

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100615\
 Data File : BF082078.D
 Acq On : 6 Oct 2015 12:04
 Operator : UM/IZ
 Sample : PB85739BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85739BL

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-BF092815.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	5.074	254	257	263	rBV	607835	877458	33.54%	4.715%
2	5.486	290	293	295	rBV	1696307	1751952	66.97%	9.415%
3	6.515	380	383	385	rBV	1665821	1649305	63.05%	8.863%
4	6.652	392	395	397	rBV	1862173	1900243	72.64%	10.212%
5	6.880	413	415	418	rBV	486445	431263	16.49%	2.318%
6	7.040	426	429	431	rBV	1592037	1510688	57.75%	8.118%
7	7.452	462	465	467	rBV	1032571	1079401	41.26%	5.801%
8	8.172	525	528	530	rBV	552389	548430	20.97%	2.947%
9	9.246	619	622	624	rBV	2339408	2158643	82.52%	11.600%
10	9.921	679	681	684	rBV	565218	646111	24.70%	3.472%
11	10.721	748	751	753	rBV	1168313	1392154	53.22%	7.481%
12	11.418	809	812	814	rBV	524981	651070	24.89%	3.499%
13	13.007	948	951	953	rBV	2409009	2615901	100.00%	14.058%
14	14.058	1040	1043	1045	rBV	684152	730625	27.93%	3.926%
15	15.510	1167	1170	1179	rBV	478074	664972	25.42%	3.574%

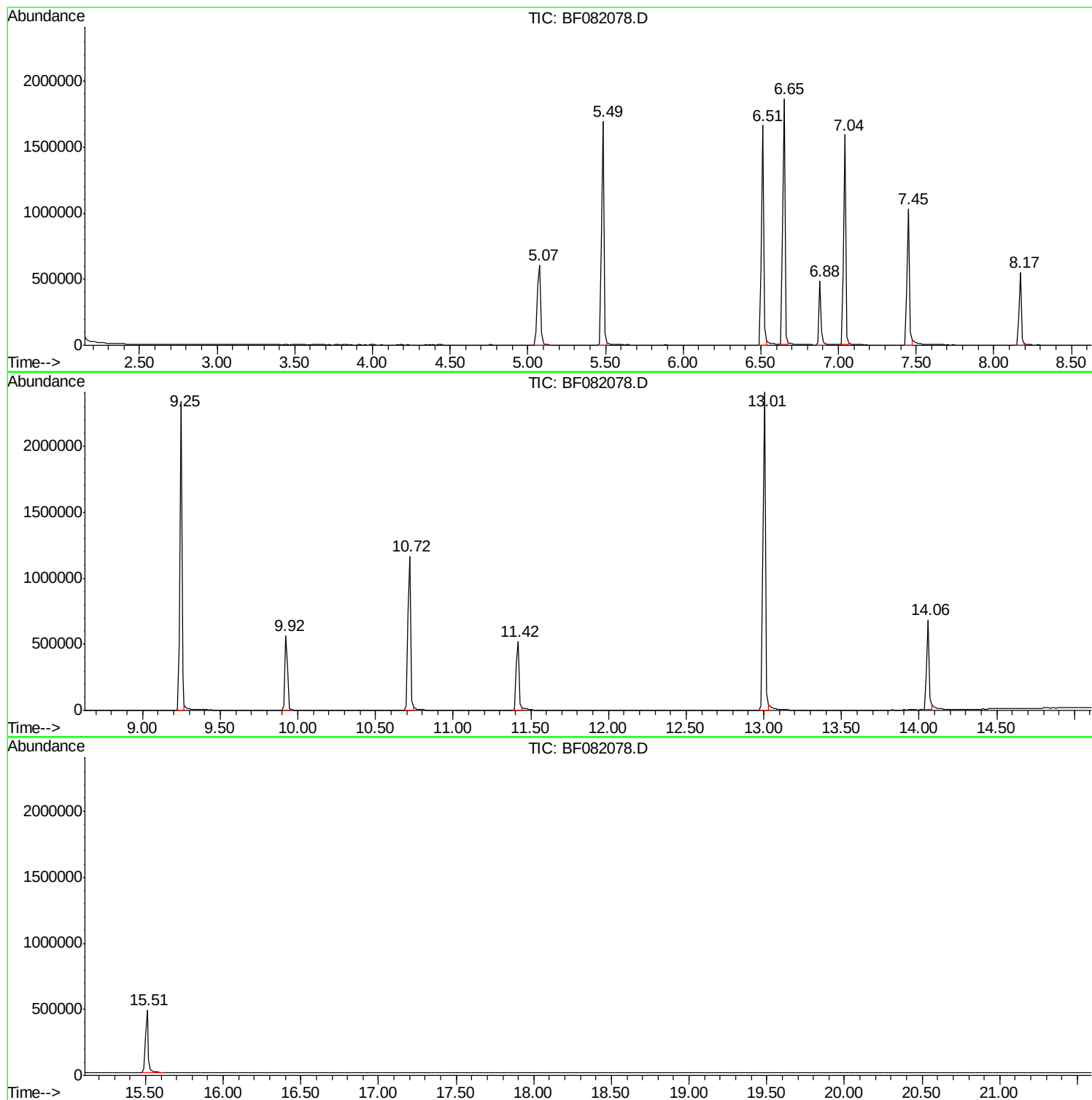
Sum of corrected areas: 18608216

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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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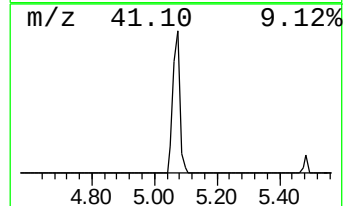
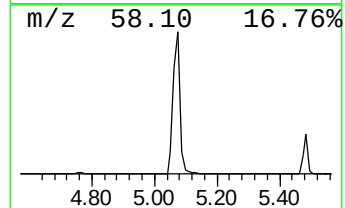
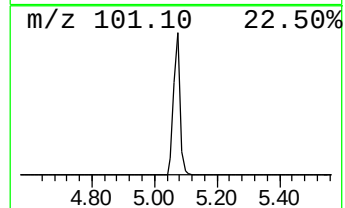
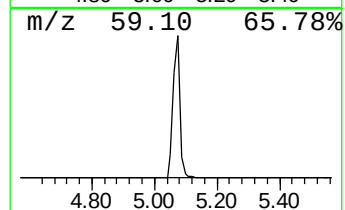
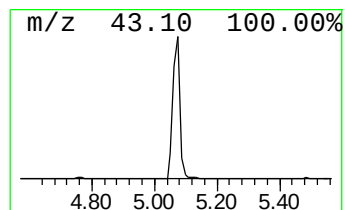
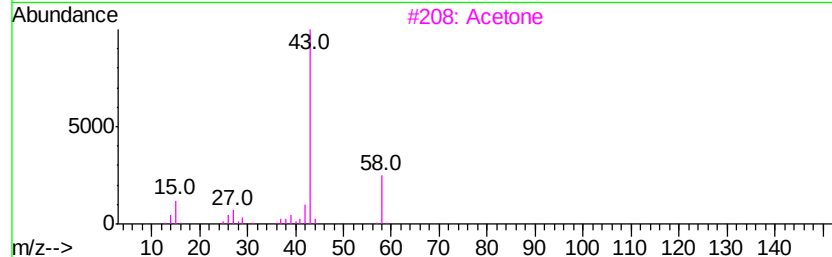
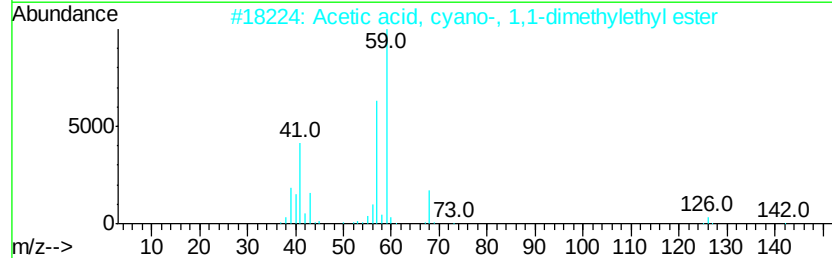
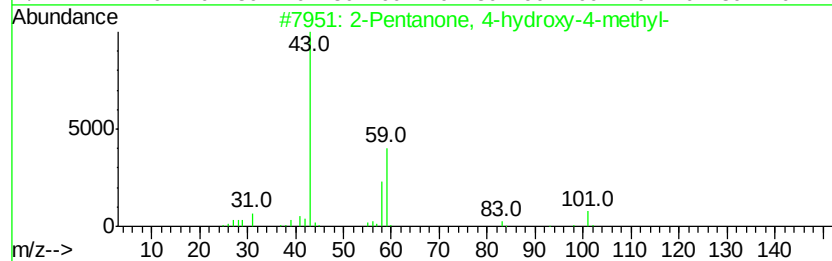
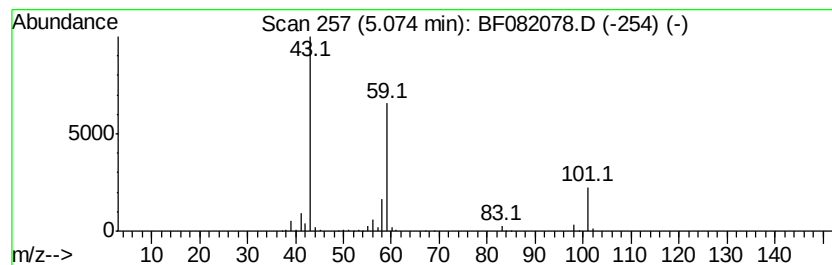
