

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062215\
 Data File : BF080102.D
 Acq On : 22 Jun 2015 19:05
 Operator : TP/IZ
 Sample : PB84129BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84129BL

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-BF061715.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	5.533	212	214	220	rBV	238905	229636	15.44%	2.220%
2	5.933	246	249	251	rBV	767951	893994	60.12%	8.642%
3	6.916	333	335	338	rBV	971910	893246	60.07%	8.635%
4	7.087	348	350	352	rBV	1067135	971012	65.30%	9.387%
5	7.316	368	370	373	rBV	261384	280903	18.89%	2.715%
6	7.476	382	384	387	rBV	781350	725140	48.77%	7.010%
7	7.876	417	419	421	rBV	687768	564442	37.96%	5.456%
8	8.607	481	483	486	rBV	415879	395853	26.62%	3.827%
9	9.682	575	577	580	rBV	934302	1130828	76.05%	10.931%
10	10.379	636	638	641	rVB2	422163	454012	30.53%	4.389%
11	11.179	705	708	710	rBV2	518510	689804	46.39%	6.668%
12	11.888	768	770	772	rBV	511968	513752	34.55%	4.966%
13	13.522	911	913	915	rBV	1749954	1486989	100.00%	14.374%
14	14.734	1017	1019	1022	rBV	536617	564051	37.93%	5.453%
15	16.563	1176	1179	1187	rBV	393798	551012	37.06%	5.327%

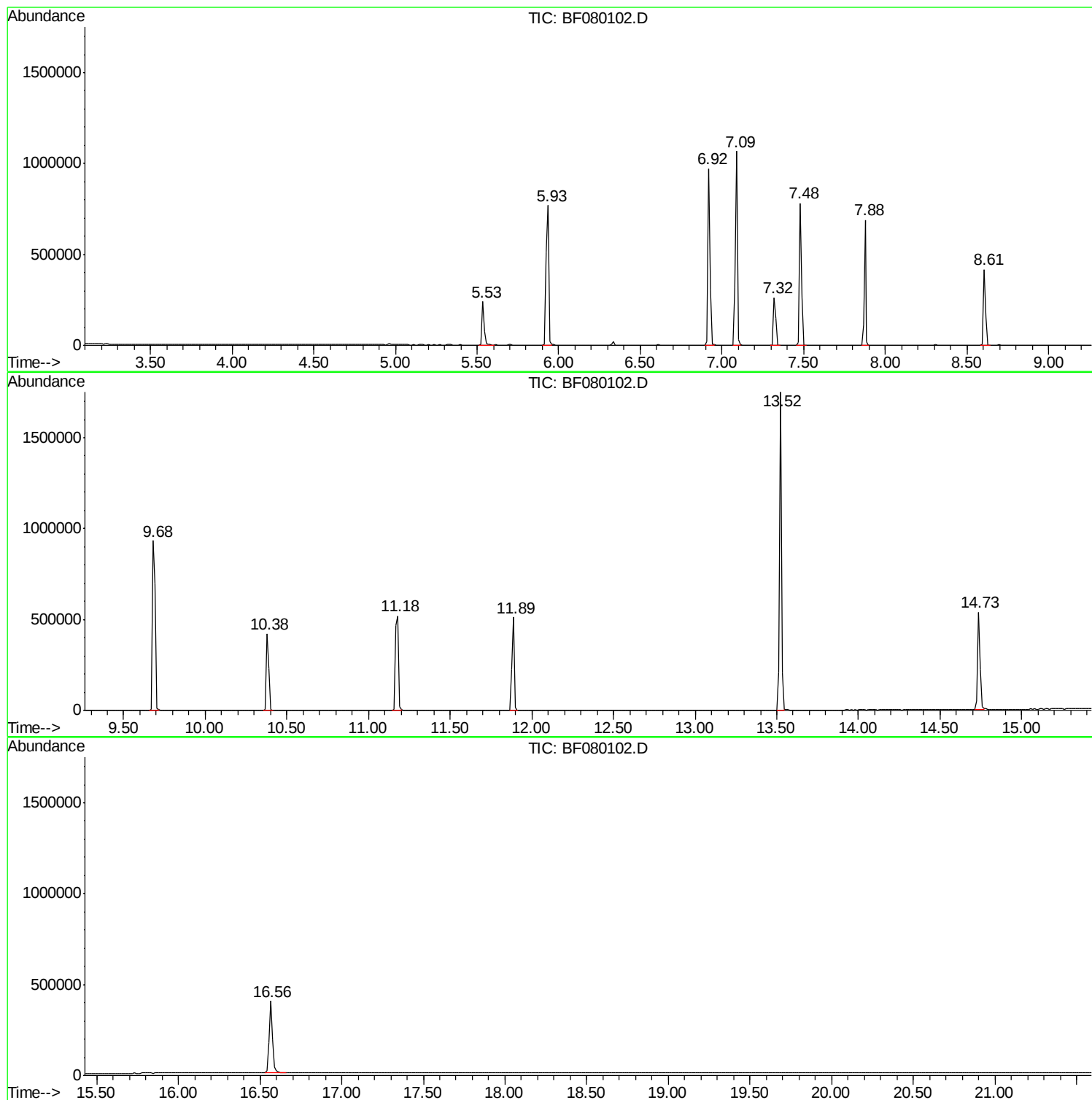
Sum of corrected areas: 10344674

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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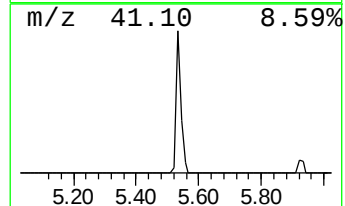
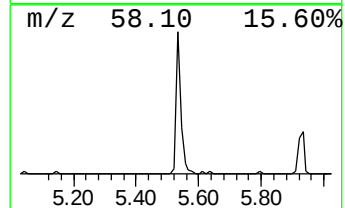
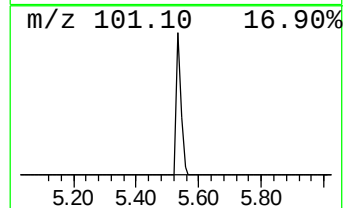
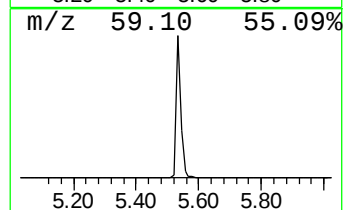
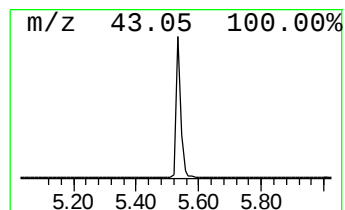
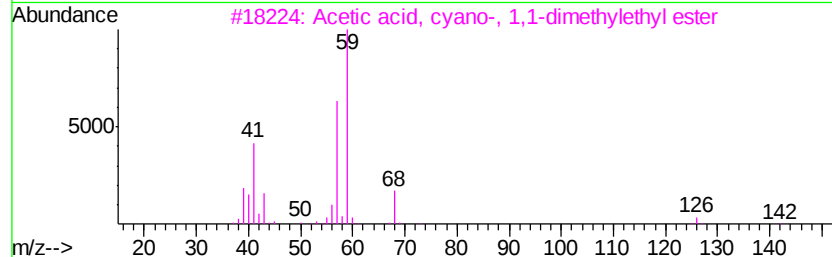
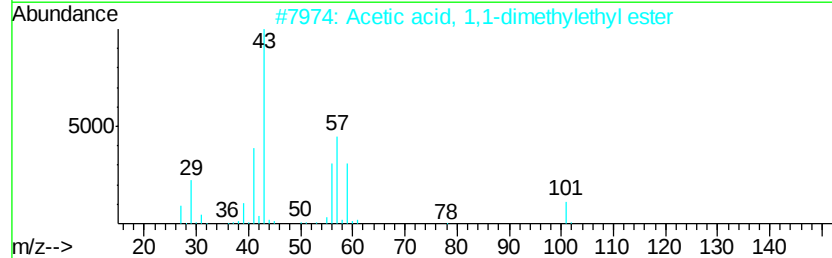
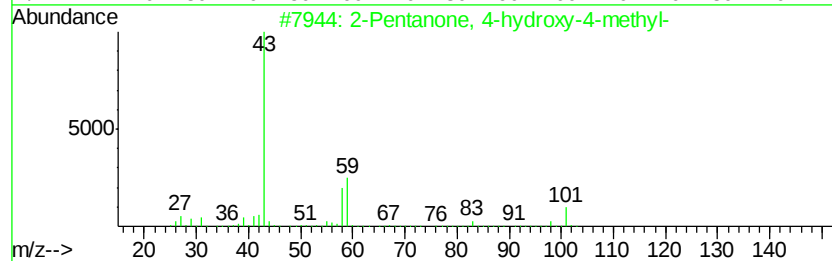
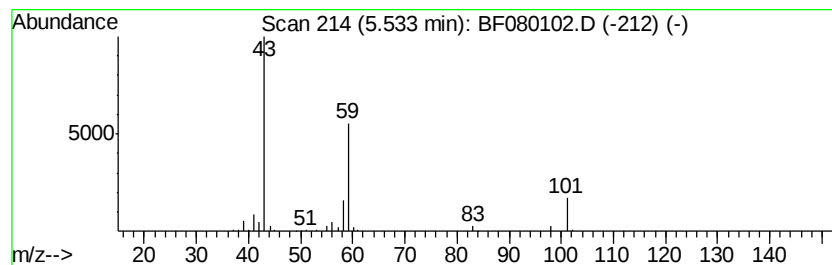
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.53	16.35 ng	229636	1,4-Dichlorobenzene-d4	7.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	37
