

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062215\
 Data File : BF080096.D
 Acq On : 22 Jun 2015 16:10
 Operator : TP/IZ
 Sample : PB84131BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84131BL

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-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	217	rBV	486249	477317	22.89%	4.042%
2	5.933	247	249	251	rBV	796571	761499	36.52%	6.449%
3	6.916	333	335	338	rBV	767282	737634	35.37%	6.247%
4	7.087	348	350	352	rBV	957740	814045	39.04%	6.894%
5	7.316	369	370	373	rBB	219871	283111	13.58%	2.398%
6	7.476	382	384	387	rBV	1014494	1060174	50.84%	8.979%
7	7.876	417	419	421	rBV	1081790	851296	40.82%	7.210%
8	8.607	481	483	486	rBV	364342	379247	18.19%	3.212%
9	9.693	575	578	580	rBV	1406603	1696982	81.38%	14.372%
10	10.379	637	638	641	rVB	365842	449663	21.56%	3.808%
11	11.179	706	708	710	rBV	589067	577282	27.68%	4.889%
12	11.899	769	771	774	rBV	602952	511038	24.51%	4.328%
13	13.648	922	924	927	rBV	1947972	2085245	100.00%	17.660%
14	14.985	1038	1041	1044	rBV	519627	546242	26.20%	4.626%
15	16.940	1209	1212	1223	rBV	369030	550383	26.39%	4.661%
16	18.814	1371	1376	1380	rBV5	6394	26558	1.27%	0.225%

