

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040216\
 Data File : BF085978.D
 Acq On : 1 Apr 2016 23:12
 Operator : UM/SJ
 Sample : PB89315BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB89315BL

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-BF040216.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	4.967	250	252	267	rBV	33201	67686	2.09%	0.279%
2	5.367	285	287	290	rBV	1630630	2322236	71.55%	9.589%
3	6.418	376	379	381	rBV	1832191	2455898	75.66%	10.141%
4	6.544	387	390	392	rBV	2226281	2624685	80.86%	10.838%
5	6.773	408	410	421	rVB	669478	607239	18.71%	2.508%
6	6.921	421	423	426	rBV	1614785	2009614	61.91%	8.298%
7	7.344	457	460	462	rBV	1260058	1641846	50.58%	6.780%
8	8.053	520	522	525	rBV	726450	793644	24.45%	3.277%
9	9.139	614	617	619	rBV	2173100	3062168	94.34%	12.645%
10	9.802	673	675	678	rBV	669947	882410	27.19%	3.644%
11	10.602	742	745	747	rBV2	1425531	1873350	57.72%	7.736%
12	11.288	803	805	808	rBV	1032388	949133	29.24%	3.919%
13	12.876	941	944	946	rBV	3368998	3245832	100.00%	13.403%
14	13.928	1034	1036	1038	rBV	814868	904724	27.87%	3.736%
15	15.334	1156	1159	1166	rBV	564492	776393	23.92%	3.206%

