

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070117\
 Data File : BF096395.D
 Acq On : 1 Jul 2017 22:52
 Operator : SJ/MA
 Sample : I3982-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 348-343-COMP

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:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.928	478	483	491	rBV	367053	470462	16.43%	1.823%
2	5.345	549	554	558	rBV	2873663	2858391	99.80%	11.074%
3	6.328	717	721	723	rBV	26571	28714	1.00%	0.111%
4	6.369	723	728	735	rVB	2841484	2864034	100.00%	11.096%
5	6.504	746	751	754	rBV	3019736	2764807	96.54%	10.712%
6	6.733	786	790	793	rBV	1010267	784938	27.41%	3.041%
7	6.892	810	817	820	rBV	2388265	2143285	74.83%	8.304%
8	7.292	880	885	888	rBV	1934366	1687113	58.91%	6.536%
9	7.386	897	901	904	rVB	45128	41714	1.46%	0.162%
10	7.716	950	957	960	rBV	46098	54523	1.90%	0.211%
11	8.016	1004	1008	1011	rBV	1088087	979873	34.21%	3.796%
12	9.092	1186	1191	1194	rBV	2780450	2524581	88.15%	9.781%
13	9.216	1209	1212	1215	rVB	40693	34807	1.22%	0.135%
14	9.486	1253	1258	1261	rBV	1129609	994528	34.72%	3.853%
15	9.580	1267	1274	1278	rBV	54809	59133	2.06%	0.229%
