

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061216\
 Data File : BF087952.D
 Acq On : 12 Jun 2016 17:06
 Operator : UM/SJ
 Sample : H3490-03
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JAMAICA-FB2-6-06-16

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-BF060716.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	1.736	11	13	22	rVB	121742	211801	6.49%	0.864%
2	1.861	22	24	62	rVB	1317766	2338606	71.65%	9.540%
3	4.959	293	295	299	rBV	205551	234090	7.17%	0.955%
4	5.382	330	332	335	rBV	1030413	1234438	37.82%	5.036%
5	6.422	421	423	426	rBV	600127	797777	24.44%	3.254%
6	6.570	433	436	438	rBV	1604365	2301567	70.51%	9.389%
7	6.799	453	456	458	rBV	752625	637202	19.52%	2.599%
8	6.959	467	470	472	rVB	2094688	2195180	67.26%	8.955%
9	7.370	502	506	508	rBV	1607644	1945666	59.61%	7.937%
10	8.090	566	569	571	rVB	815211	864493	26.49%	3.527%
11	9.165	660	663	666	rBV	2727339	3147982	96.45%	12.842%
12	9.839	719	722	729	rBV	1098762	1027166	31.47%	4.190%
13	10.628	788	791	794	rBV	1709424	1881105	57.63%	7.674%
14	11.325	849	852	854	rBV	801752	1024483	31.39%	4.179%
15	12.914	988	991	993	rBV	2822784	3263942	100.00%	13.315%
16	13.954	1079	1082	1084	rBV	761848	791207	24.24%	3.228%
17	15.360	1202	1205	1221	rBV	324520	617224	18.91%	2.518%

