

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082416\
 Data File : BF089877.D
 Acq On : 24 Aug 2016 16:28
 Operator : UM/SJ
 Sample : PB93061BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB93061BL

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-BF081916.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	1.655	3	6	14	rBV	1878549	4252739	22.97%	3.893%
2	1.770	14	16	73	rVB	7427358	18512480	100.00%	16.947%
3	4.810	279	282	288	rBV	1076725	1931270	10.43%	1.768%
4	5.244	317	320	323	rBV	7344146	8539825	46.13%	7.818%
5	6.307	410	413	415	rBV	8572582	9462701	51.12%	8.663%
6	6.433	421	424	426	rBV	9746689	9466739	51.14%	8.666%
7	6.662	442	444	447	rBV	2313115	2332803	12.60%	2.136%
8	6.822	455	458	460	rBV	8619683	7315139	39.51%	6.697%
9	7.233	491	494	496	rBV	6042658	5945113	32.11%	5.442%
10	7.953	554	557	559	rBV	3820765	3359946	18.15%	3.076%
11	9.039	649	652	654	rBV	9214736	10747302	58.05%	9.839%
12	9.702	708	710	712	rBV	4117762	3454238	18.66%	3.162%
13	10.491	776	779	781	rBV	5215035	5793387	31.29%	5.304%
14	11.176	837	839	842	rBV	3847682	3366684	18.19%	3.082%
15	12.777	976	979	981	rBV	9067260	9543374	51.55%	8.737%
16	13.851	1070	1073	1075	rBV	3124553	2972415	16.06%	2.721%
17	15.348	1201	1204	1207	rBV	1967015	2238857	12.09%	2.050%

