

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082817\
 Data File : BF098075.D
 Acq On : 28 Aug 2017 19:27
 Operator : SJ/JU
 Sample : PB101859BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB101859BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % 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:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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	4.939	480	485	495	rBV	260046	350250	11.05%	1.523%
2	5.351	550	555	558	rBV	2115427	2054205	64.78%	8.934%
3	6.375	724	729	732	rBV	2261575	2061569	65.01%	8.966%
4	6.510	747	752	755	rBV	2249408	2041211	64.37%	8.878%
5	6.739	787	791	794	rBV	847083	664639	20.96%	2.891%
6	6.898	813	818	821	rBV	2494017	2266566	71.48%	9.858%
7	7.304	882	887	890	rBV	1838134	1883697	59.40%	8.193%
8	8.022	1005	1009	1013	rBV	947665	861425	27.16%	3.747%
9	9.104	1187	1193	1196	rBV	2844344	3171097	100.00%	13.792%
10	9.774	1302	1307	1317	rVB	965954	897088	28.29%	3.902%
11	10.563	1436	1441	1460	rBV	1314617	1256525	39.62%	5.465%
12	11.257	1555	1559	1566	rBV	1065212	918759	28.97%	3.996%
13	12.845	1824	1829	1832	rBV	2953914	2949152	93.00%	12.826%
14	13.892	2002	2007	2010	rBV	972612	872116	27.50%	3.793%
15	15.315	2243	2249	2259	rBV	560452	744490	23.48%	3.238%

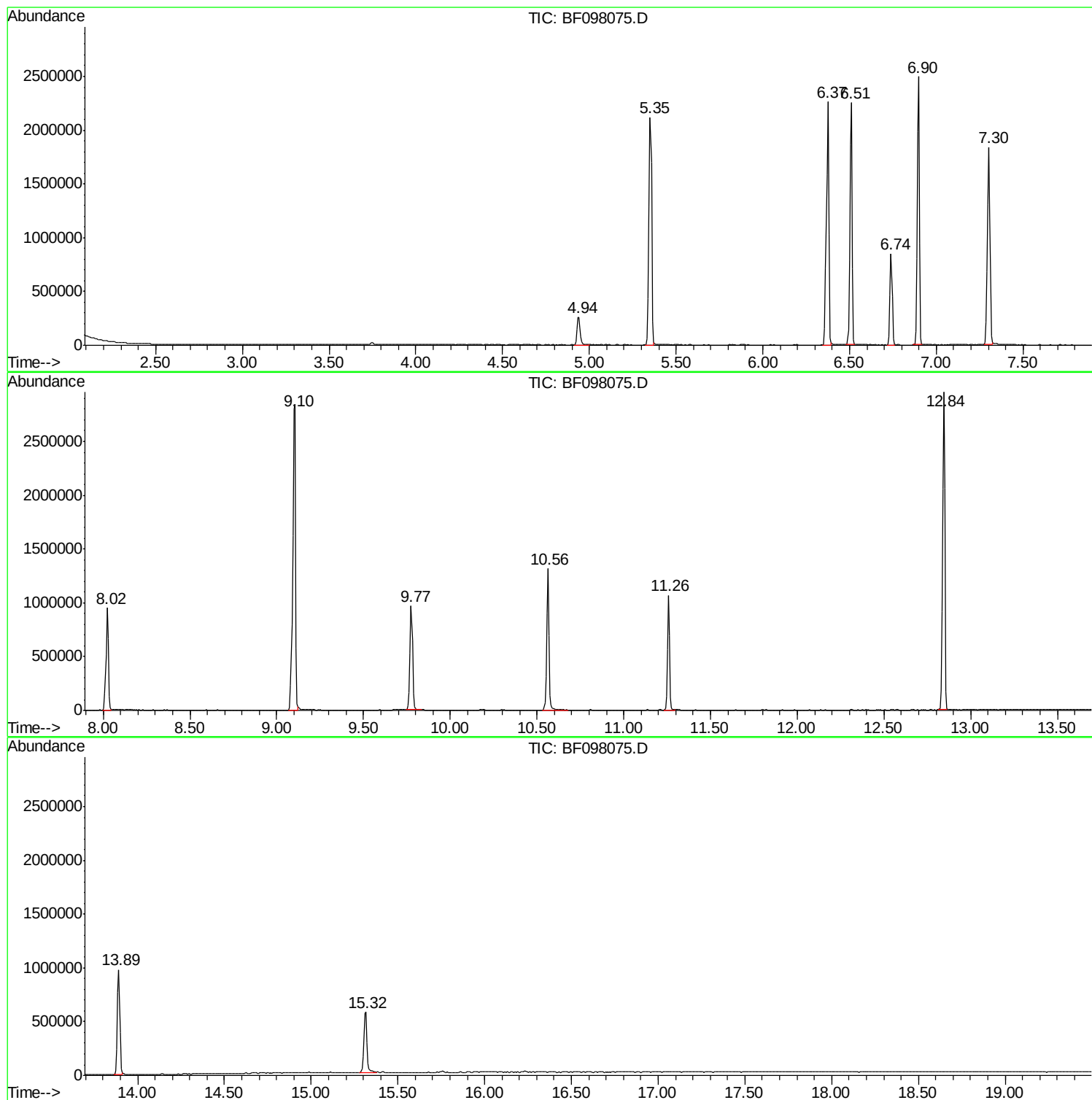
Sum of corrected areas: 22992789

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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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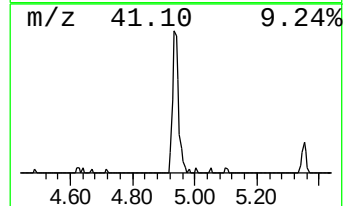
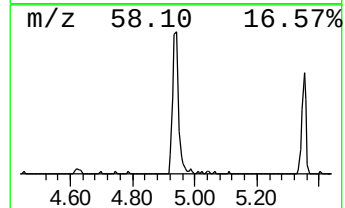
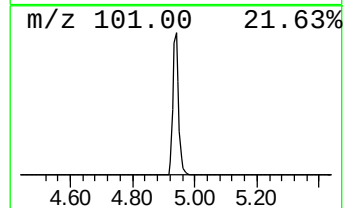
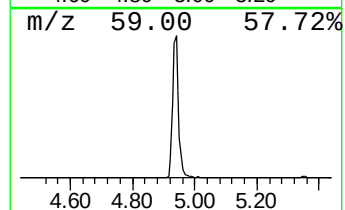
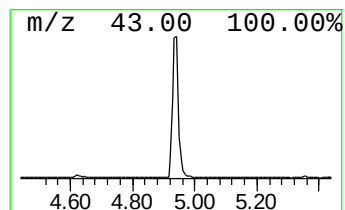
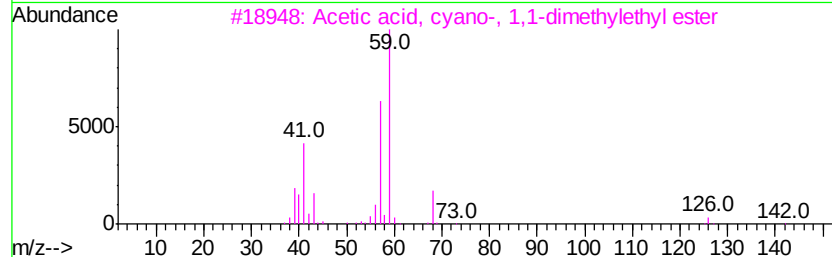
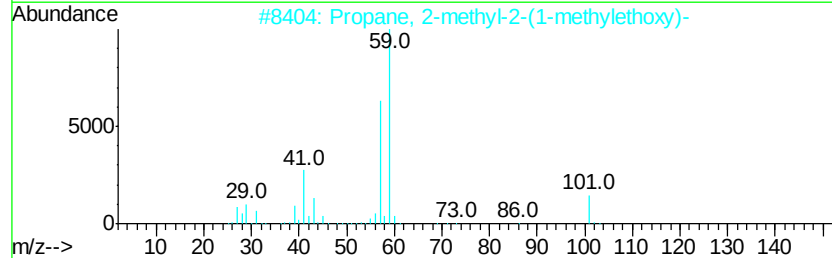
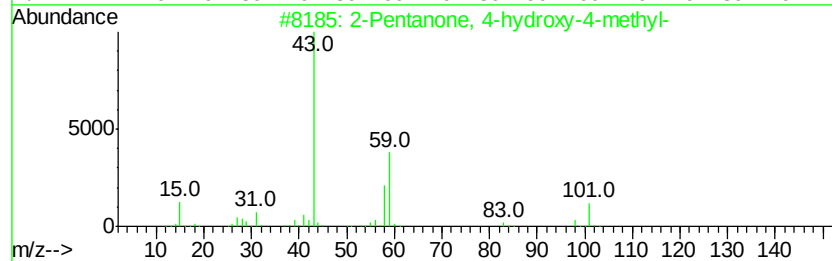
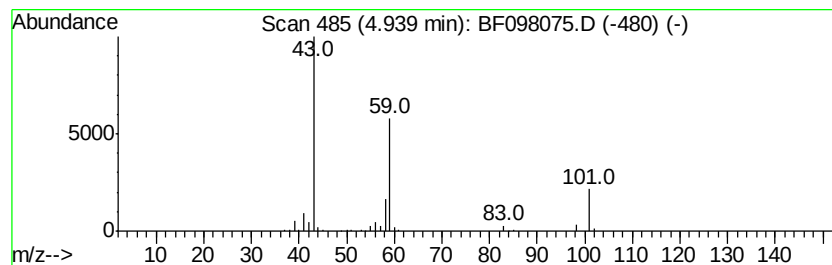
