

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081616\
 Data File : BF089747.D
 Acq On : 17 Aug 2016 6:42
 Operator : UM/SJ
 Sample : H4399-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FB080416

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-BF081616.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	1.667	6	7	16	rVV	1749143	3037106	8.51%	1.646%
2	1.793	16	18	73	rVB	13567244	35688821	100.00%	19.347%
3	4.833	282	284	287	rBV	1558351	2417470	6.77%	1.311%
4	5.279	320	323	325	rBV	7915452	8330967	23.34%	4.516%
5	6.330	412	415	416	rBV	3711879	4870902	13.65%	2.641%
6	6.467	424	427	429	rBV	11406482	15127972	42.39%	8.201%
7	6.696	445	447	449	rBV	5171253	4361708	12.22%	2.365%
8	6.856	458	461	463	rVB	14164117	13156943	36.87%	7.133%
9	7.267	493	497	499	rBV	10230435	11648737	32.64%	6.315%
10	7.988	557	560	562	rBV	6095265	6373417	17.86%	3.455%
11	9.073	651	655	657	rBV	16861278	21456830	60.12%	11.632%
12	9.736	710	713	715	rBV	8103170	7135384	19.99%	3.868%
13	10.525	778	782	784	rBV	9484069	11859380	33.23%	6.429%
14	11.211	839	842	845	rBV	7936552	6924161	19.40%	3.754%
15	12.548	957	959	962	rBV	1812026	1968129	5.51%	1.067%
16	12.822	979	983	985	rBV	15340008	18504527	51.85%	10.032%
17	13.920	1076	1079	1081	rBV	6633692	6471341	18.13%	3.508%
18	15.405	1206	1209	1212	rBV	4491377	5128987	14.37%	2.780%

