

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061116\
 Data File : BF087912.D
 Acq On : 11 Jun 2016 20:03
 Operator : UM/SJ
 Sample : H3387-05
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP16-LF2MW3

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	1.735	12	13	22	rVB	222772	340251	8.28%	1.170%
2	1.861	22	24	65	rVB	2028405	4096032	99.72%	14.084%
3	4.959	292	295	301	rVB	89038	103479	2.52%	0.356%
4	5.381	330	332	335	rBV	980115	1269397	30.90%	4.365%
5	6.422	421	423	426	rBV	557656	780945	19.01%	2.685%
6	6.570	433	436	438	rBV	2281823	2641191	64.30%	9.082%
7	6.799	454	456	458	rBV	791377	658343	16.03%	2.264%
8	6.959	467	470	472	rVB	2554243	2412152	58.72%	8.294%
9	7.370	503	506	508	rBV	1950749	2150234	52.35%	7.394%
10	8.090	566	569	571	rBV	963759	914557	22.27%	3.145%
11	9.176	660	664	666	rBV	2801213	4107582	100.00%	14.124%
12	9.839	720	722	725	rBB	1049892	1034700	25.19%	3.558%
13	10.628	788	791	794	rBV2	1683503	2366504	57.61%	8.137%
14	11.325	849	852	854	rBV	1273430	1129597	27.50%	3.884%
15	12.914	988	991	994	rBV	2810569	3325553	80.96%	11.435%
16	13.954	1080	1082	1085	rBV	1038874	1005348	24.48%	3.457%
17	15.371	1203	1206	1209	rBV	642395	746047	18.16%	2.565%