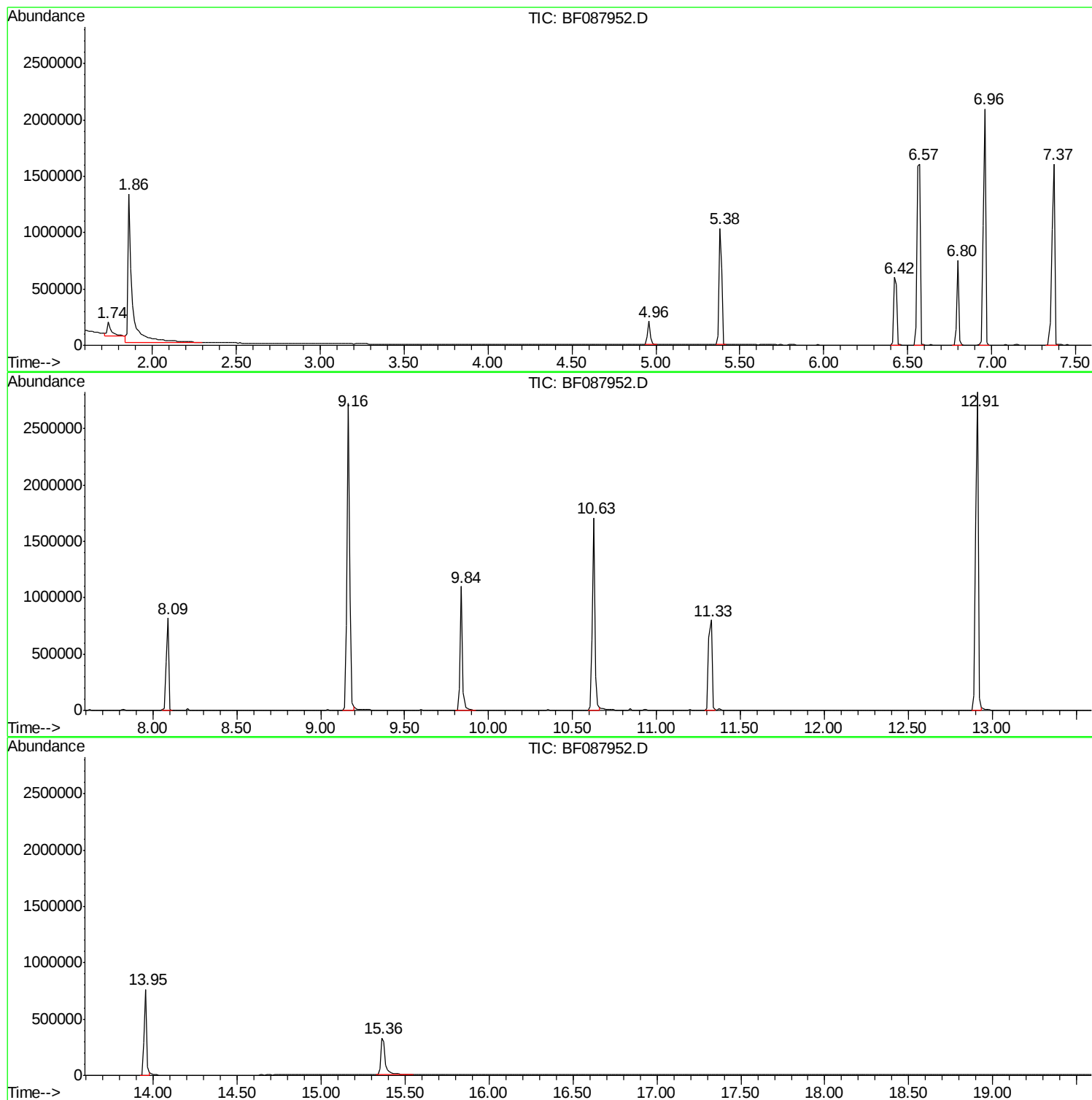
Sum of corrected areas: 24513929

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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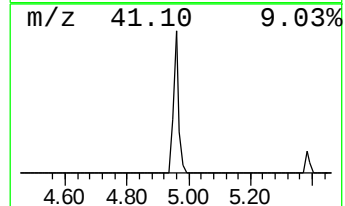
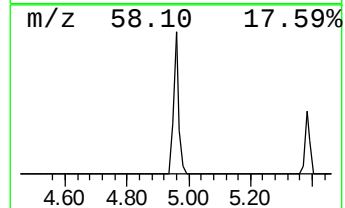
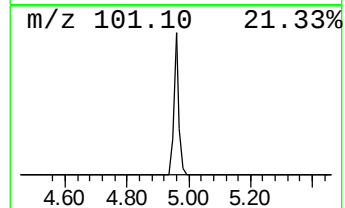
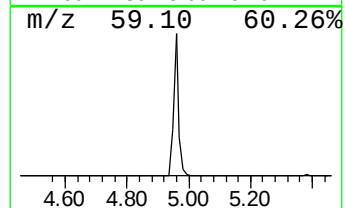
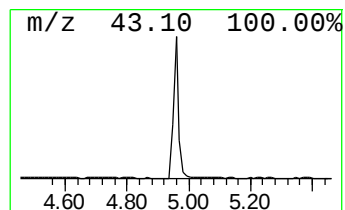
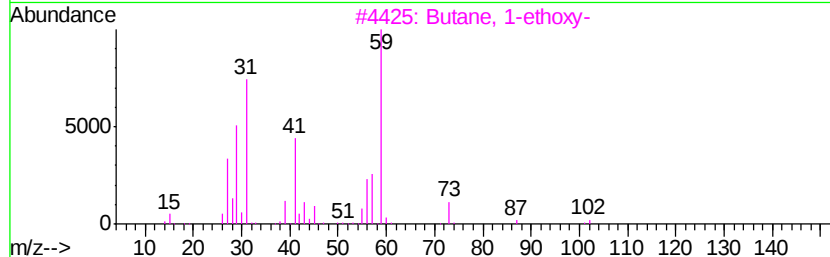
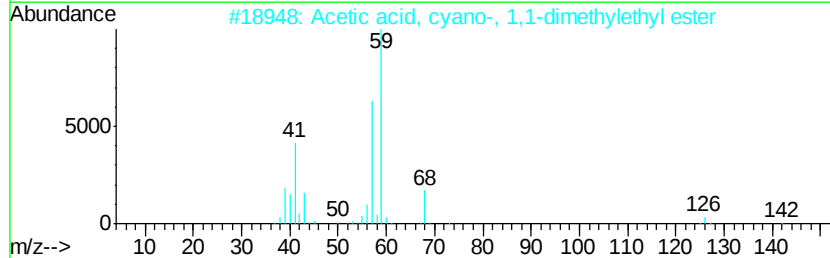
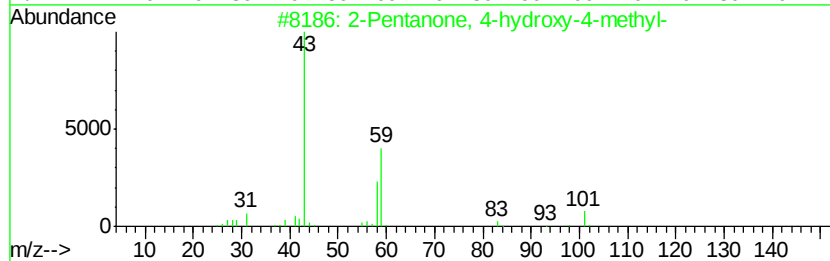
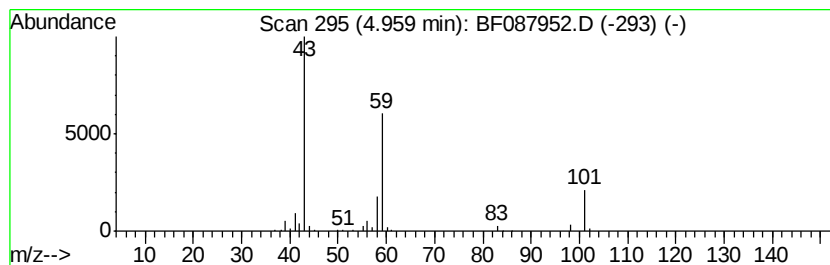
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	7.35 ng	234090	1,4-Dichlorobenzene-d4	6.80