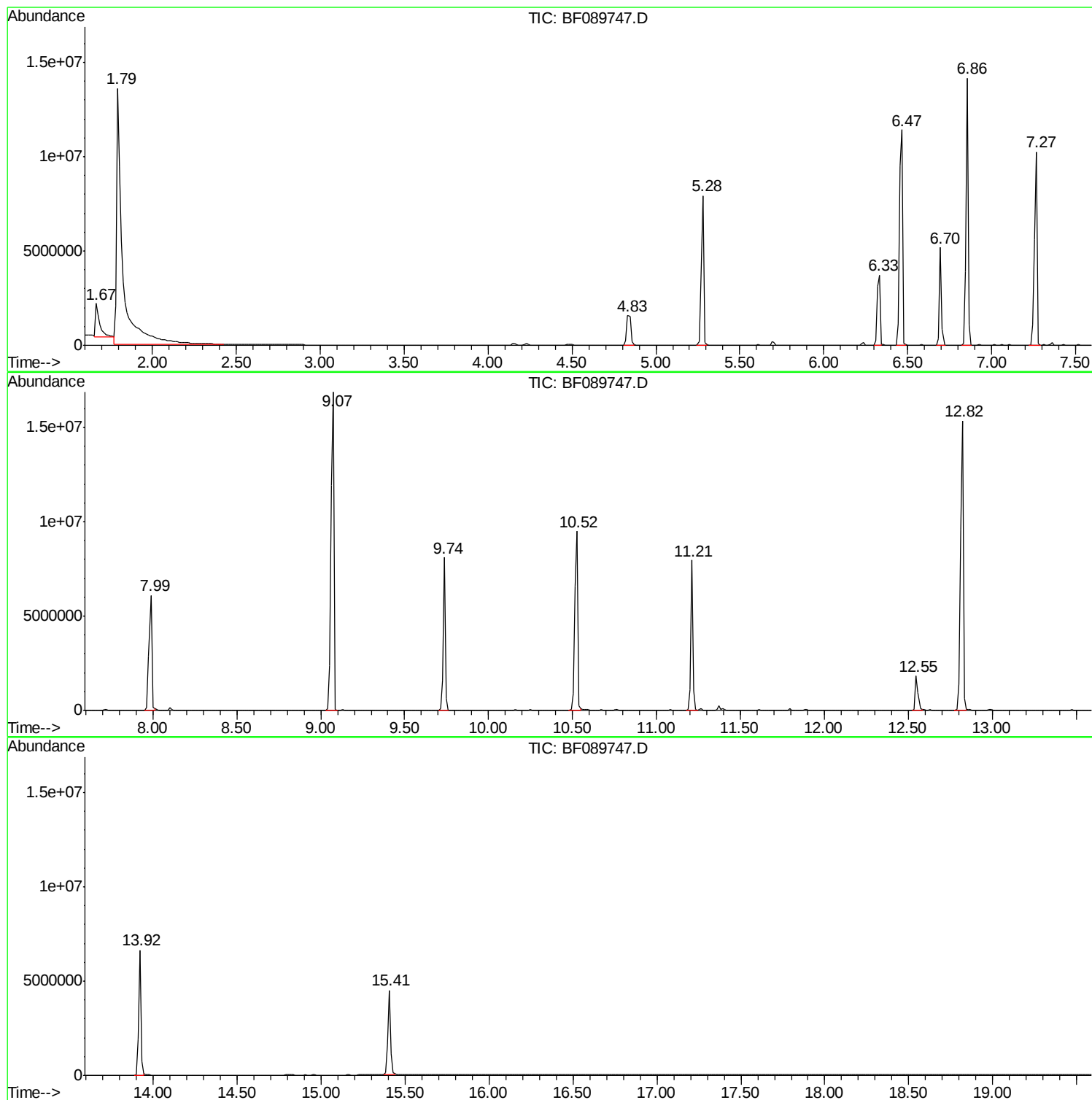
Sum of corrected areas: 184462782

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081616.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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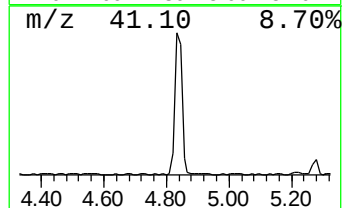
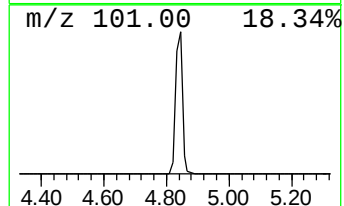
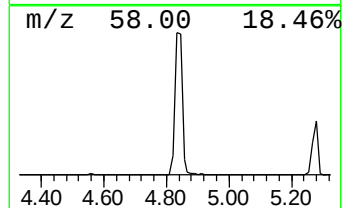
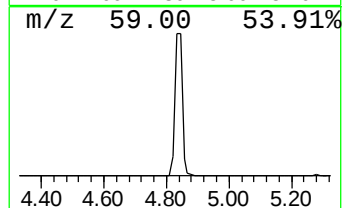
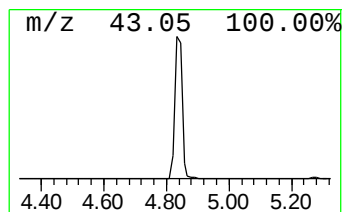
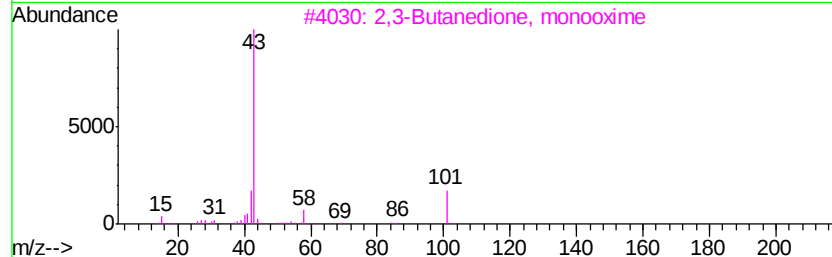
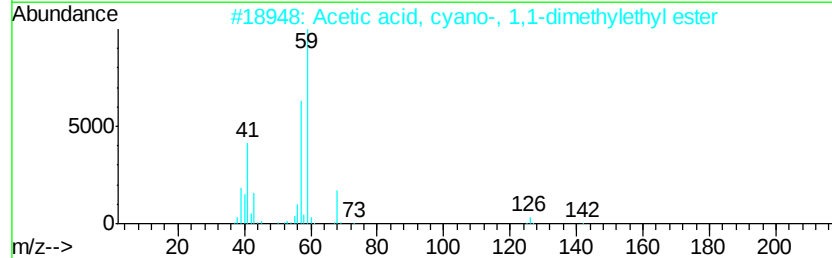
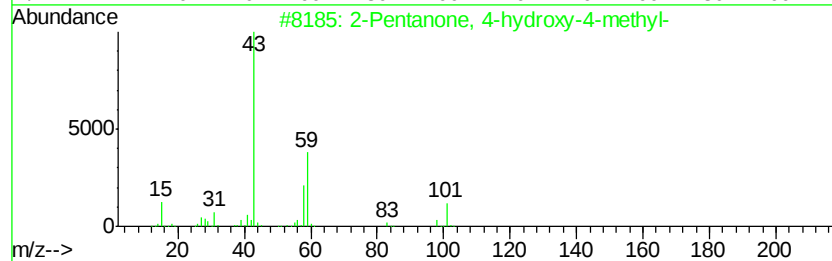
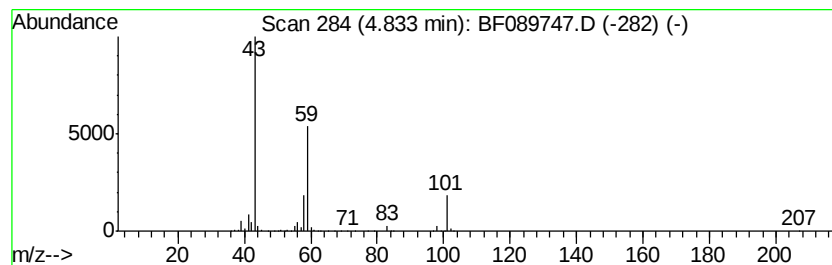
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	11.09 ng	2417470	1,4-Dichlorobenzene-d4	6.70

