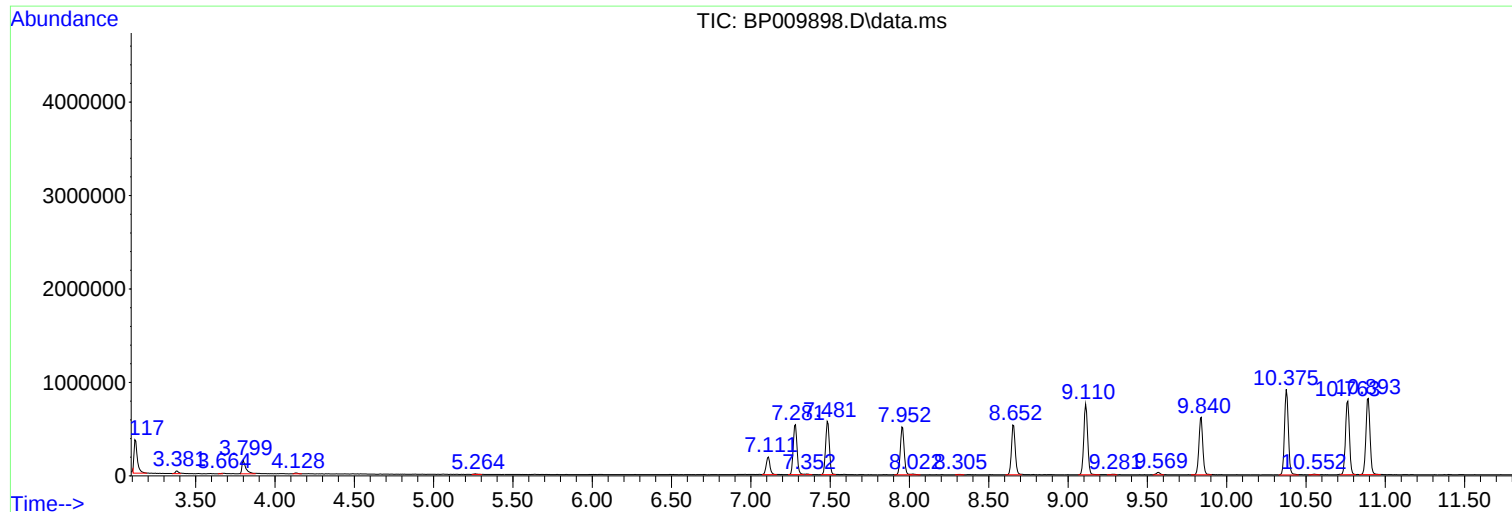
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



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Sample : N2421-17
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Instrument :
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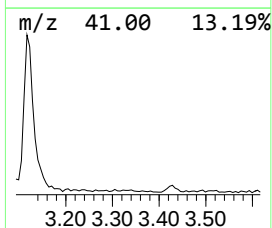
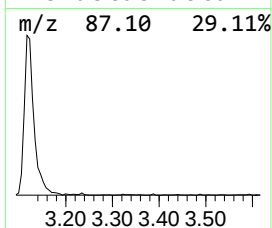
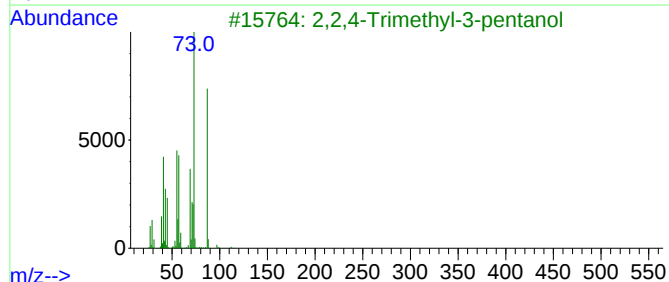
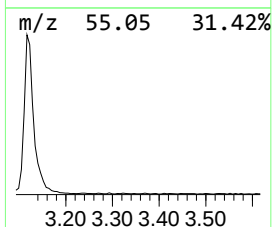
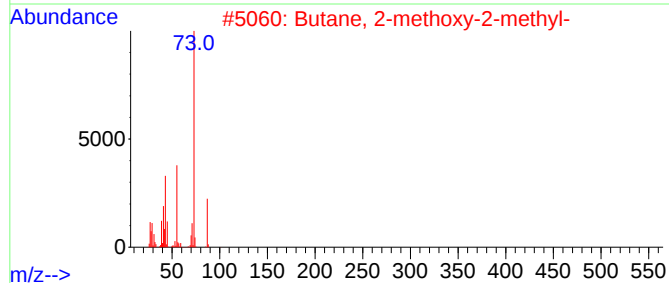
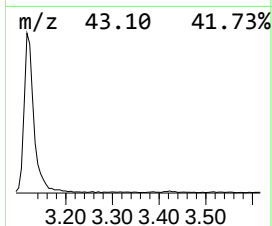
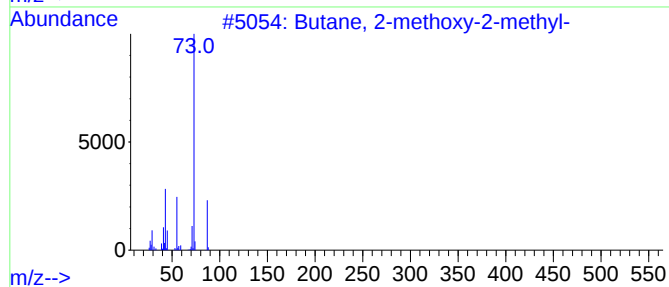
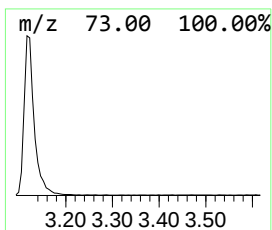
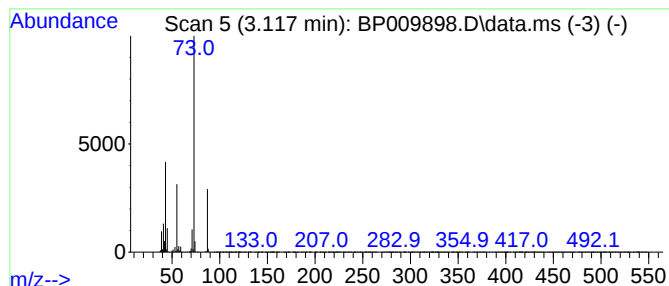
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.117	11.57 ng/ul	498375	1,4-Dichlorobenzene-d4	7.958