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	40.69 ng	877458	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		Acetone	58	C3H6O	000067-64-1	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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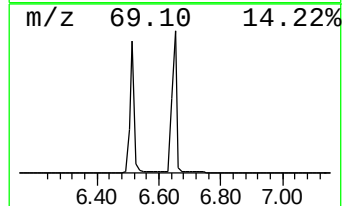
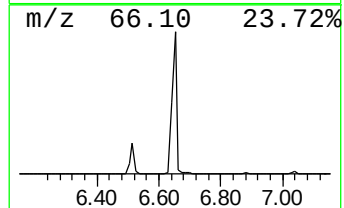
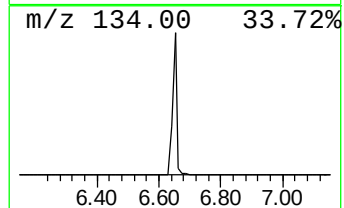
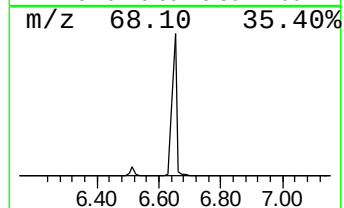
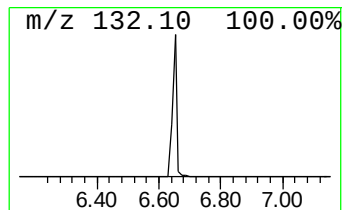
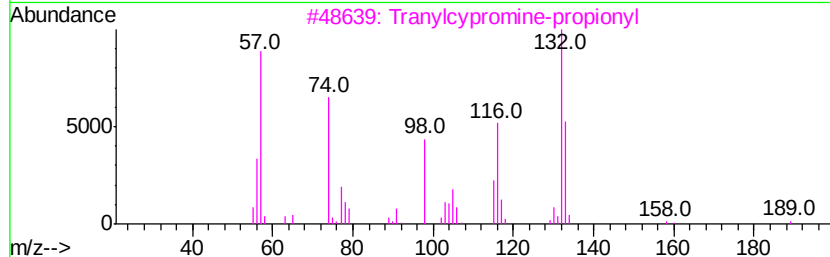
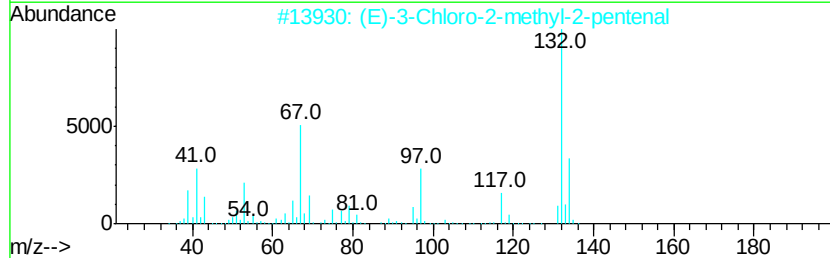
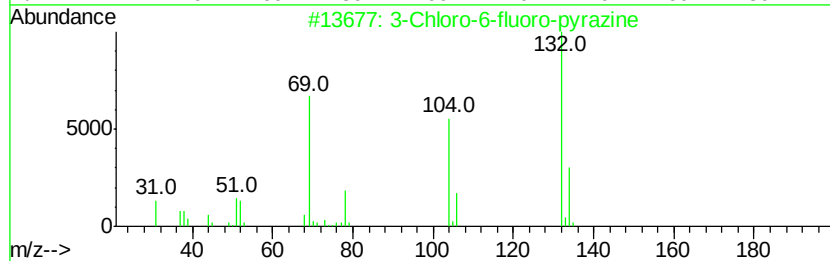
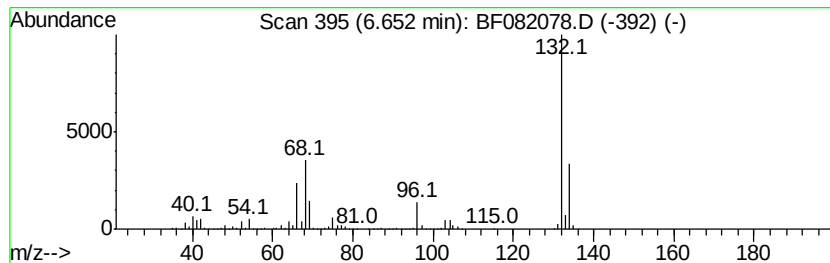
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Peak Number 2 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	88.12 ng	1900240	1,4-Dichlorobenzene-d4	6.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	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...	5.07	40.7	ng	877458	1	6.88	431263	20.0
unknown6.65	6.65	88.1	ng	1900240	1	6.88	431263	20.0