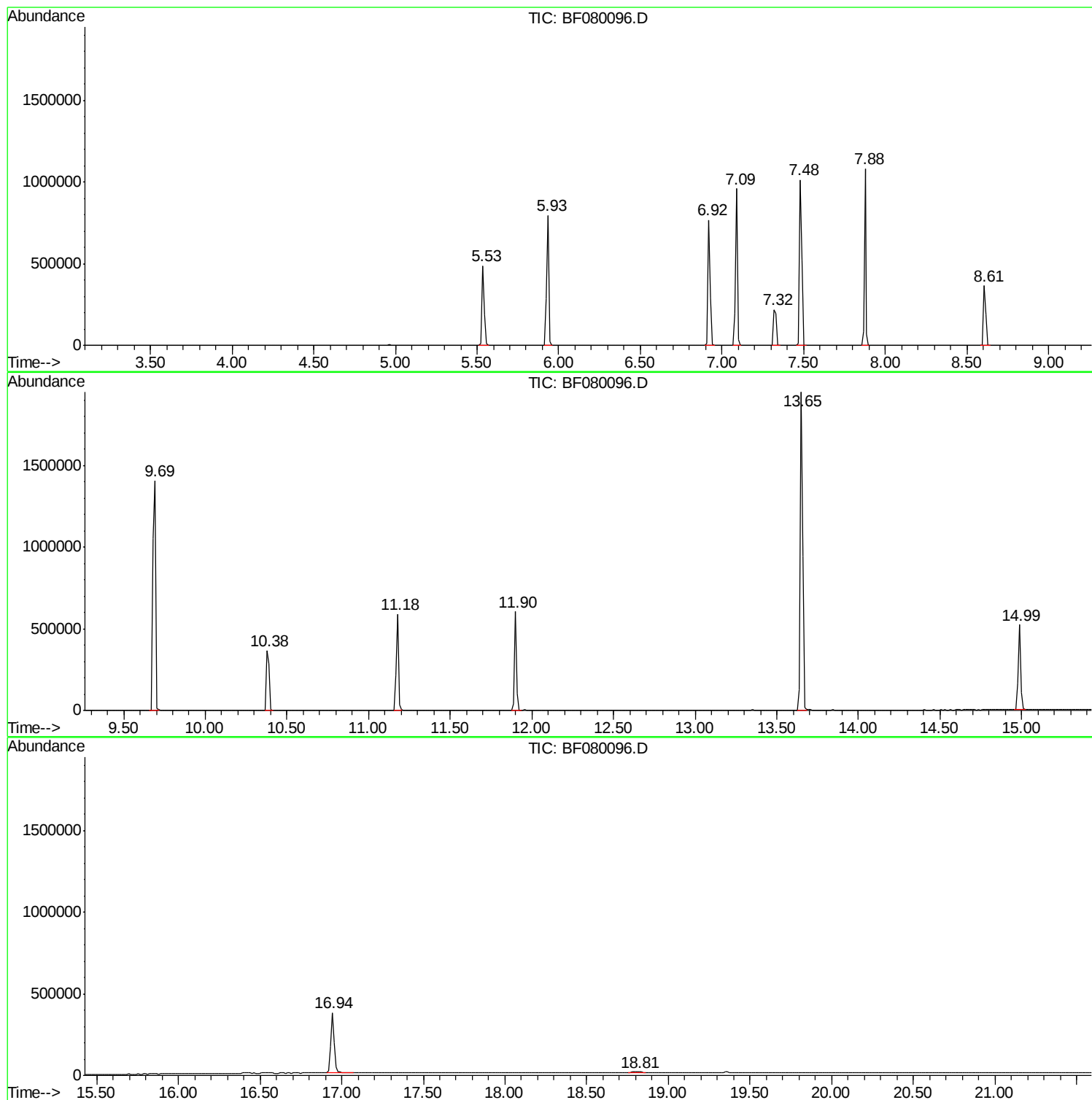
Sum of corrected areas: 11807716

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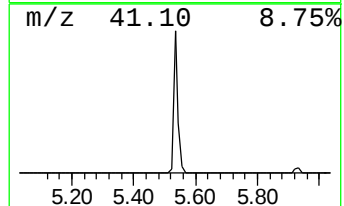
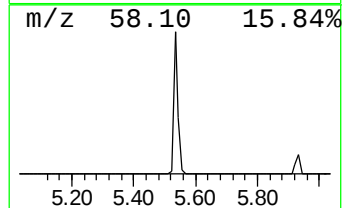
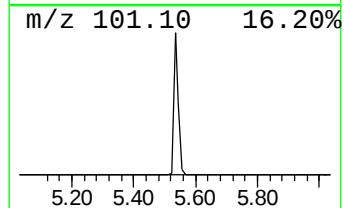
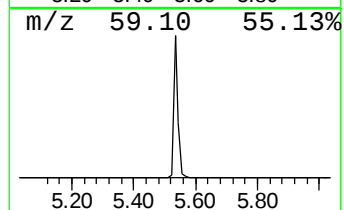
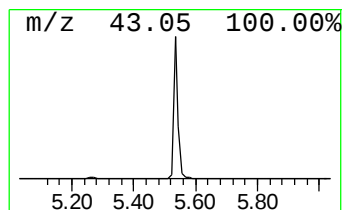
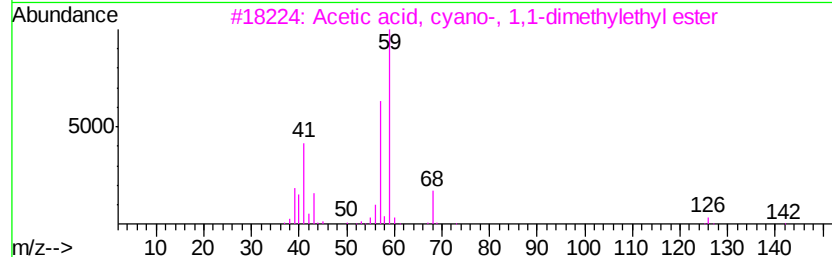
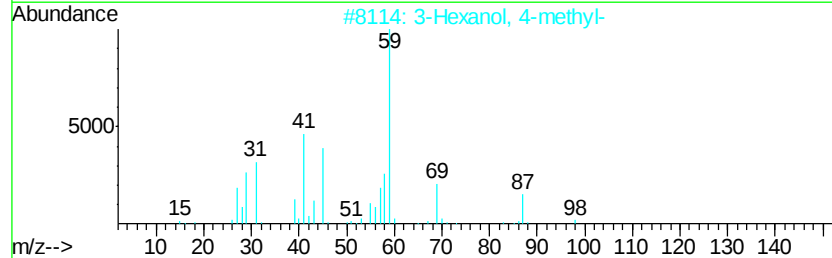
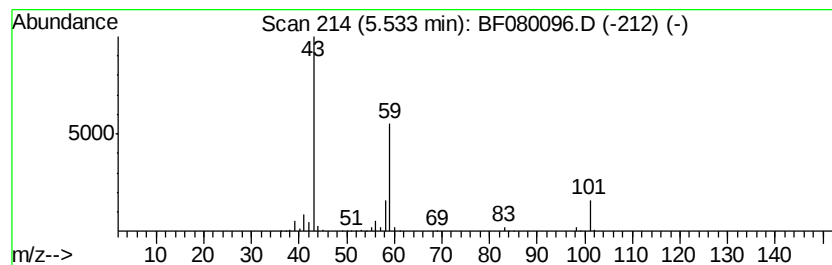
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.53	33.72 ng	477317	1,4-Dichlorobenzene-d4	7.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	17