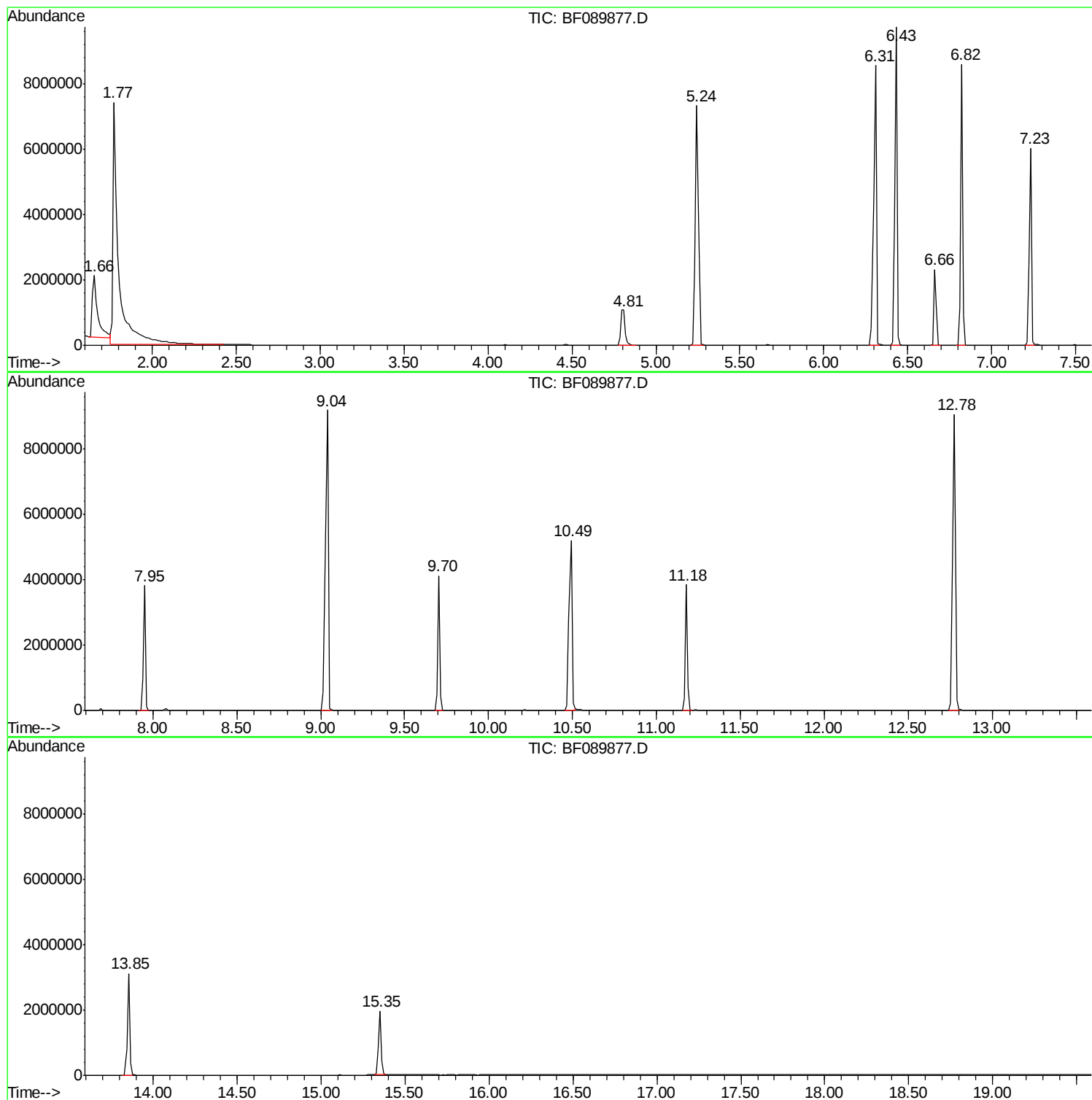
Sum of corrected areas: 109235012

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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.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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Instrument :
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ClientSampled :
PB93061BL

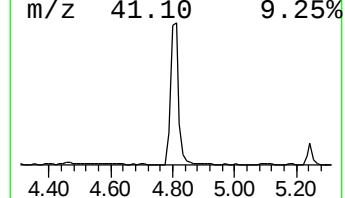
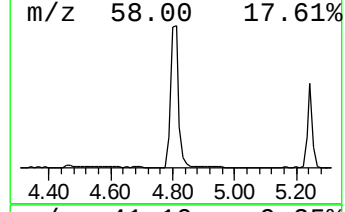
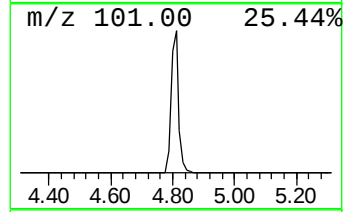
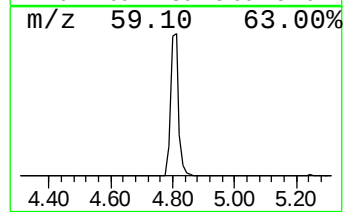
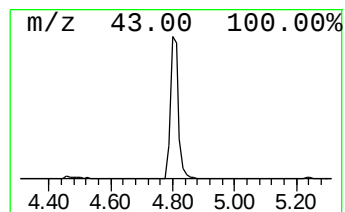
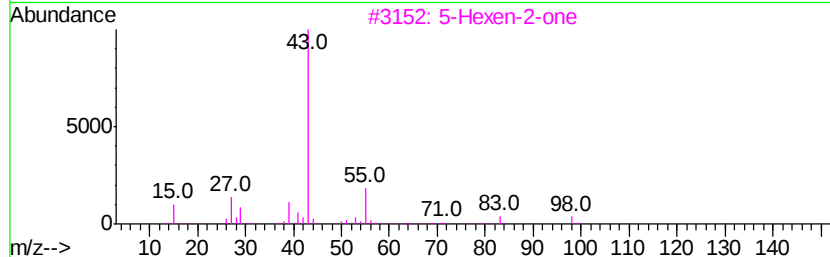
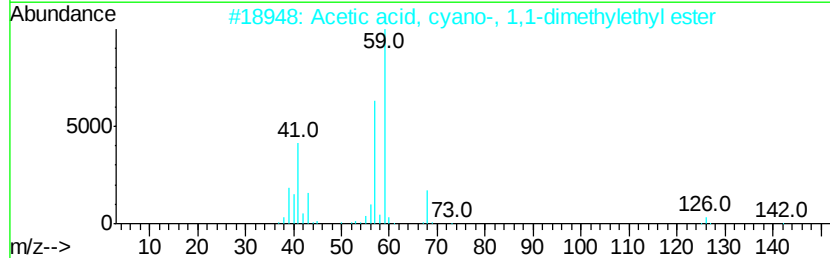
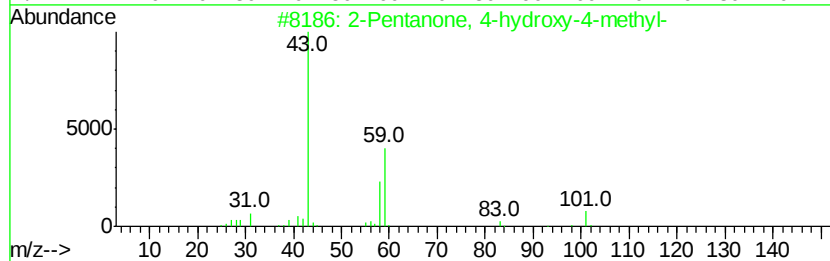
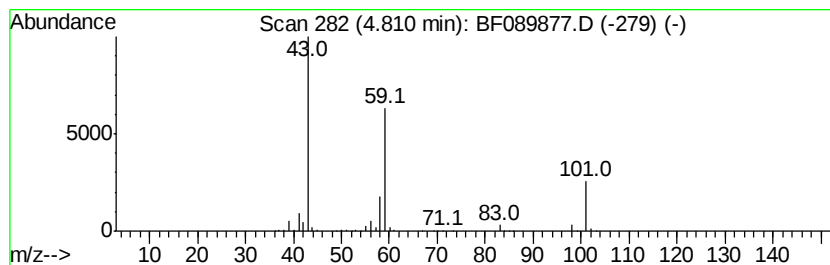
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.81	16.56 ng	1931270	1,4-Dichlorobenzene-d4	6.66