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.94	10.54 ng	350250	1,4-Dichlorobenzene-d4	6.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	23
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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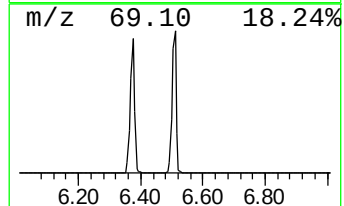
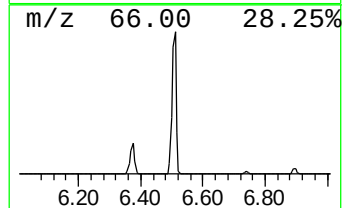
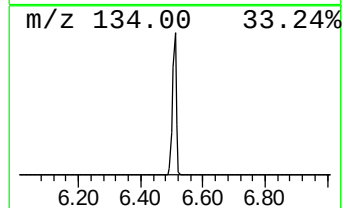
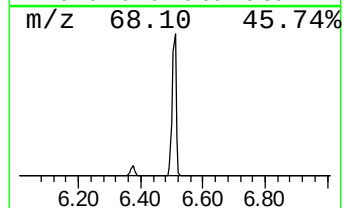
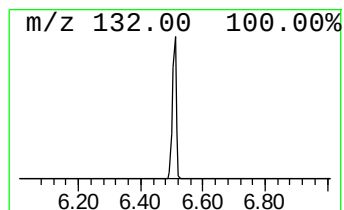
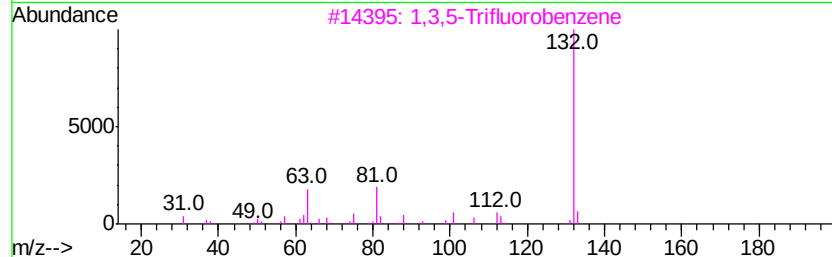
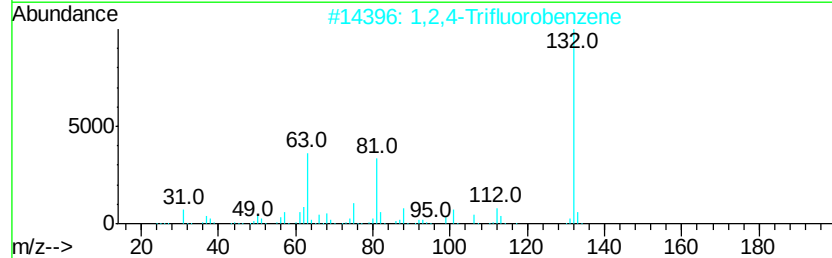
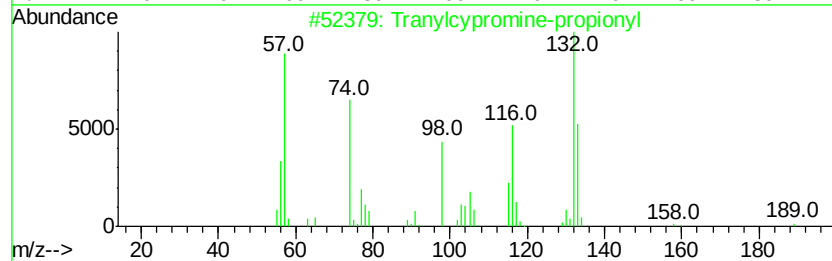
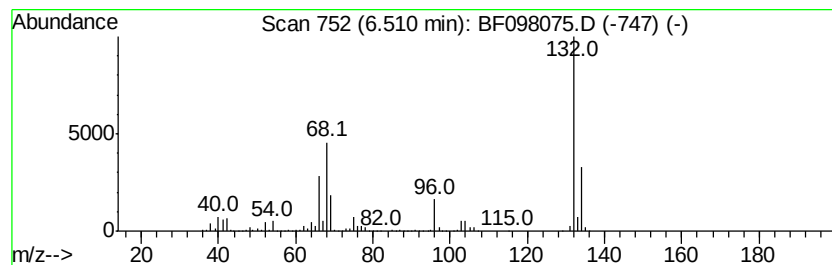
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Peak Number 2 unknown6.51 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	61.42 ng	2041210	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
2		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
3		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.94	10.5	ng	350250	1	6.74	664639	20.0
unknown6.51	6.51	61.4	ng	2041210	1	6.74	664639	20.0