

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
 Data File : BF080240.D
 Acq On : 30 Jun 2015 17:05
 Operator : TP/IZ
 Sample : PB84281BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84281BL

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	5.498	209	211	219	rVB	110794	111298	8.37%	1.222%
2	5.898	243	246	248	rBV	654085	837788	62.99%	9.195%
3	6.893	330	333	335	rBV	773516	853857	64.20%	9.372%
4	7.053	345	347	349	rBV	924898	901499	67.78%	9.894%
5	7.281	365	367	370	rVB	263156	222404	16.72%	2.441%
6	7.441	379	381	383	rBV	902970	720538	54.18%	7.908%
7	7.842	413	416	418	rBV	661162	547373	41.16%	6.008%
8	8.573	478	480	483	rBV	388568	309555	23.28%	3.398%
9	9.647	572	574	577	rBV	1259036	1093592	82.23%	12.003%
10	10.345	633	635	638	rBV	438627	357547	26.88%	3.924%
11	11.133	702	704	707	rBV2	526845	649549	48.84%	7.129%
12	11.853	764	767	769	rBV	279238	370635	27.87%	4.068%
13	13.511	910	912	915	rBV	1186366	1329989	100.00%	14.597%
14	14.756	1018	1021	1024	rBV	409195	409615	30.80%	4.496%
15	16.585	1178	1181	1188	rBV	260039	395917	29.77%	4.345%

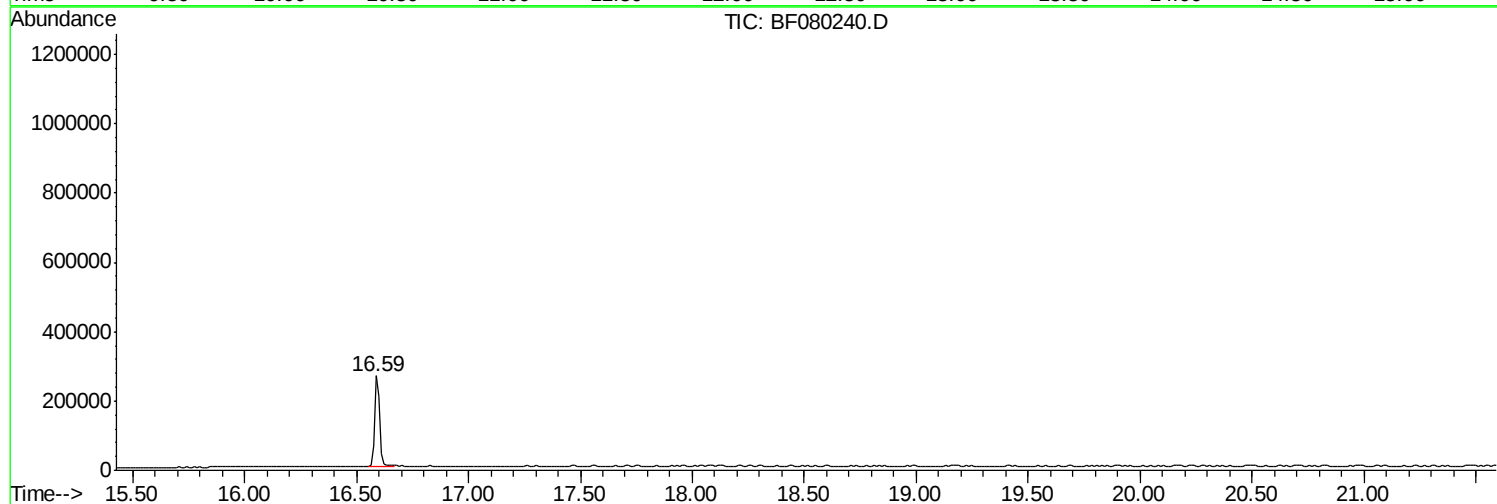
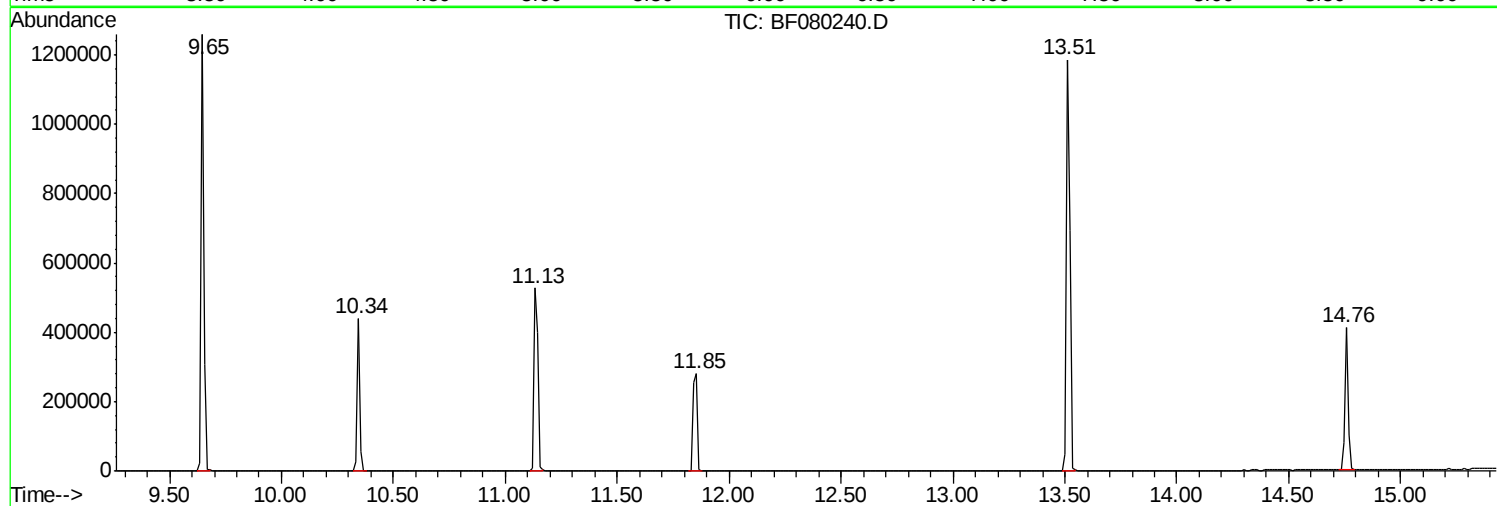
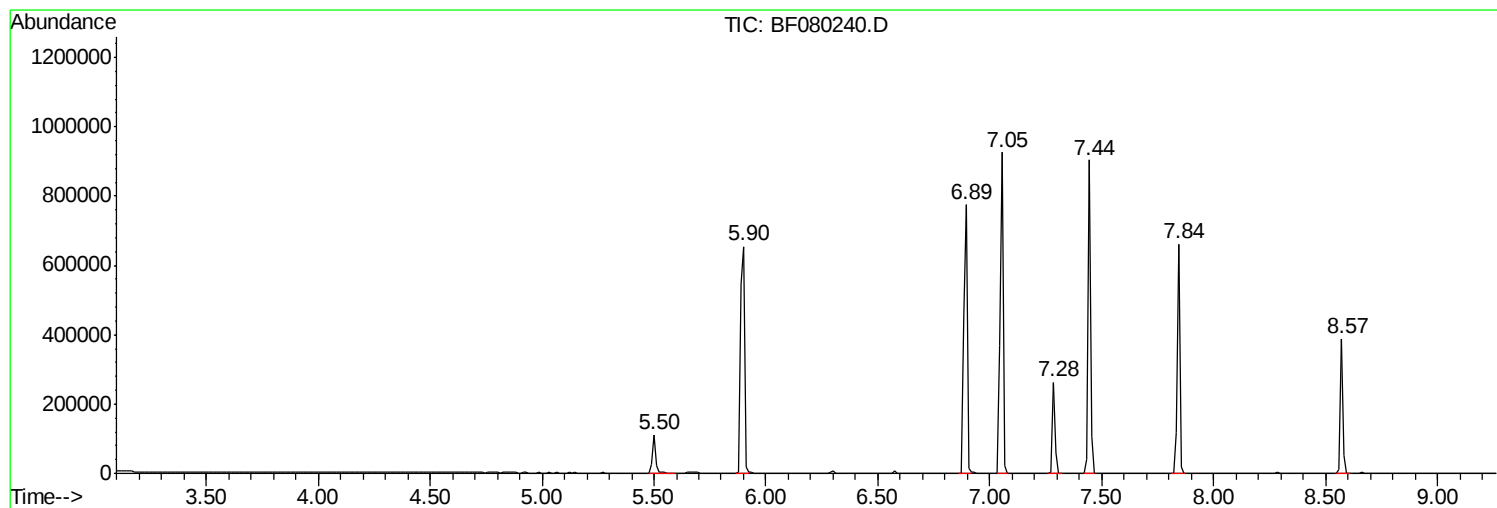
Sum of corrected areas: 9111156

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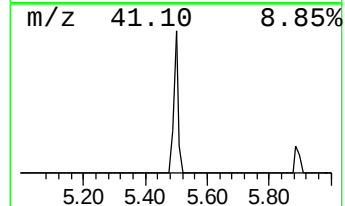