16	9.710	1293	1296	1299	rVB	49356	39272	1.37%	0.152%
17	9.763	1301	1305	1309	rVV2	1010710	886713	30.96%	3.435%
18	10.210	1378	1381	1384	rVB	64565	51950	1.81%	0.201%
19	10.551	1434	1439	1442	rBV	1311326	1285206	44.87%	4.979%
20	10.680	1458	1461	1470	rVB	64329	71634	2.50%	0.278%
21	11.127	1534	1537	1540	rBV2	43423	39849	1.39%	0.154%
22	11.245	1553	1557	1560	rBV	899749	779982	27.23%	3.022%
23	11.780	1644	1648	1652	rBV2	83143	77129	2.69%	0.299%
24	12.451	1759	1762	1766	rBV	51917	51452	1.80%	0.199%
25	12.674	1795	1800	1805	rBV	51228	52940	1.85%	0.205%
26	12.827	1822	1826	1830	rBV	2247140	2115527	73.87%	8.196%
27	13.727	1976	1979	1984	rVB	219172	217665	7.60%	0.843%
28	13.874	1999	2004	2007	rBV	901115	956692	33.40%	3.707%
29	13.898	2007	2008	2012	rVB	51490	28994	1.01%	0.112%
30	14.580	2122	2124	2127	rBV3	23485	35257	1.23%	0.137%
31	14.880	2172	2175	2178	rBV	64666	95821	3.35%	0.371%
32	15.162	2220	2223	2226	rBV	47991	57833	2.02%	0.224%
33	15.280	2238	2243	2250	rBV	552003	676323	23.61%	2.620%

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

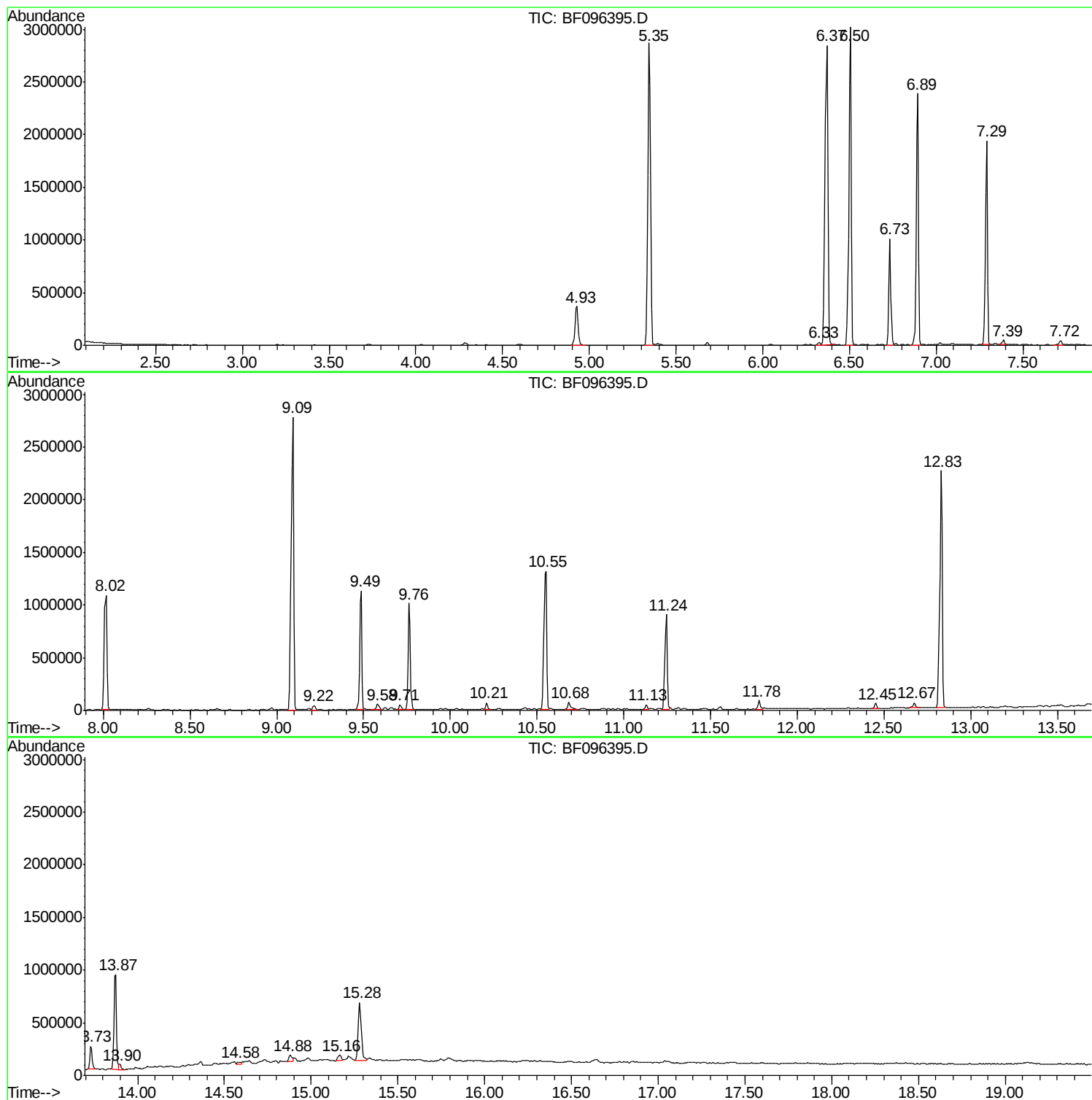
