

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111915\
 Data File : BF083052.D
 Acq On : 19 Nov 2015 23:35
 Operator : UM/IZ
 Sample : PB86723BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB86723BL

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-BF111915.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	4.853	239	242	252	rVB	1323050	1907537	44.46%	5.786%
2	5.299	279	281	284	rBV	2860944	3090127	72.03%	9.372%
3	6.350	370	373	375	rBV	3115234	3070179	71.56%	9.312%
4	6.476	381	384	386	rBV	2774500	3449405	80.40%	10.462%
5	6.704	401	404	406	rVB	593268	705070	16.43%	2.138%
6	6.853	415	417	420	rBV	2228875	2514244	58.60%	7.626%
7	7.265	450	453	455	rBV	2172105	2110791	49.20%	6.402%
8	7.985	514	516	519	rBV	904160	891716	20.79%	2.705%
9	9.070	608	611	613	rBV	4143675	3906946	91.07%	11.850%
10	9.745	667	670	672	rBV	956706	1088312	25.37%	3.301%
11	10.533	736	739	741	rBV	1991368	2626211	61.21%	7.965%
12	11.219	796	799	801	rBV	1407508	1173891	27.36%	3.560%
13	12.808	935	938	941	rBV	3677861	4290170	100.00%	13.012%
14	13.848	1026	1029	1031	rBV	1387959	1186381	27.65%	3.598%
15	15.231	1147	1150	1153	rBV	871259	960016	22.38%	2.912%

