

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021017\
 Data File : BF092886.D
 Acq On : 10 Feb 2017 16:45
 Operator : SJ/MA
 Sample : PB96842BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB96842BL

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-BF013017.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	5.082	260	262	269	rBV	1091666	1650943	33.77%	4.028%
2	5.504	296	299	301	rBV	3877752	4004768	81.91%	9.771%
3	6.533	386	389	391	rBV	3533664	3778367	77.28%	9.218%
4	6.659	398	400	403	rBV	3081200	4194930	85.80%	10.235%
5	6.887	418	420	423	rBV	871007	1030261	21.07%	2.514%
6	7.047	431	434	437	rBV	3043367	3125085	63.92%	7.624%
7	7.459	467	470	473	rBV	2427915	2546350	52.08%	6.213%
8	8.179	530	533	535	rBV	1514176	1428378	29.21%	3.485%
9	9.253	624	627	630	rBV	3387582	4797580	98.12%	11.705%
10	9.848	669	679	684	rBV	129750	307777	6.29%	0.751%
11	9.939	684	687	689	rVB	1134982	1675213	34.26%	4.087%
12	10.305	711	719	727	rBV2	14081	50262	1.03%	0.123%
13	10.728	752	756	758	rBV	2243113	2865684	58.61%	6.992%
14	11.414	814	816	819	rBV	1269329	1644517	33.64%	4.012%
15	13.014	952	956	958	rBV	3500728	4889278	100.00%	11.929%
16	14.054	1045	1047	1050	rBV	1577368	1489343	30.46%	3.634%
17	15.505	1171	1174	1183	rBV	1086074	1399186	28.62%	3.414%
18	15.940	1209	1212	1215	rBV	30419	52718	1.08%	0.129%
19	16.431	1252	1255	1260	rVB	31687	56884	1.16%	0.139%

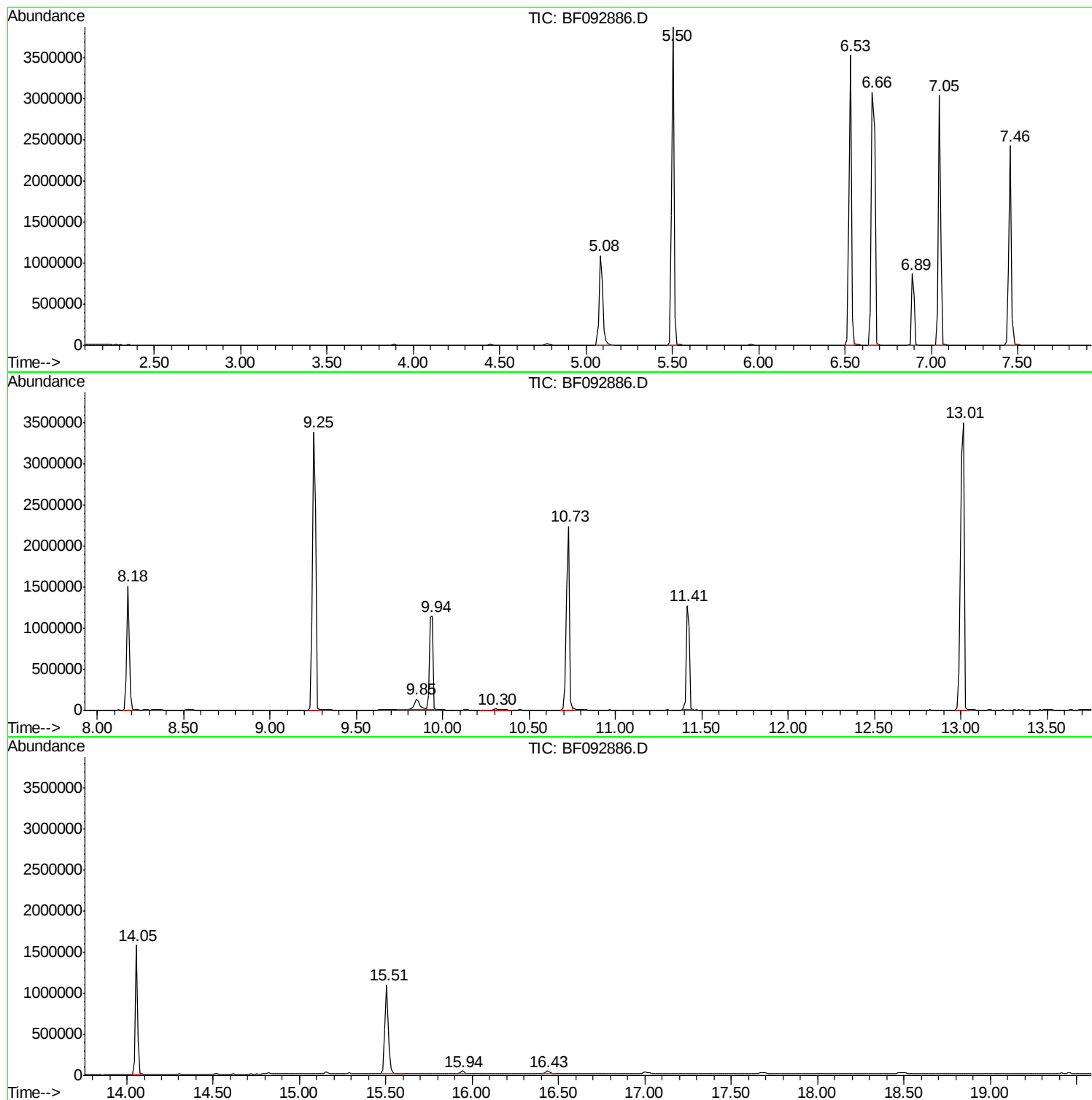
Sum of corrected areas: 40987524

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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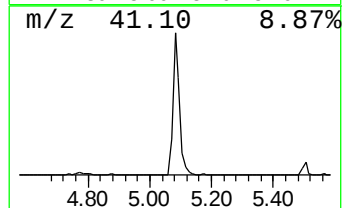