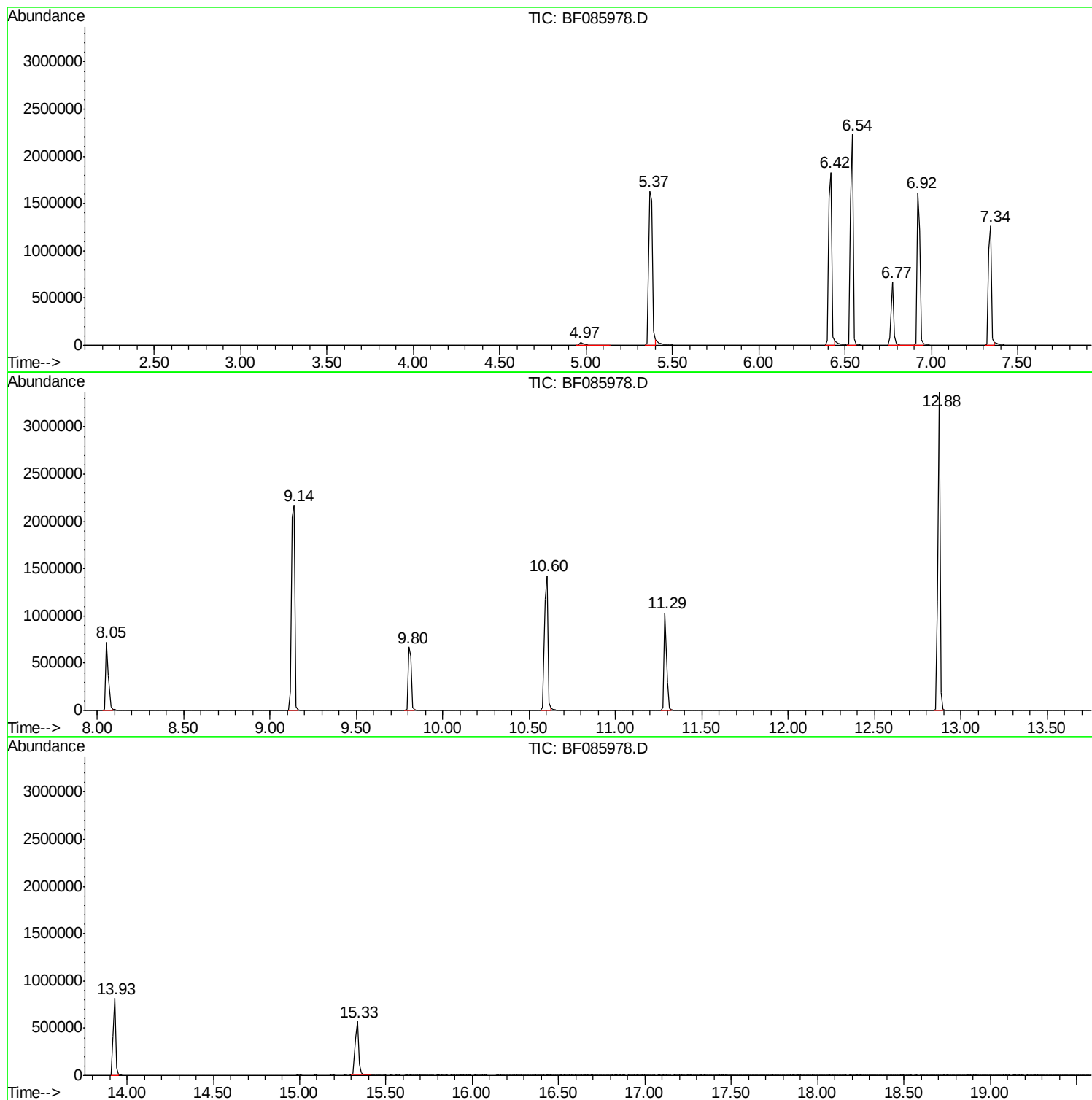
Sum of corrected areas: 24216858

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.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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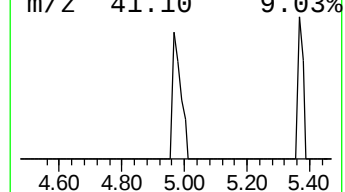
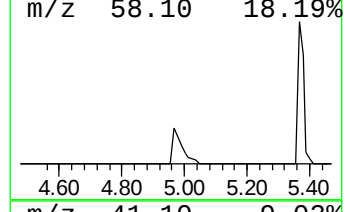
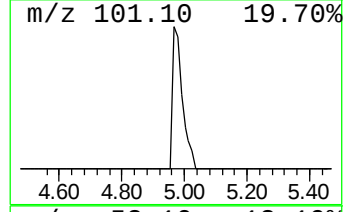
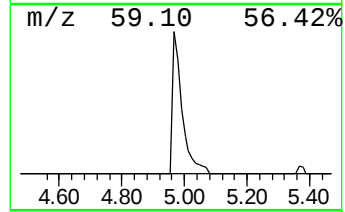
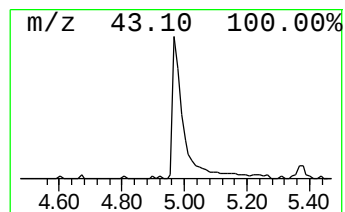
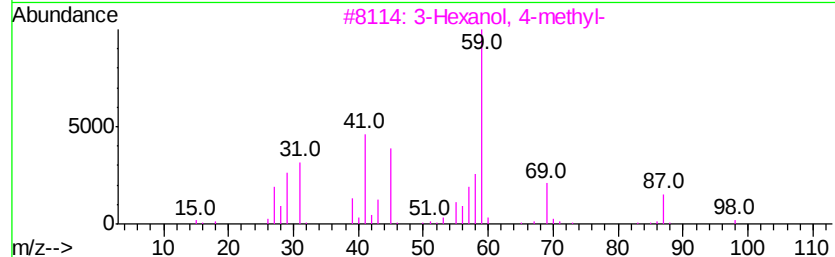
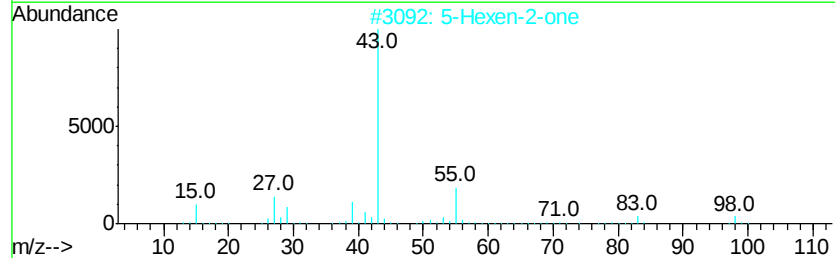
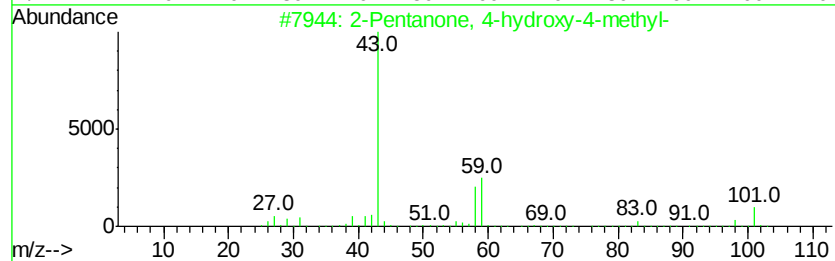
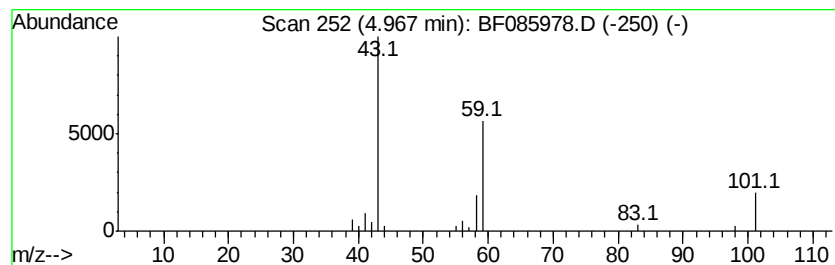
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	2.23 ng	67686	1,4-Dichlorobenzene-d4	6.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			5-Hexen-2-one	98	C6H10O	000109-49-9	9
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7