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	25
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Acetone	58	C3H6O	000067-64-1	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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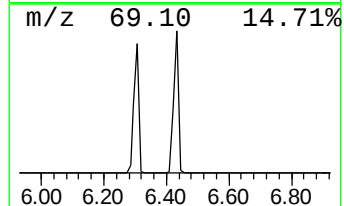
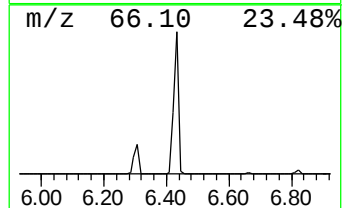
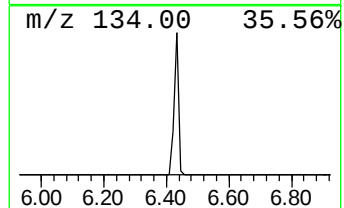
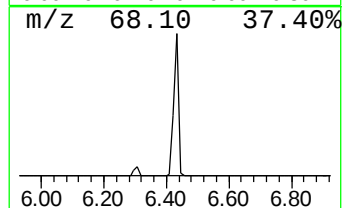
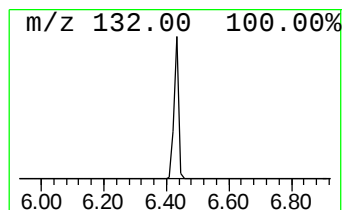
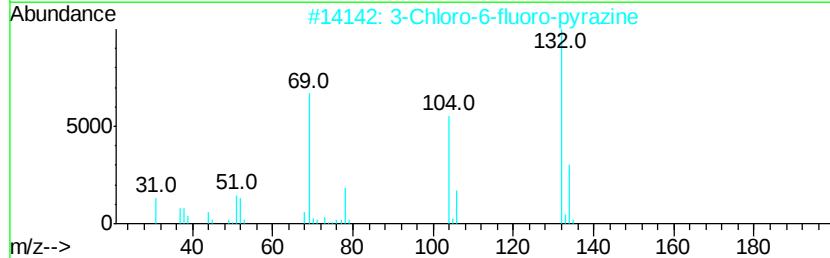
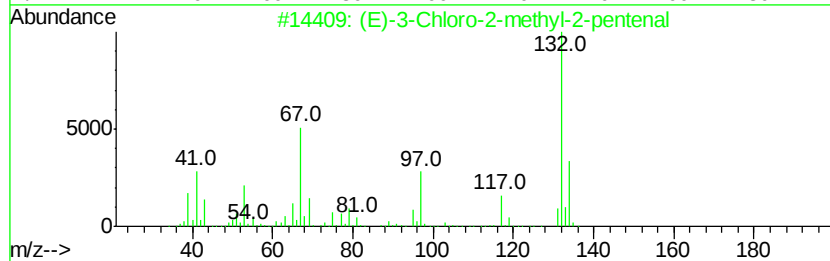
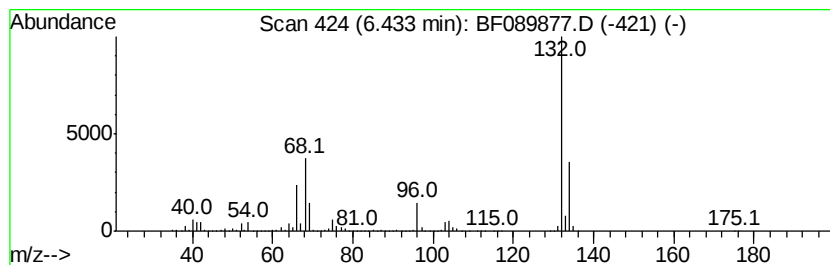
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Peak Number 4 unknown6.43 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.43	81.16 ng	9466740	1,4-Dichlorobenzene-d4	6.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
3			Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	12
4			Amphetaminil	250	C17H18N2	017590-01-1	12
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...	4.81	16.6	ng	1931270	1	6.66	2332800	20.0
unknown6.43	6.43	81.2	ng	9466740	1	6.66	2332800	20.0