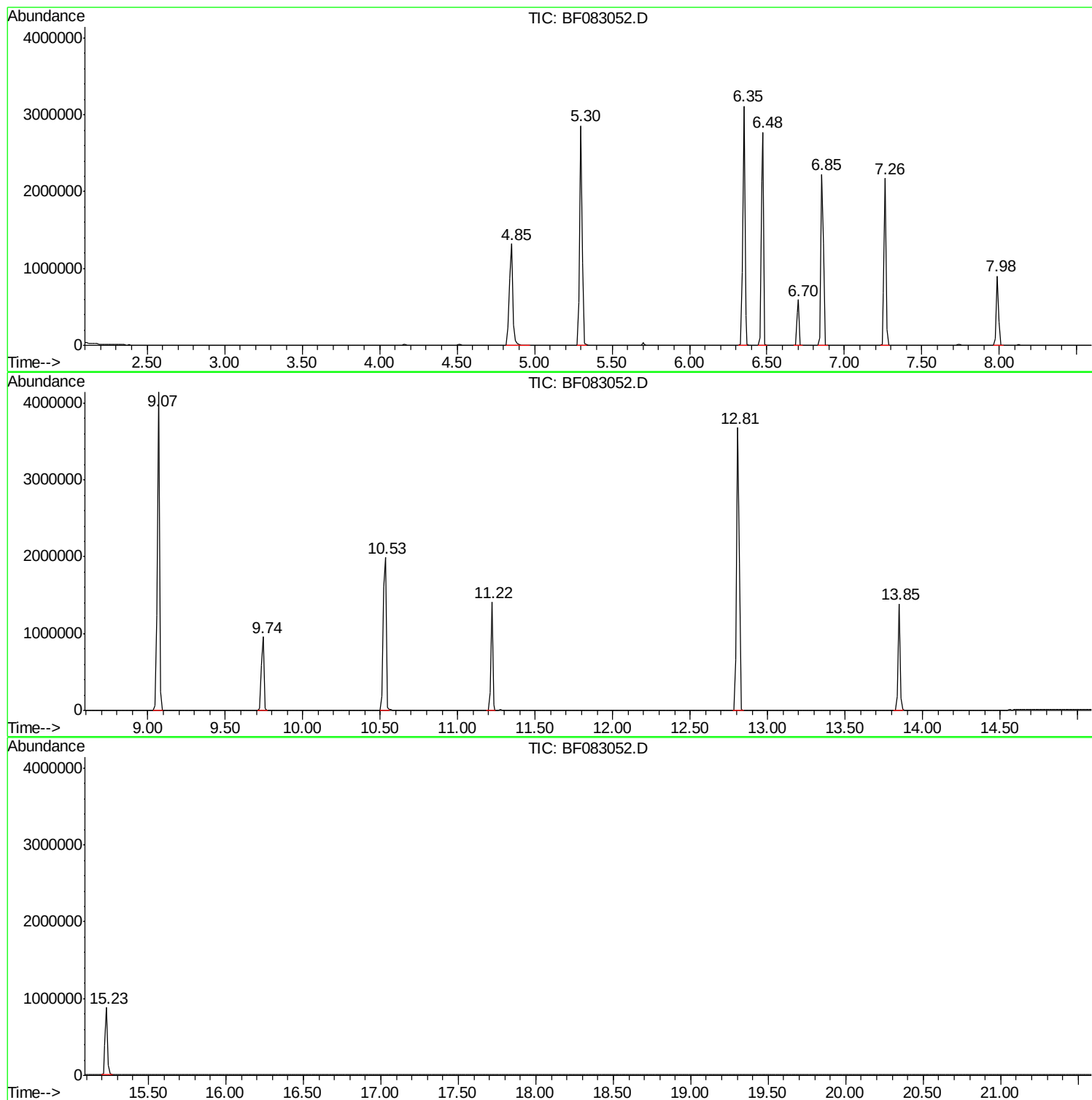
Sum of corrected areas: 32970996

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF111915.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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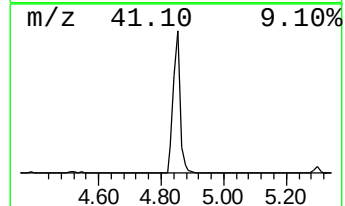
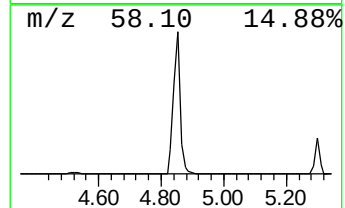
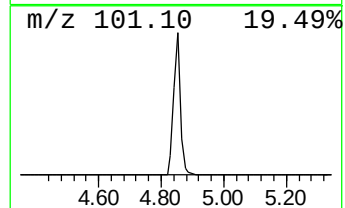
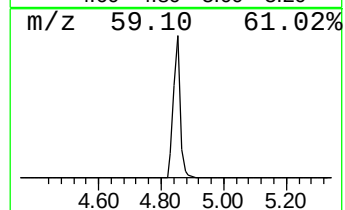
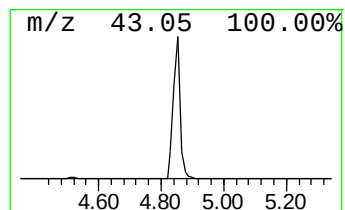
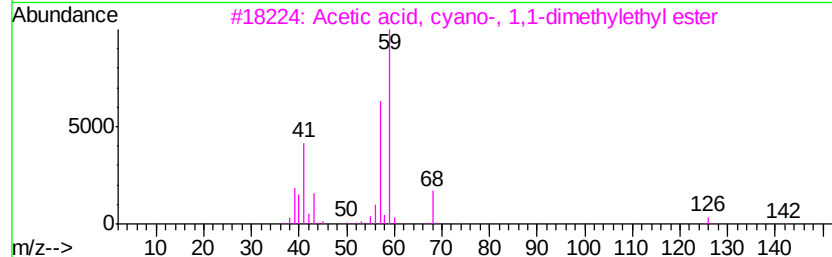
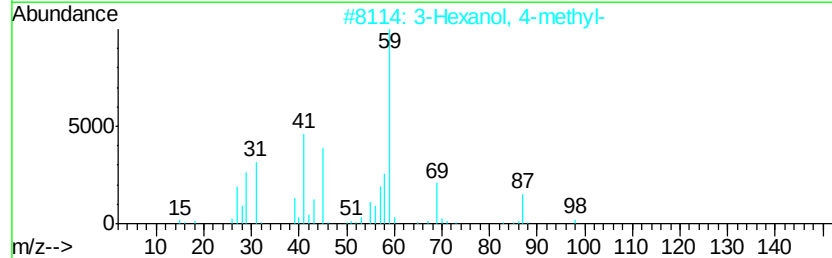
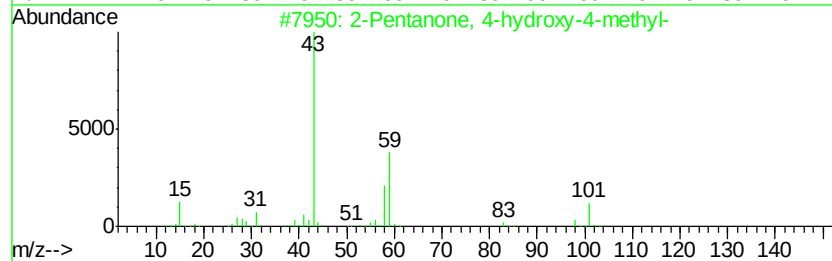
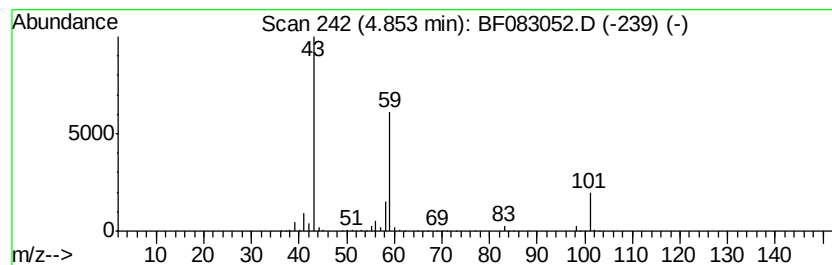
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.85	54.11 ng	1907540	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9