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		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Acetone	58	C3H6O	000067-64-1	9



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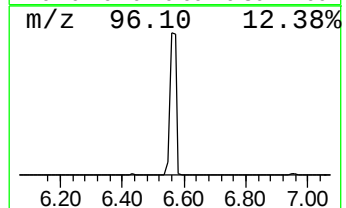
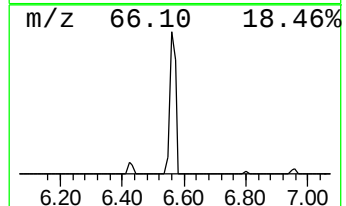
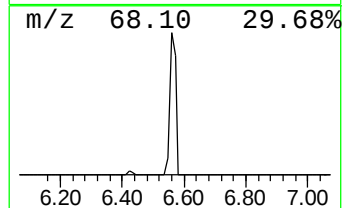
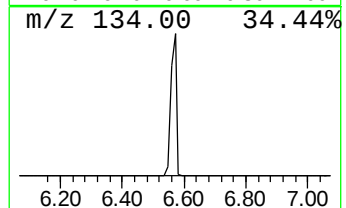
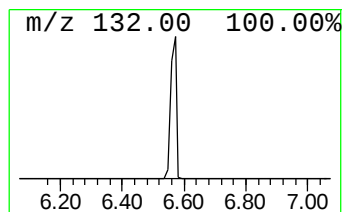
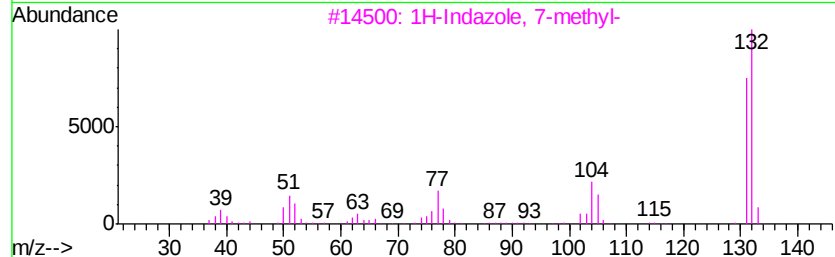
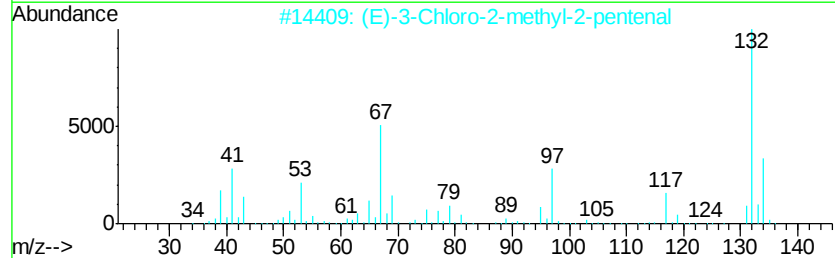
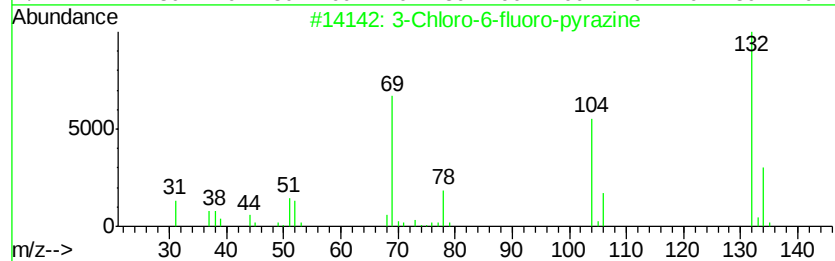
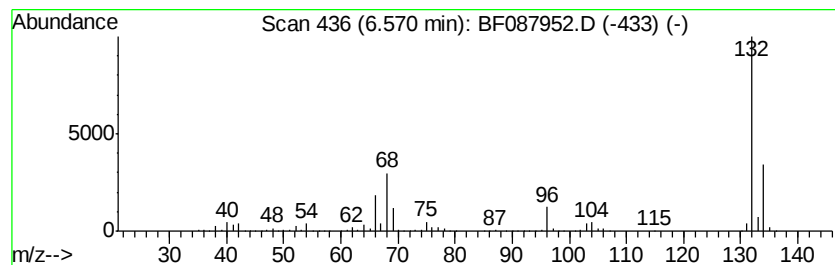
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Peak Number 4 3-Chloro-6-fluoro-pyrazine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	72.24 ng	2301570	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	52
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	28
3		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	25
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	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...	4.96	7.3	ng	234090	1	6.80	637202	20.0
3-Chloro-6-fluoro...	6.57	72.2	ng	2301570	1	6.80	637202	20.0