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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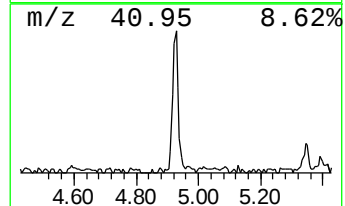
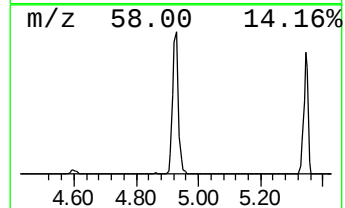
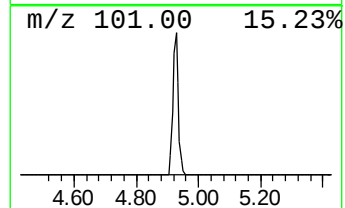
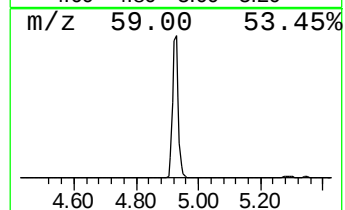
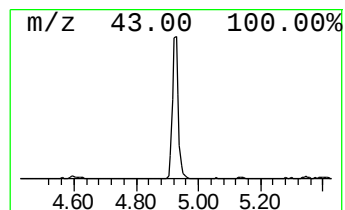
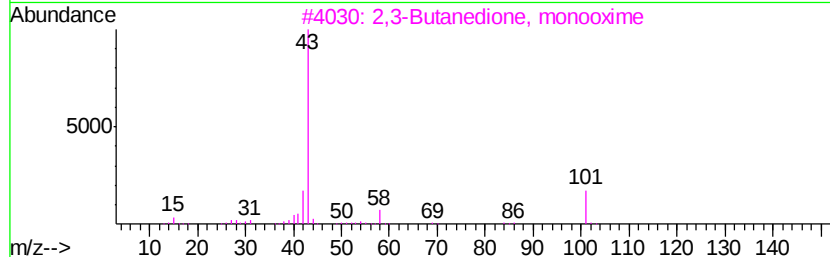
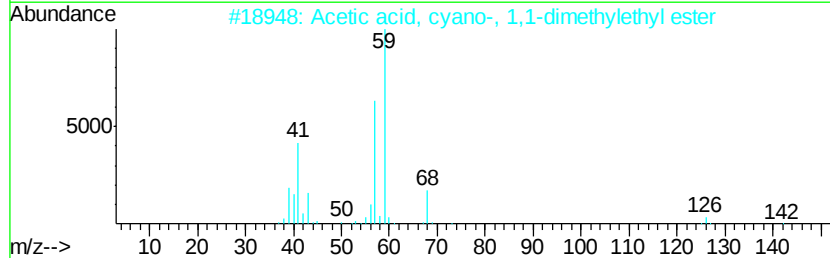
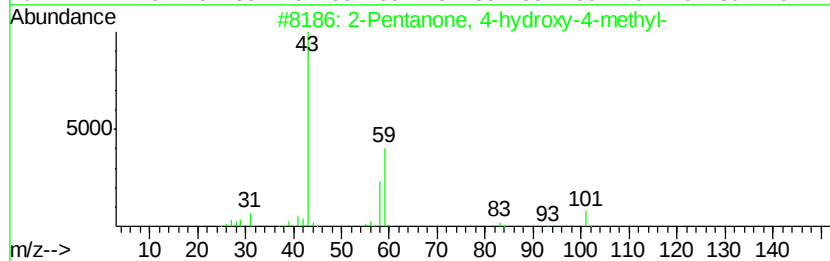
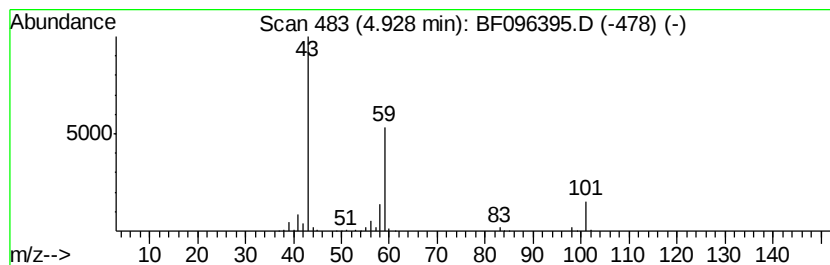
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	11.99 ng	470462	1,4-Dichlorobenzene-d4	6.73

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			5-Oxohexanenitrile	111	C6H9NO	010412-98-3	9
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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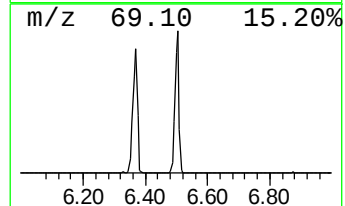
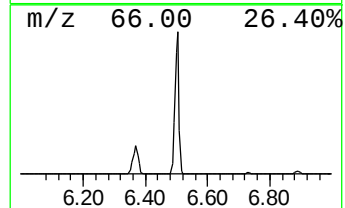
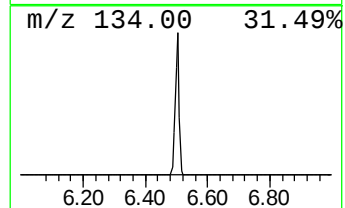
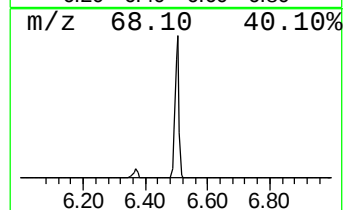
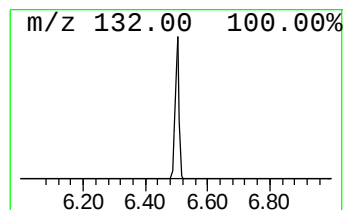
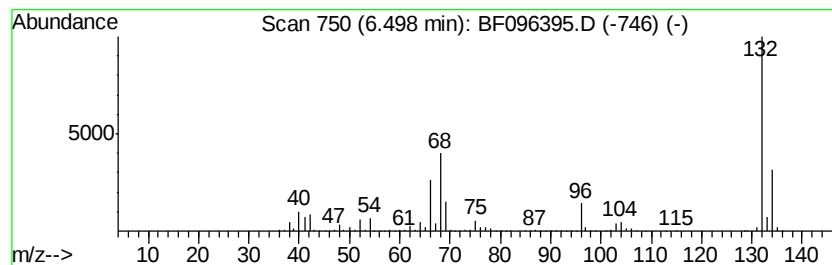
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Peak Number 2 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	70.45 ng	2764810	1,4-Dichlorobenzene-d4	6.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Amphetaminil	250	C17H18N2	017590-01-1	10