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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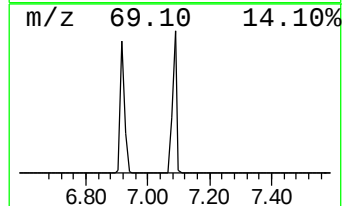
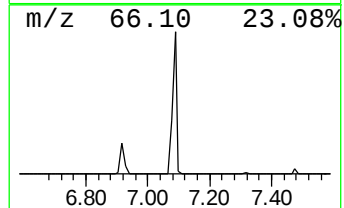
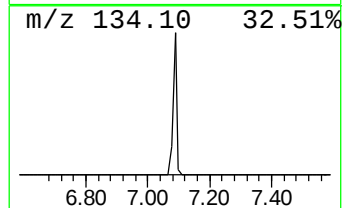
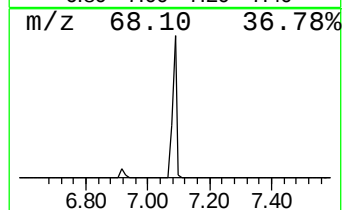
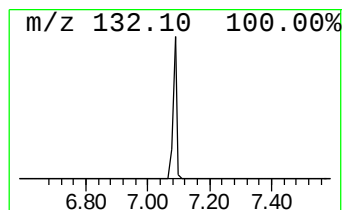
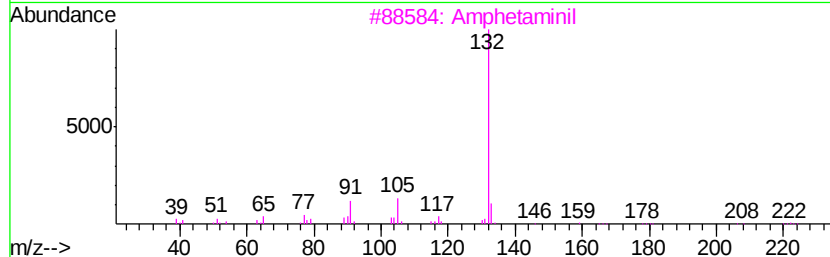
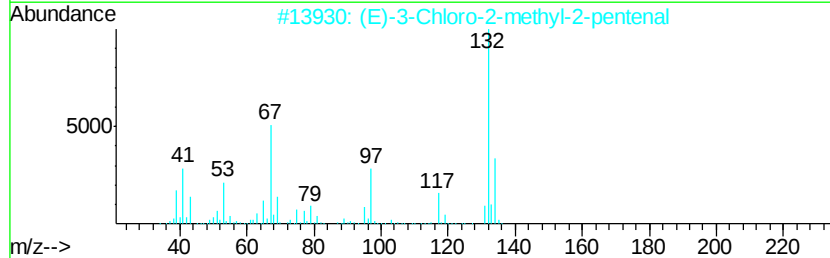
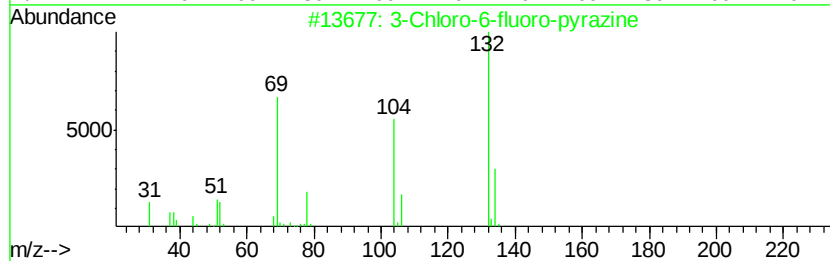
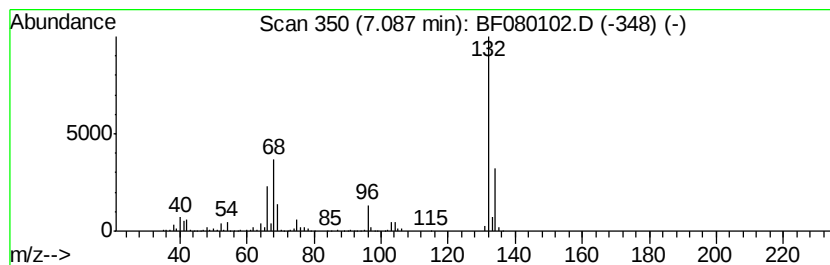
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Peak Number 2 unknown7.09 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.09	69.14 ng	971012	1,4-Dichlorobenzene-d4	7.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	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...	5.53	16.4	ng	229636	1	7.32	280903	20.0
unknown7.09	7.09	69.1	ng	971012	1	7.32	280903	20.0