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39		
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27		
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25		



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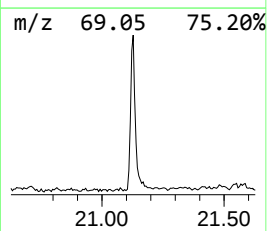
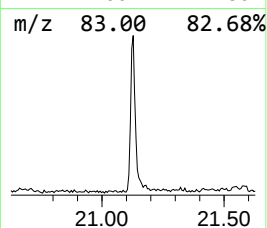
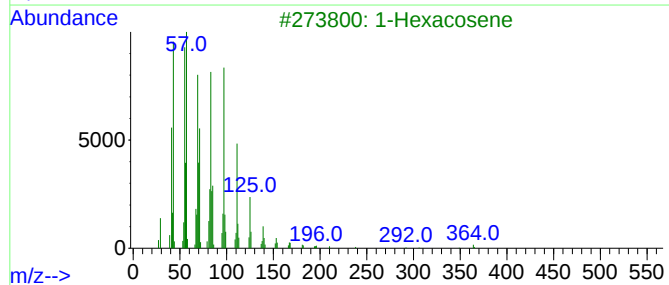
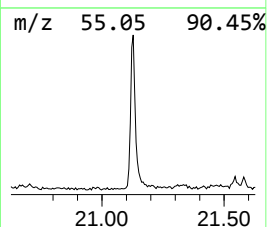
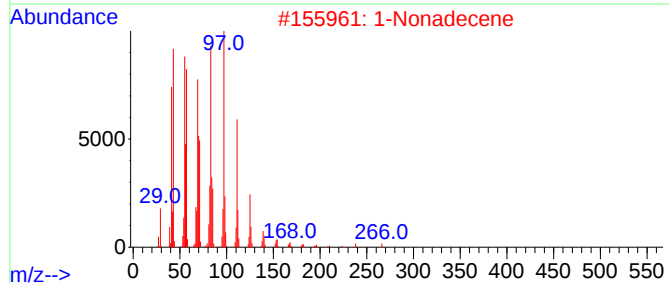
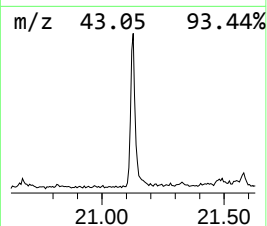
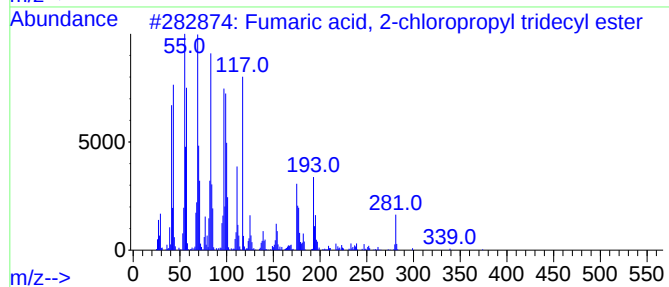
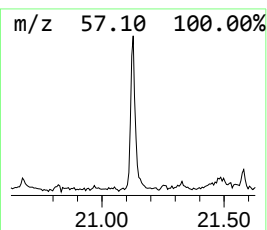
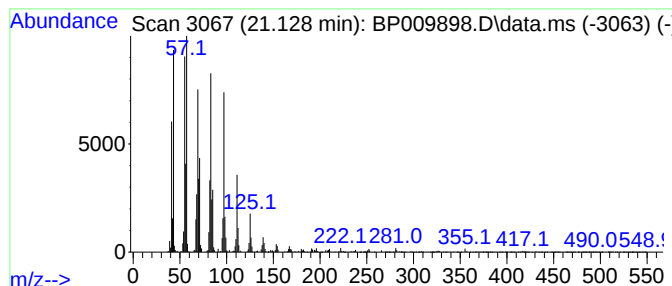
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Fumaric acid, 2-chloropropyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	3.22 ng/ul	453769	Chrysene-d12	21.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fumaric acid, 2-chloropropyl tri...	374	C20H35ClO4	1010348-57-1	97
2			1-Nonadecene	266	C19H38	018435-45-5	95
3			1-Hexacosene	364	C26H52	018835-33-1	95
4			1-Tetracosene	336	C24H48	010192-32-2	94
5			n-Nonadecanol-1	284	C19H40O	001454-84-8	94



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.117	11.6	ng/ul	498375	1	7.958	861549	20.0
Fumaric acid, 2...	21.128	3.2	ng/ul	453769	5	21.416	2817090	20.0