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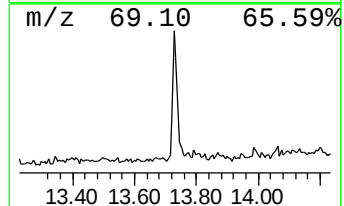
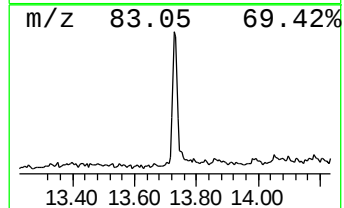
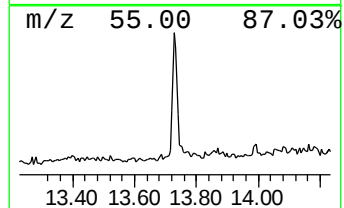
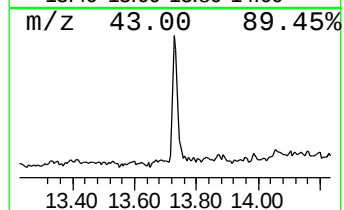
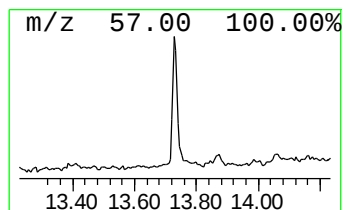
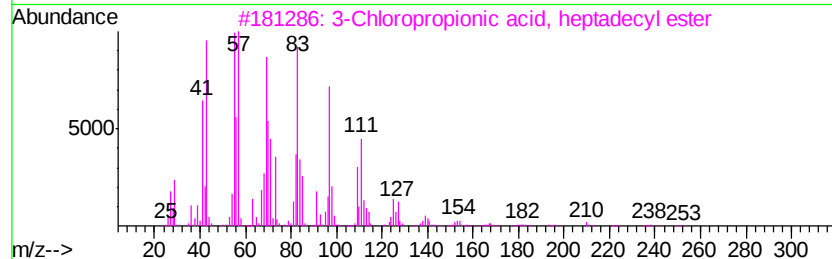
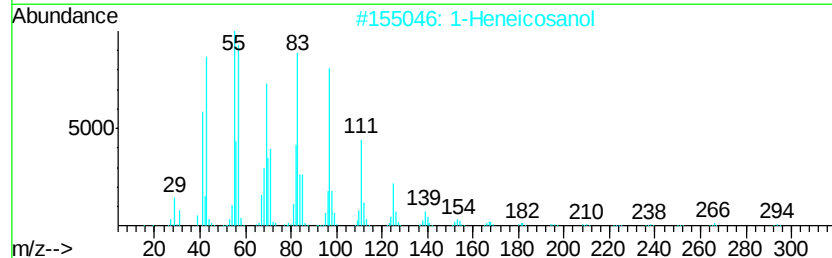
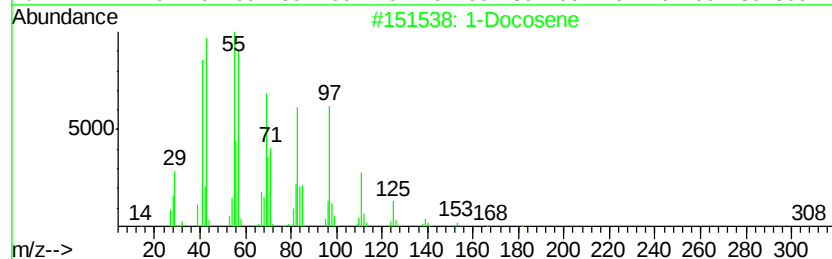
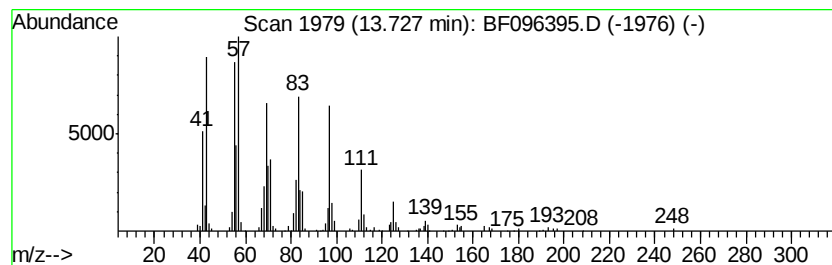
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Peak Number 4 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	4.55 ng	217665	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	94
2		1-Heneicosanol	312	C21H44O	015594-90-8	93
3		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	91
4		1-Nonadecene	266	C19H38	018435-45-5	91
5		1-Hexadecanol	242	C16H34O	036653-82-4	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.93	12.0	ng	470462	1	6.73	784938	20.0
unknown6.50	6.50	70.5	ng	2764810	1	6.73	784938	20.0
1-Docosene	13.73	4.5	ng	217665	5	13.87	956692	20.0