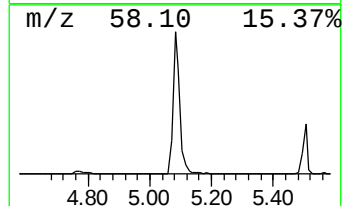
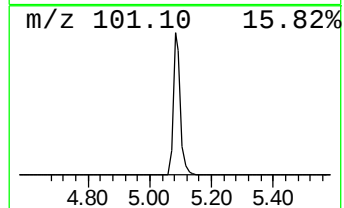
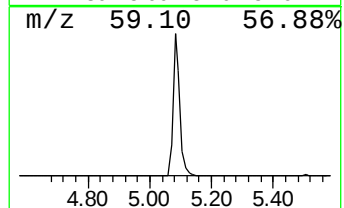
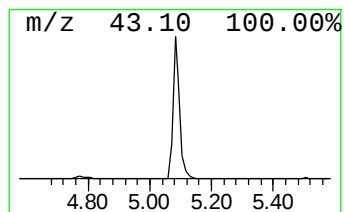
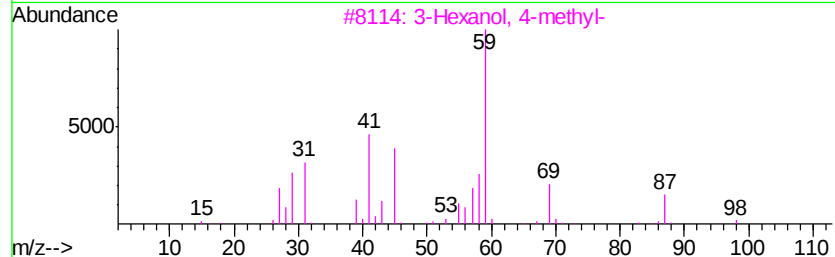
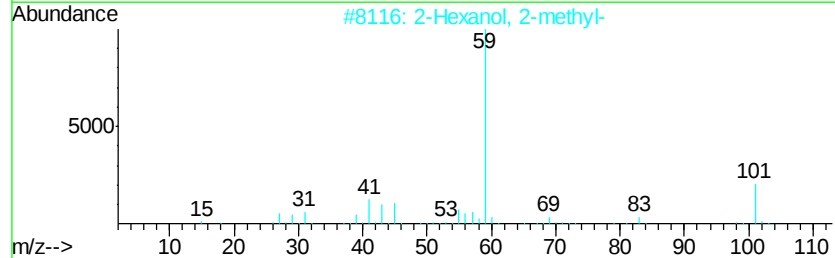
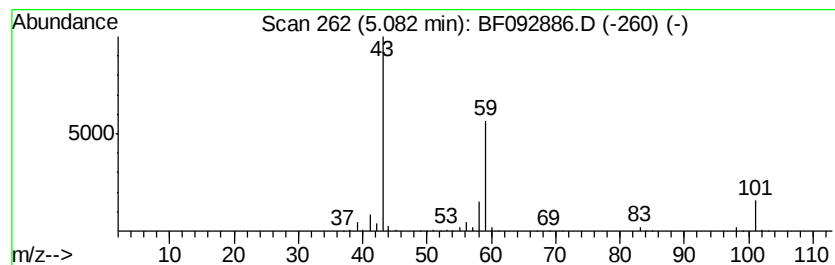
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	32.05 ng	1650940	1,4-Dichlorobenzene-d4	6.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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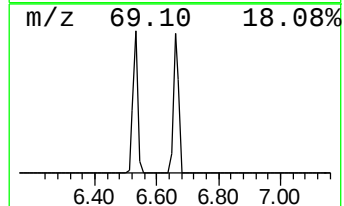
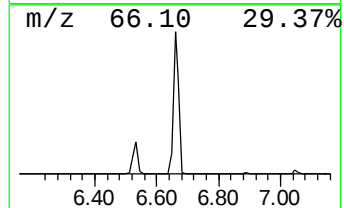
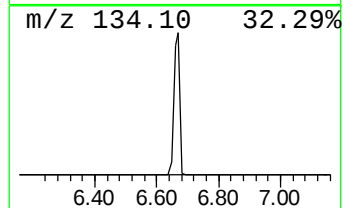
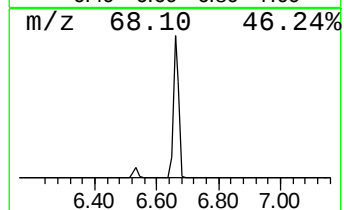
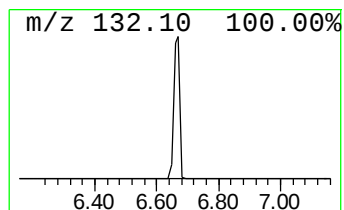
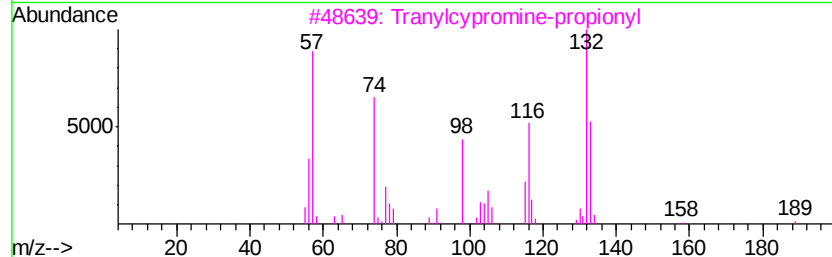
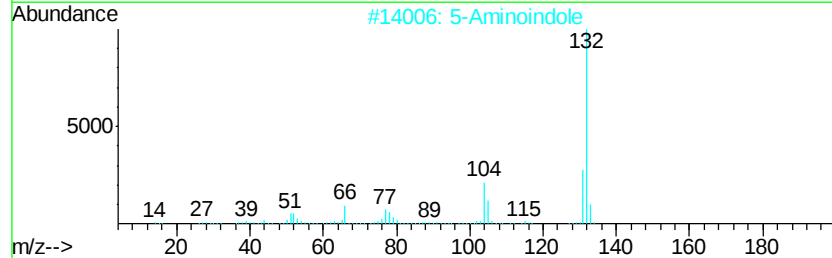
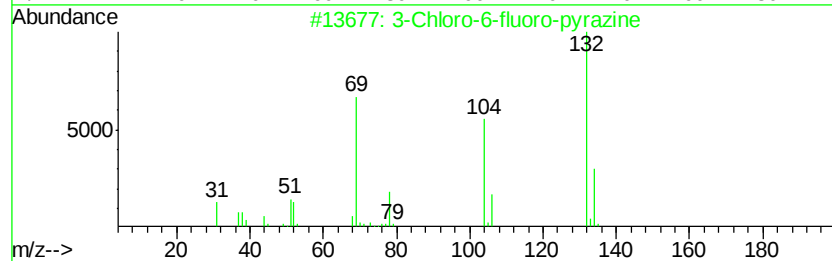
