

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102315\
 Data File : BF082476.D
 Acq On : 23 Oct 2015 20:24
 Operator : UM/IZ
 Sample : PB86188BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB86188BL

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-BF101015.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	4.914	240	243	254	rBV	595739	1017579	33.42%	4.826%
2	5.360	279	282	284	rBV	1383637	1767932	58.07%	8.385%
3	6.400	370	373	381	rBV	1718069	1723480	56.61%	8.174%
4	6.526	381	384	386	rBV	1501768	1899265	62.38%	9.008%
5	6.754	401	404	406	rBV	451122	547273	17.97%	2.596%
6	6.903	415	417	420	rBV	1605738	1495438	49.12%	7.093%
7	7.326	451	454	456	rBV	1073705	1175233	38.60%	5.574%
8	8.035	514	516	519	rBV	581755	748455	24.58%	3.550%
9	9.120	608	611	614	rBV	2418633	2412672	79.24%	11.443%
10	9.795	667	670	673	rBV	932418	955240	31.37%	4.531%
11	10.595	737	740	742	rBV	1048602	1397669	45.91%	6.629%
12	11.281	798	800	803	rBV	874723	994260	32.66%	4.716%
13	12.881	937	940	942	rBV	2701314	3044661	100.00%	14.441%
14	13.944	1030	1033	1036	rBV	1048983	1049882	34.48%	4.980%
15	15.418	1159	1162	1171	rBV	677634	854766	28.07%	4.054%

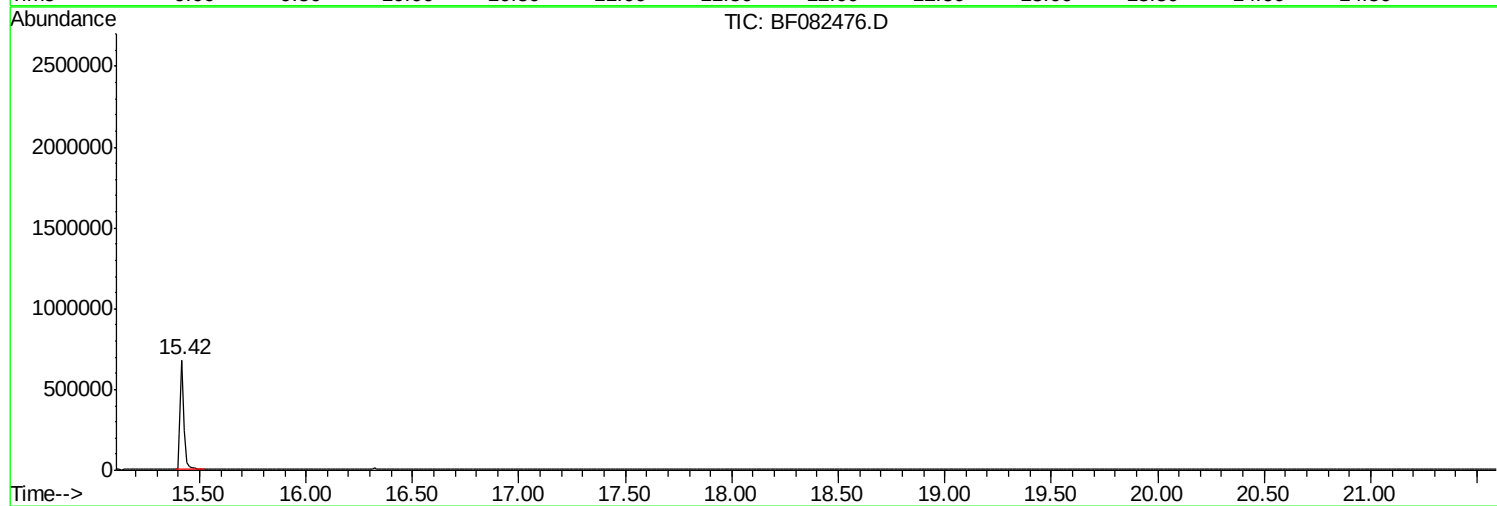
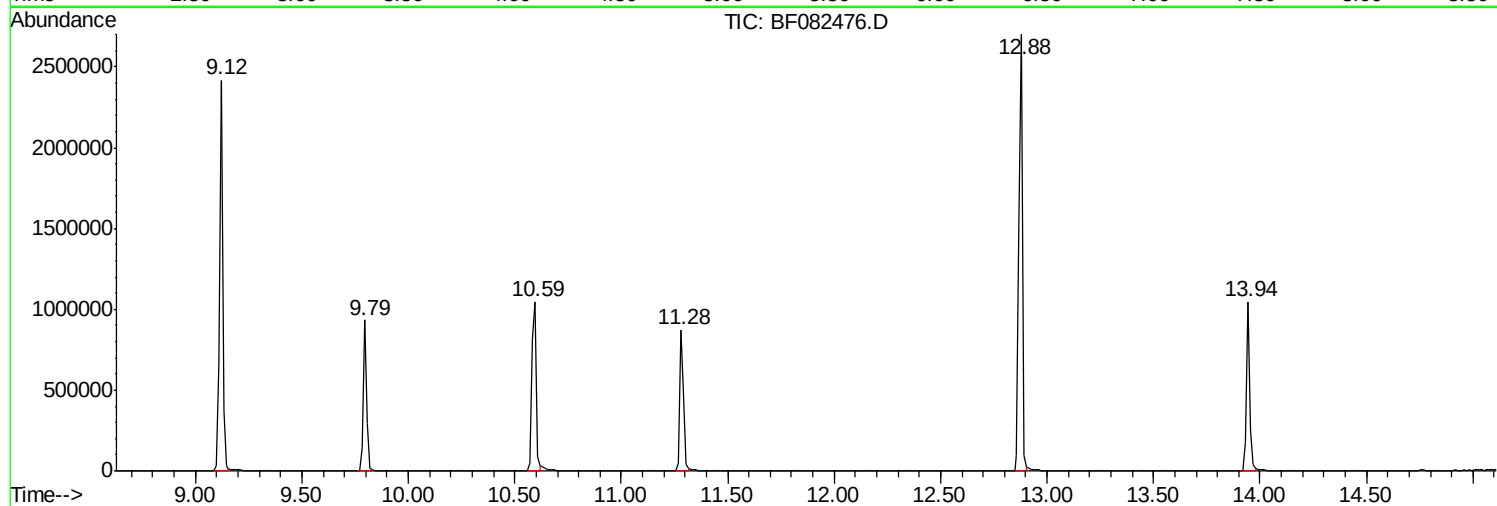
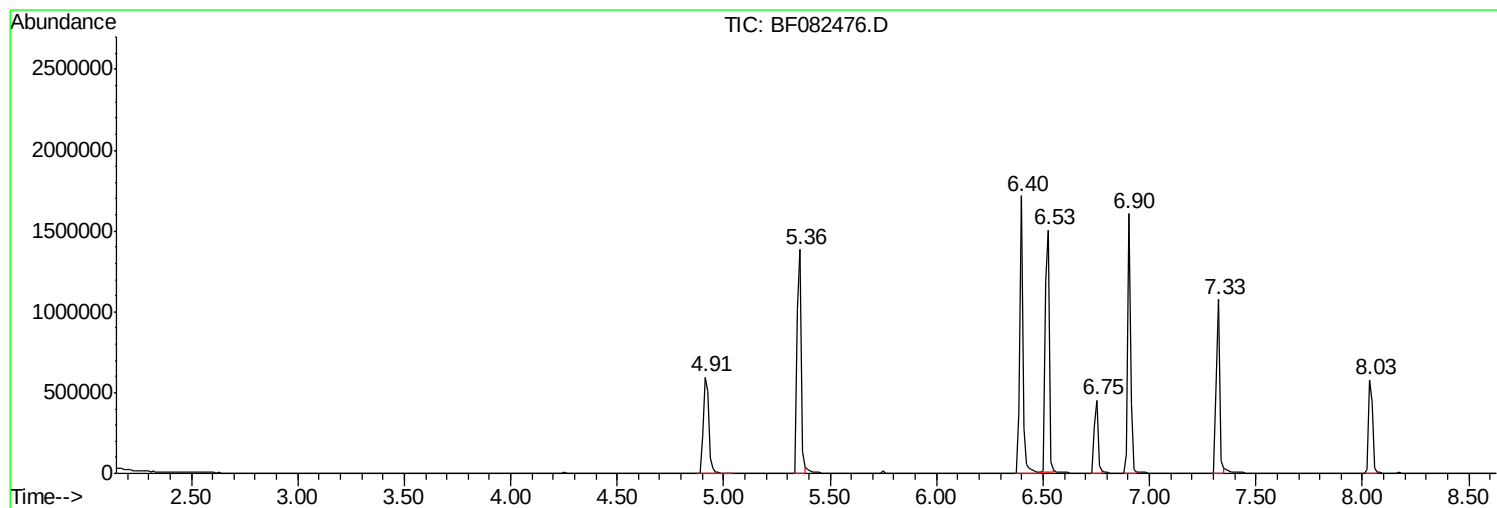
Sum of corrected areas: 21083805