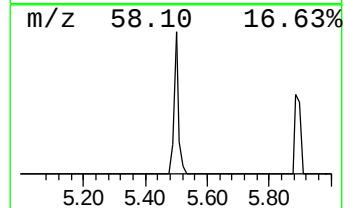
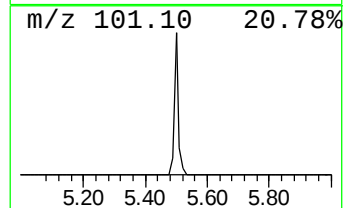
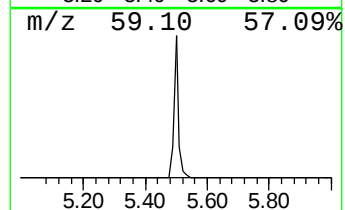
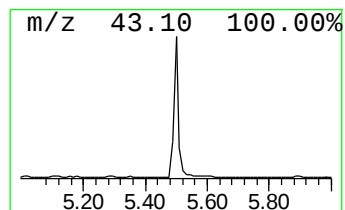
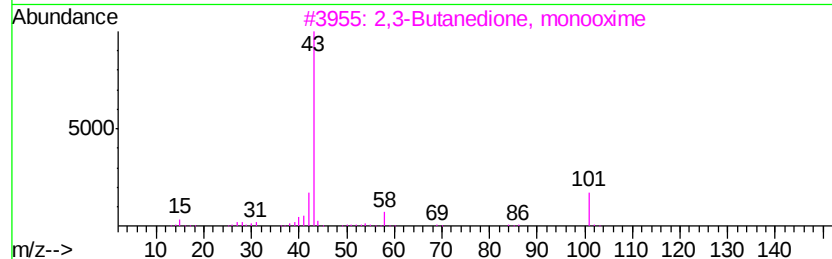
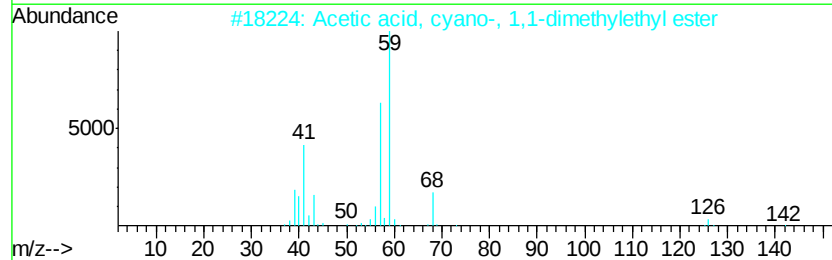
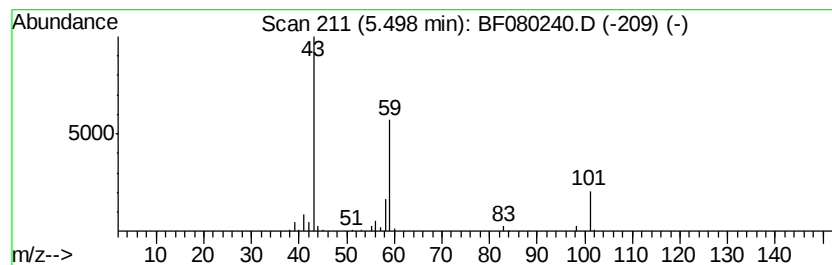
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.50	10.01 ng	111298	1,4-Dichlorobenzene-d4	7.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9



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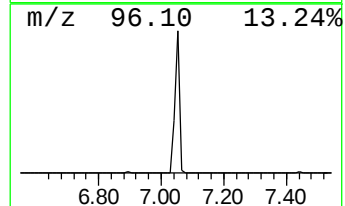
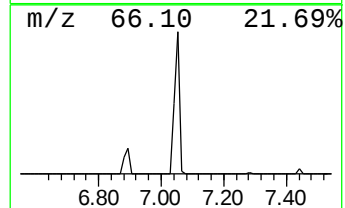
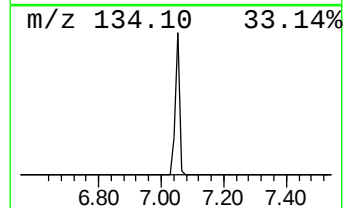
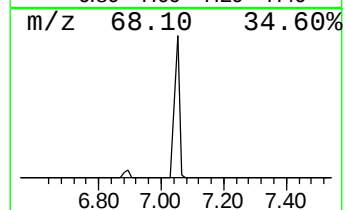