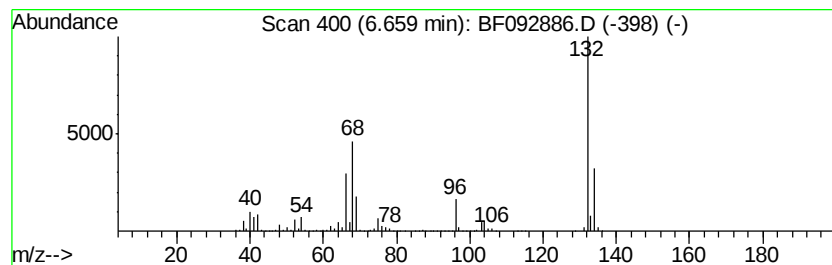
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Peak Number 2 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	81.43 ng	4194930	1,4-Dichlorobenzene-d4	6.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole			132	C8H8N2	005192-03-0	12
3	Tranylcypromine-propionyl			189	C12H15NO	1000123-86-3	10
4	Imidazole, 4,5-dicyano-1-methyl-			132	C6H4N4	019485-35-9	9
5	Benzene, 1,3,5-trifluoro-			132	C6H3F3	000372-38-3	9



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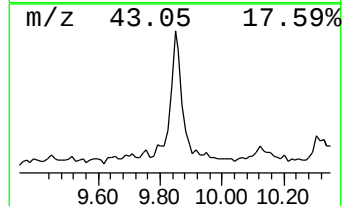
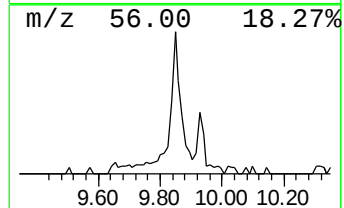
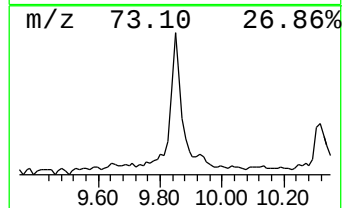
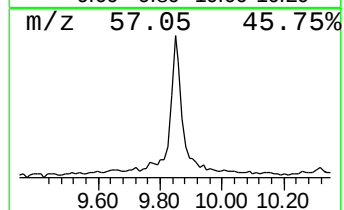
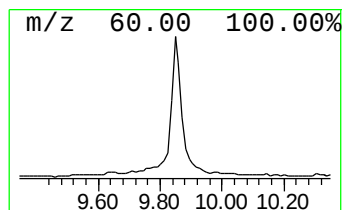
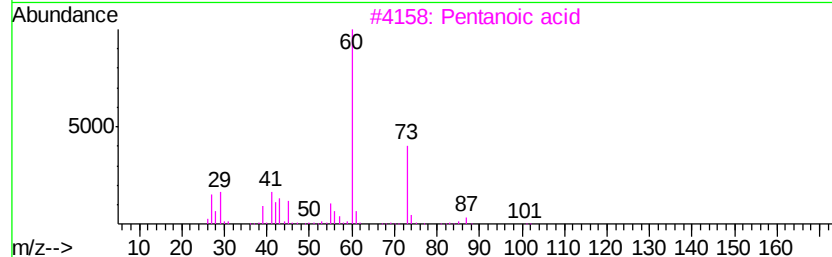
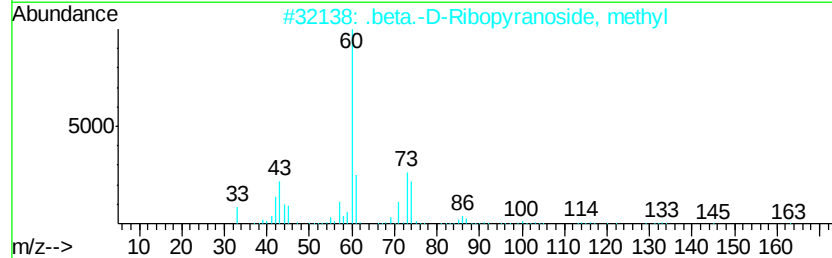