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.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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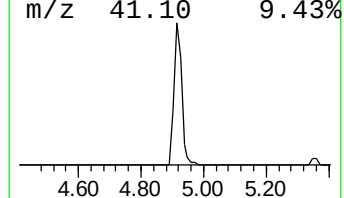
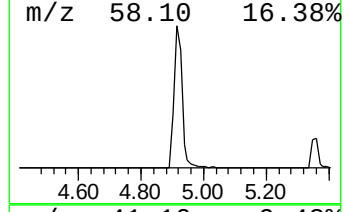
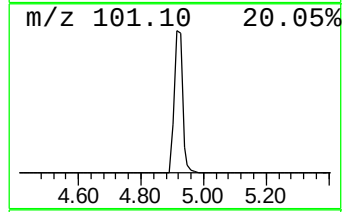
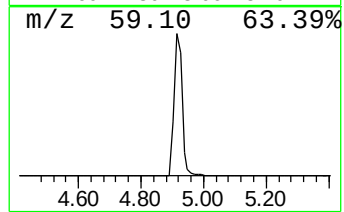
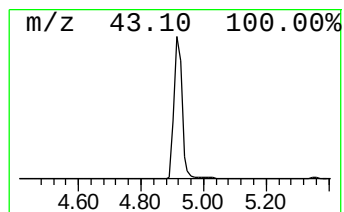
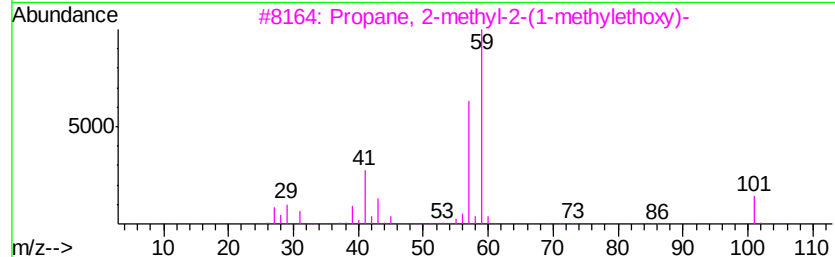
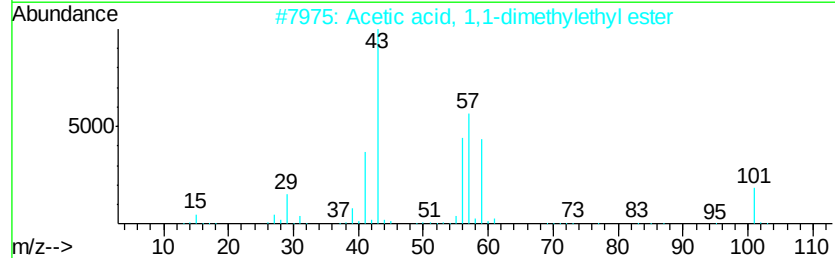
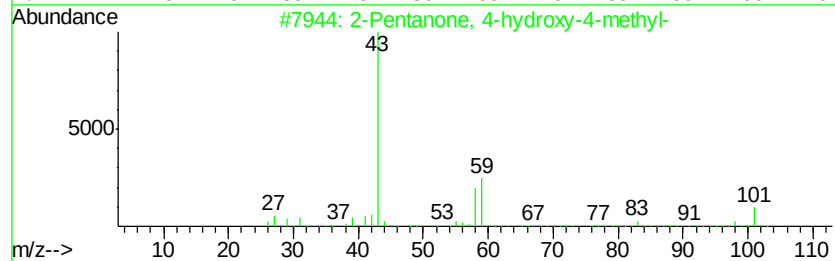
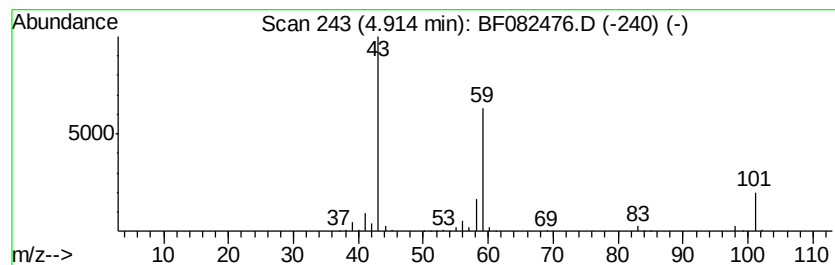
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.91	37.19 ng	1017580	1,4-Dichlorobenzene-d4	6.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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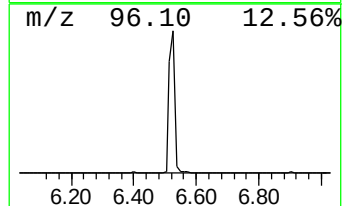