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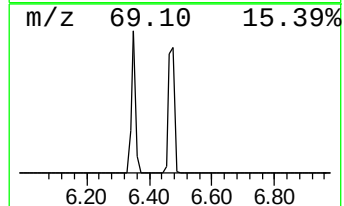
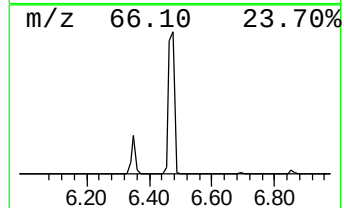
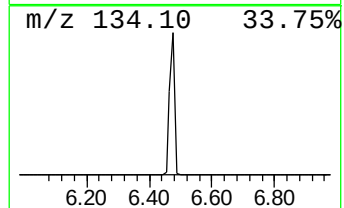
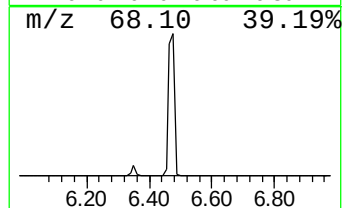
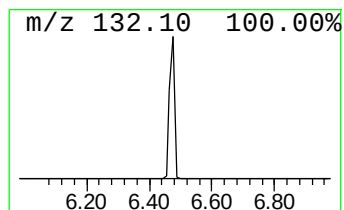
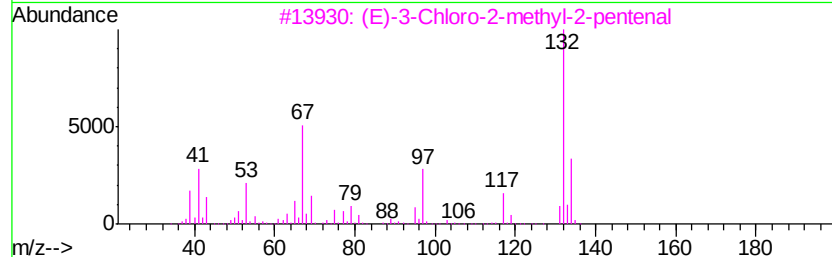
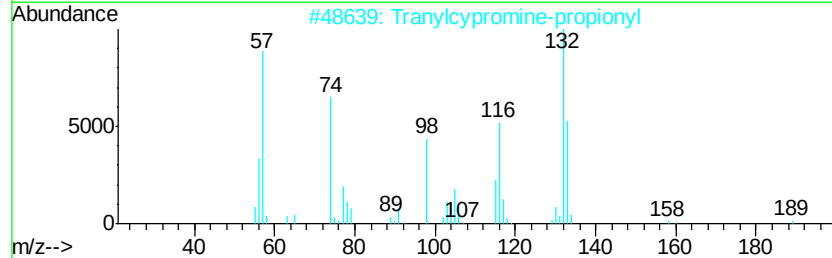
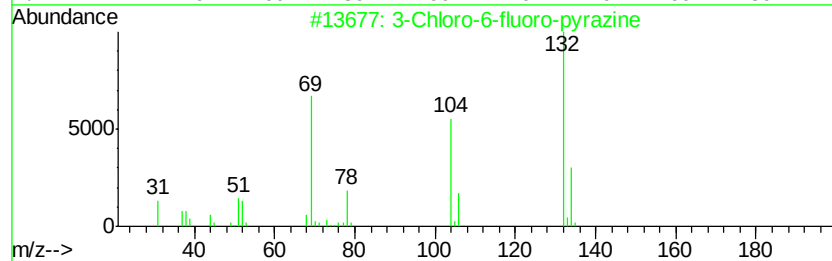
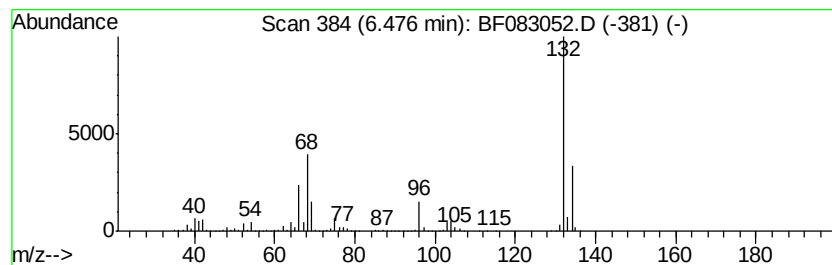
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Peak Number 2 unknown6.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	97.85 ng	3449410	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.85	54.1	ng	1907540	1	6.70	705070	20.0
unknown6.48	6.48	97.8	ng	3449410	1	6.70	705070	20.0