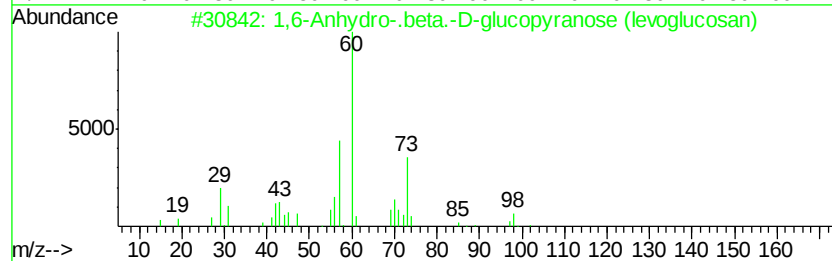
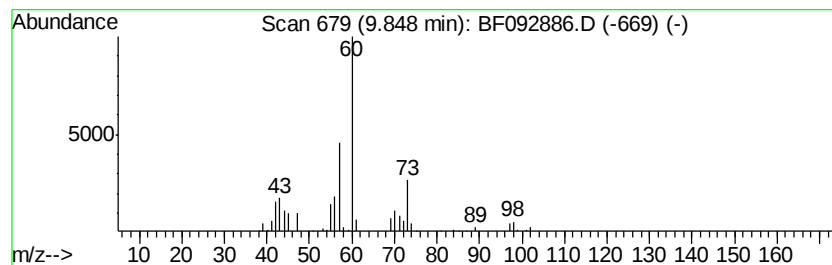
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Peak Number 4 1,6-Anhydro-.beta.-D-glucop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.85	3.67 ng	307777	Acenaphthene-d10	9.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,6-Anhydro-.beta.-D-glucopyrano...	162	C6H10O5	000498-07-7	72
2		.beta.-D-Ribopyranoside, methyl	164	C6H12O5	017289-61-1	59
3		Pentanoic acid	102	C5H10O2	000109-52-4	50
4		Heptanoic acid	130	C7H14O2	000111-14-8	45
5		3,4-Altrosan	162	C6H10O5	1000129-76-4	40



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.08	32.0	ng	1650940	1	6.89	1030260	20.0
unknown6.66	6.66	81.4	ng	4194930	1	6.89	1030260	20.0
1,6-Anhydro-.beta...	9.85	3.7	ng	307777	3	9.94	1675210	20.0