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59	
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25	
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16	
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10	
5	Butane	58	C4H10	000106-97-8	9	



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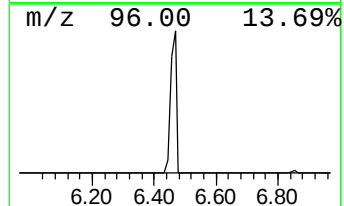
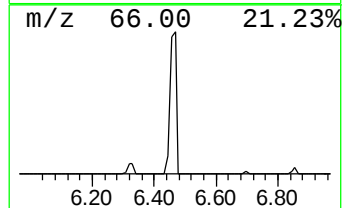
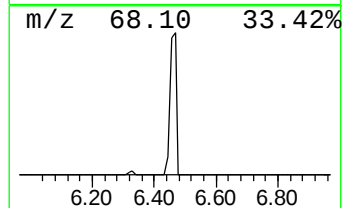
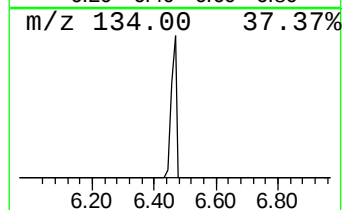
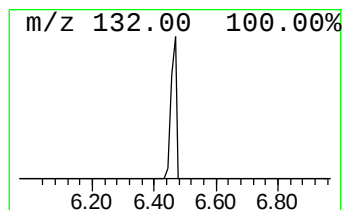
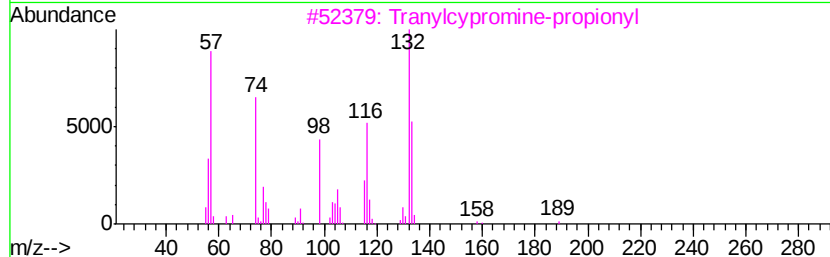
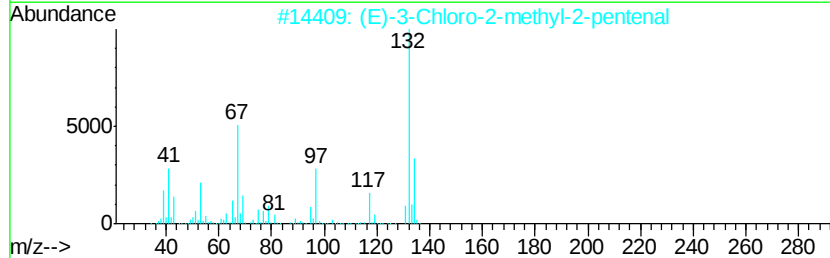
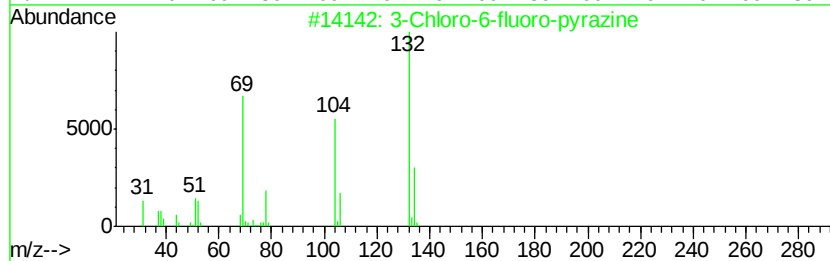
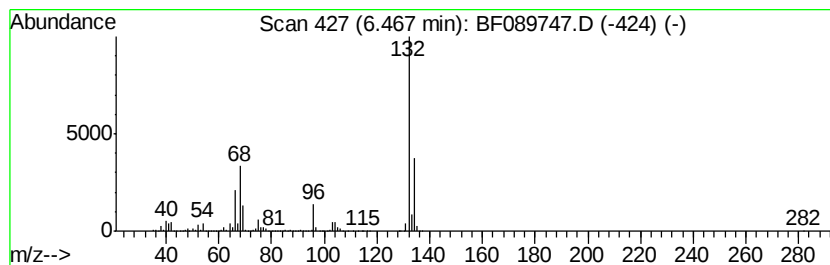
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Peak Number 4 unknown6.47 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	69.37 ng	15128000	1,4-Dichlorobenzene-d4	6.70

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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	16
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	12



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.83	11.1	ng	2417470	1	6.70	4361710	20.0
unknown6.47	6.47	69.4	ng	15128000	1	6.70	4361710	20.0