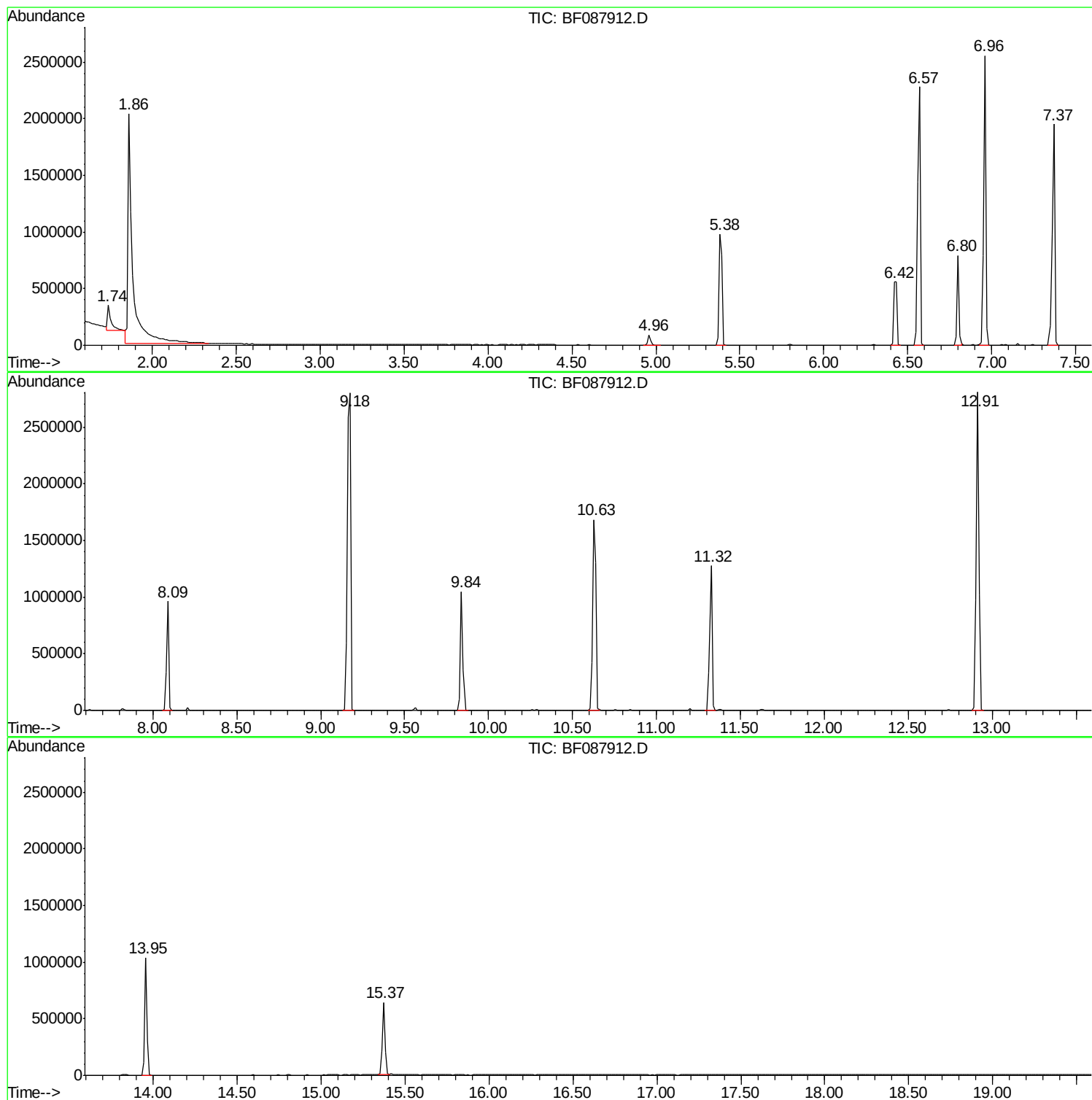
Sum of corrected areas: 29081912

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



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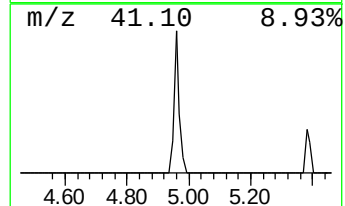
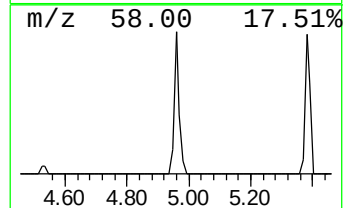
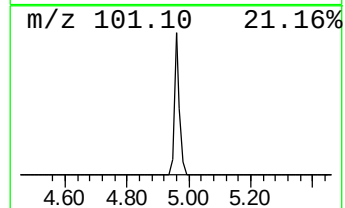
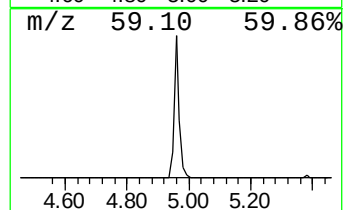
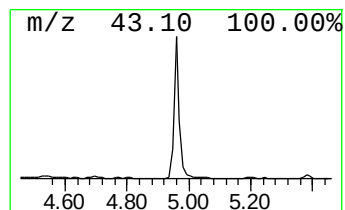
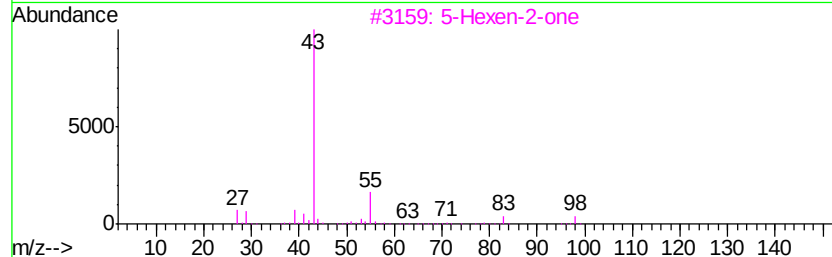
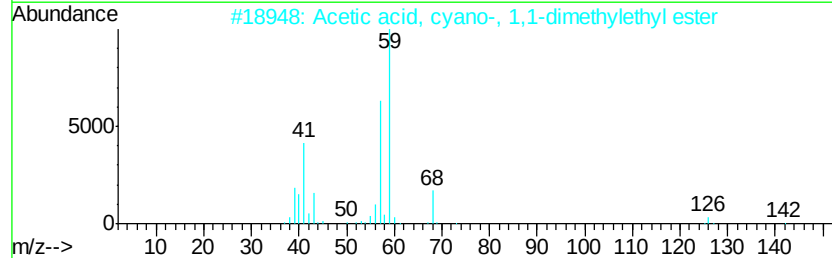
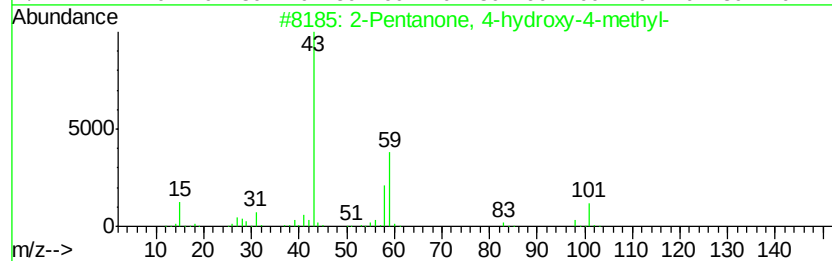
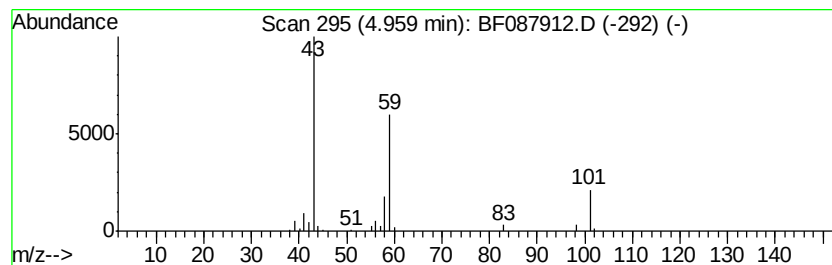
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	3.14 ng	103479	1,4-Dichlorobenzene-d4	6.80