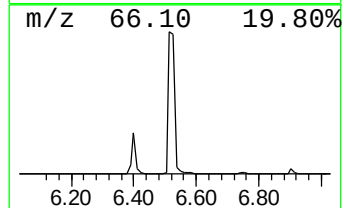
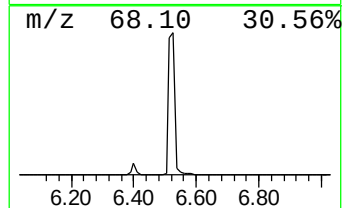
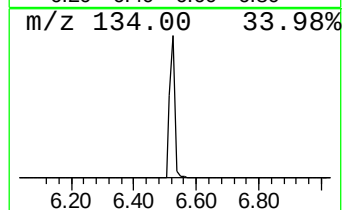
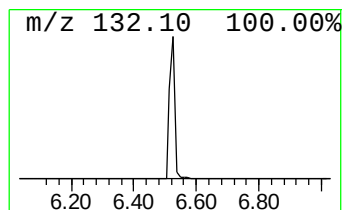
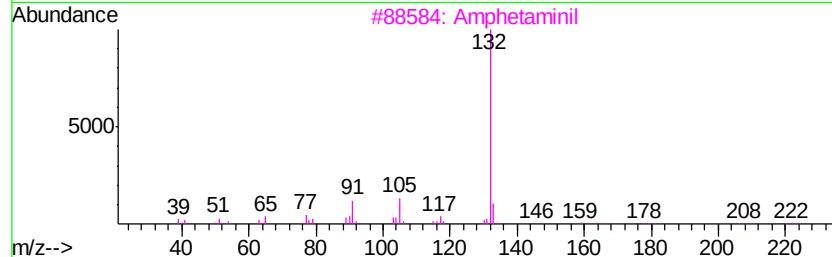
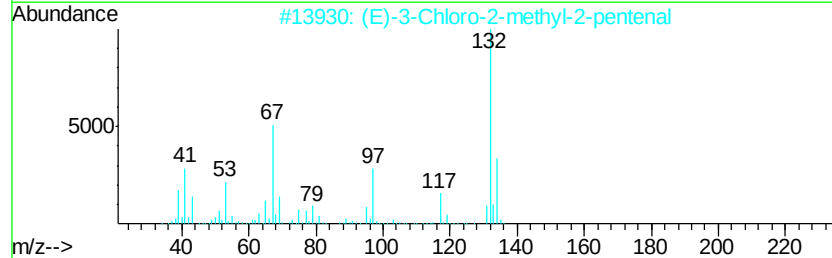
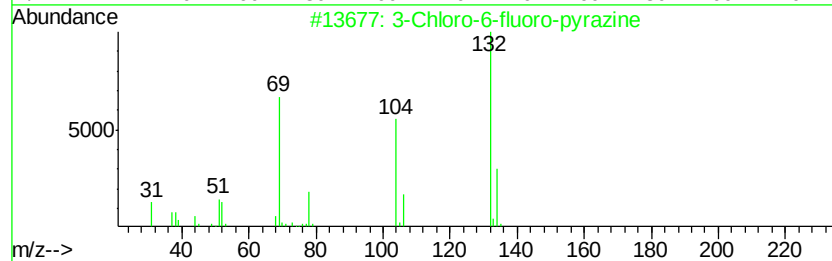
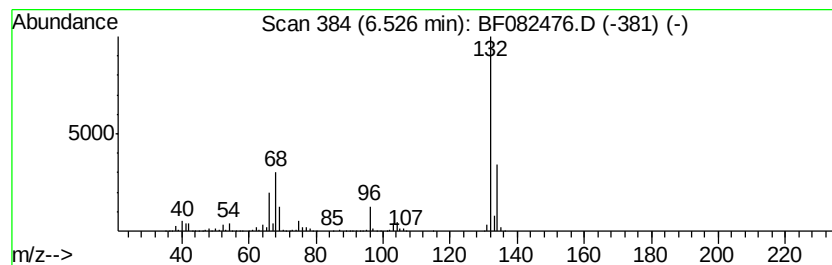
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Peak Number 2 unknown6.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	69.41 ng	1899270	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.91	37.2	ng	1017580	1	6.75	547273	20.0
unknown6.53	6.53	69.4	ng	1899270	1	6.75	547273	20.0