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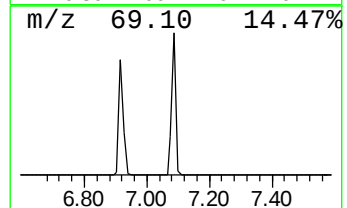
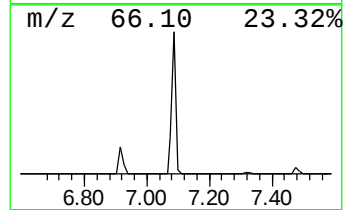
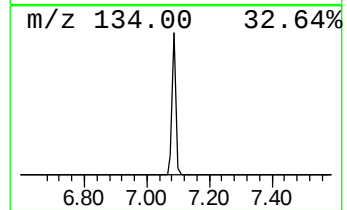
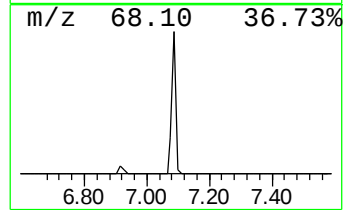
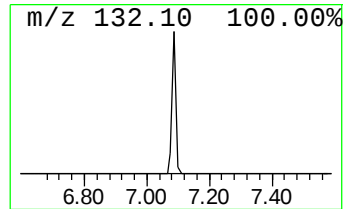
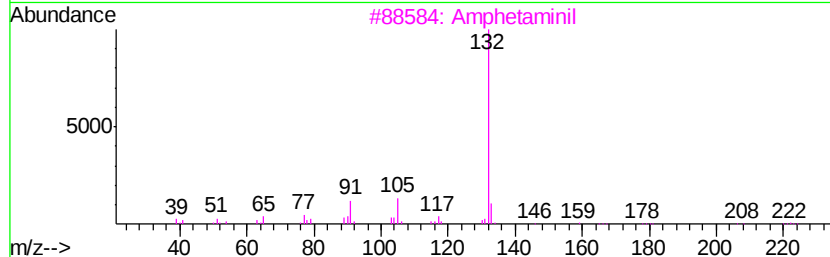
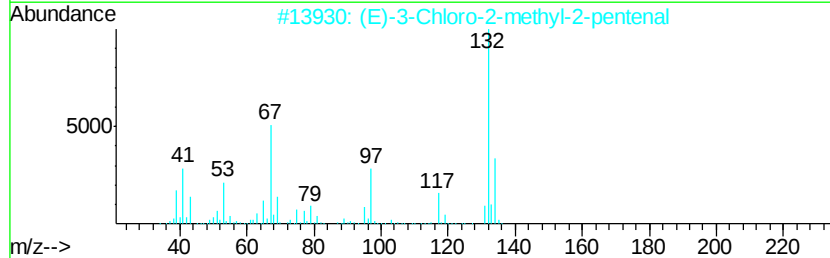
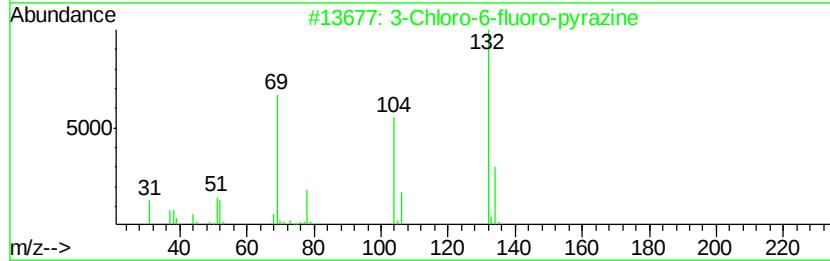
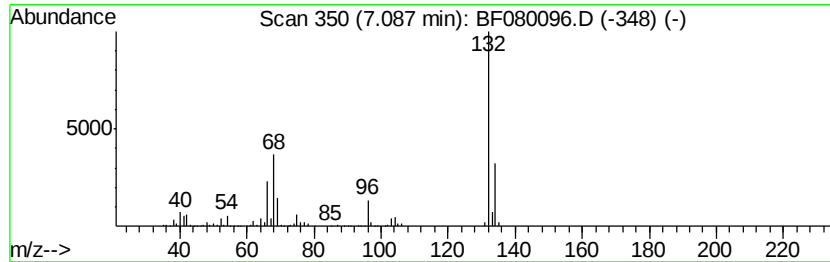
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Peak Number 2 unknown7.09 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.09	57.51 ng	814045	1,4-Dichlorobenzene-d4	7.33

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		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.53	33.7	ng	477317	1	7.33	283111	20.0
unknown7.09	7.09	57.5	ng	814045	1	7.33	283111	20.0