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-dimethyl...	141	C7H11NO2	001116-98-9	23
3	5-Hexen-2-one	98	C6H10O	000109-49-9	9
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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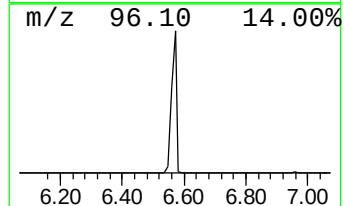
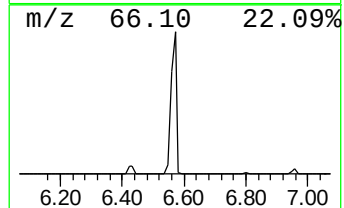
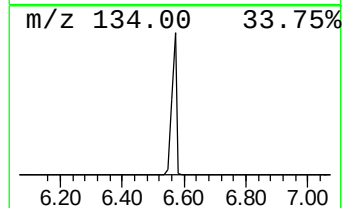
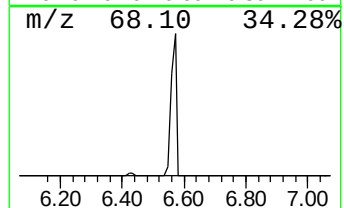
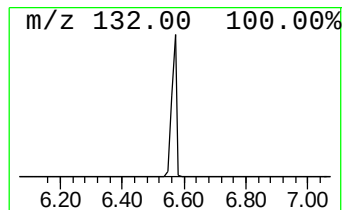
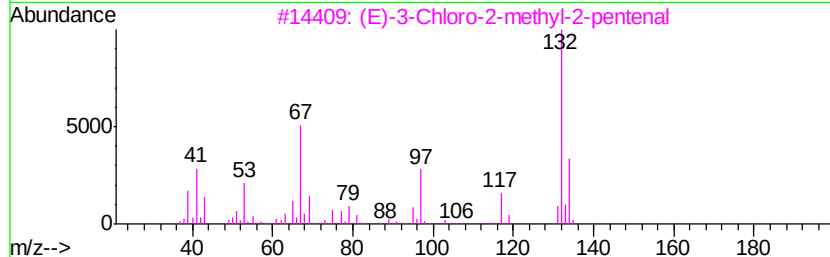
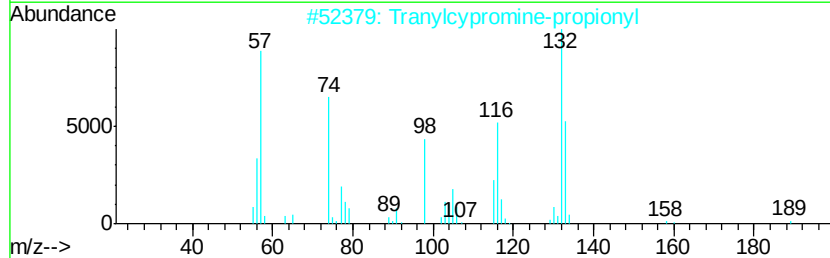
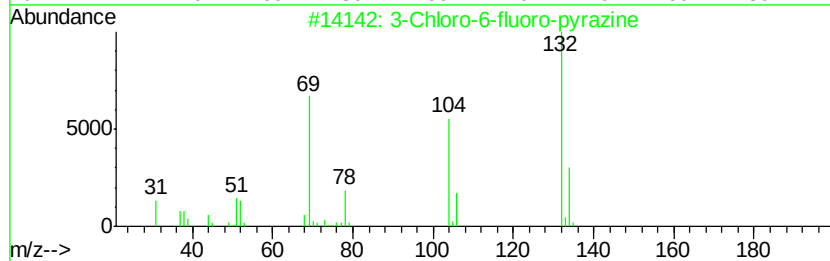
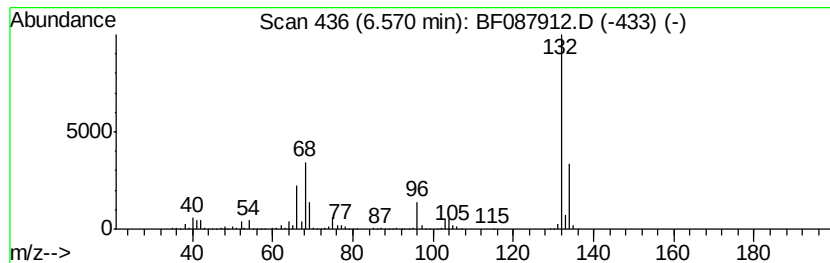
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Peak Number 4 unknown6.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	80.24 ng	2641190	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	38
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	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	3.1	ng	103479	1	6.80	658343	20.0
unknown6.57	6.57	80.2	ng	2641190	1	6.80	658343	20.0