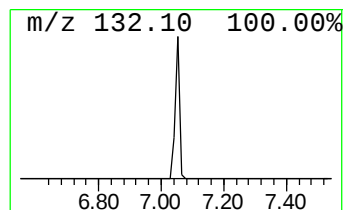
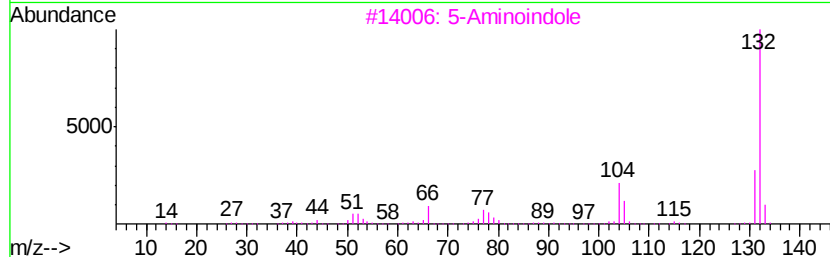
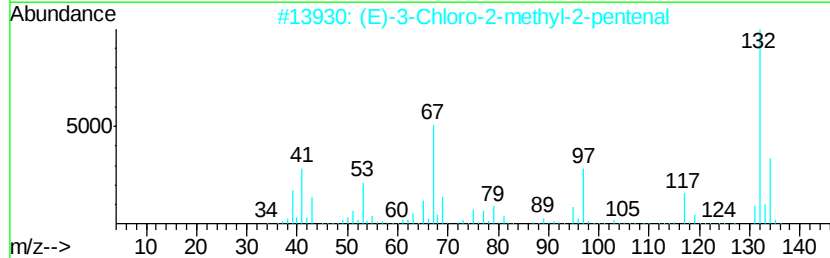
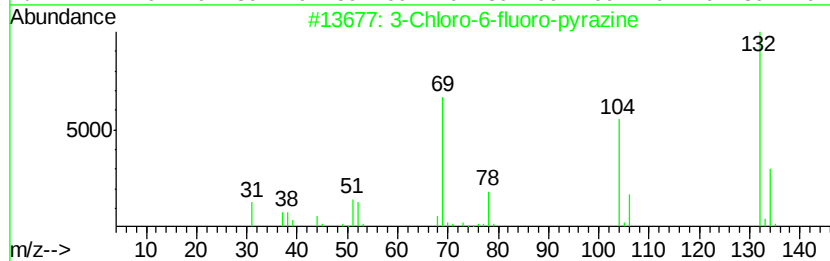
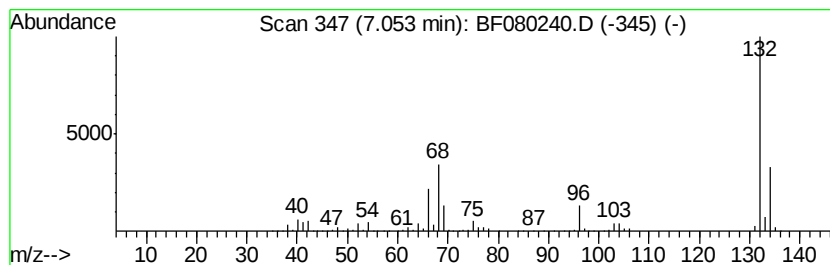
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Peak Number 2 unknown7.05 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	81.07 ng	901499	1,4-Dichlorobenzene-d4	7.28

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		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	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...	5.50	10.0	ng	111298	1	7.28	222404	20.0
unknown7.05	7.05	81.1	ng	901499	1	7.28	222404	20.0