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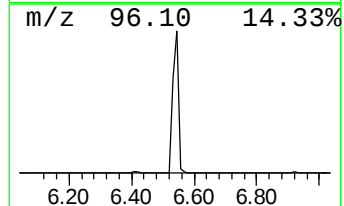
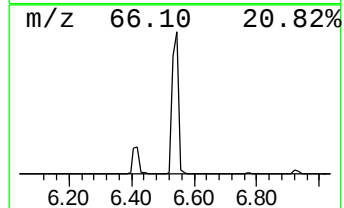
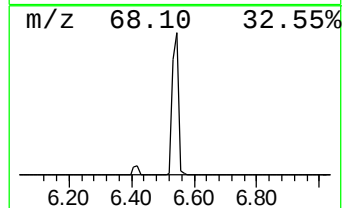
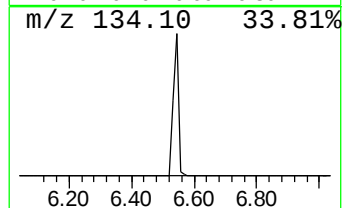
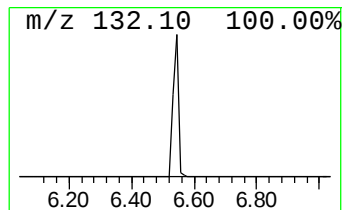
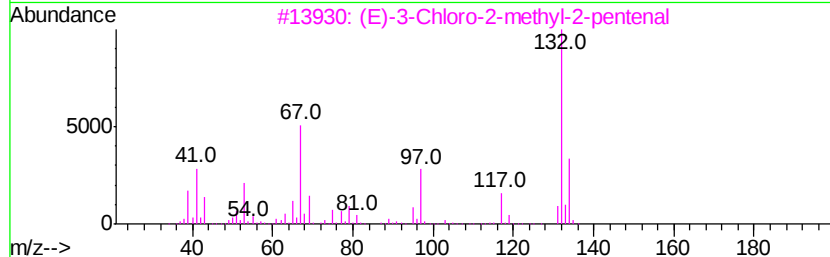
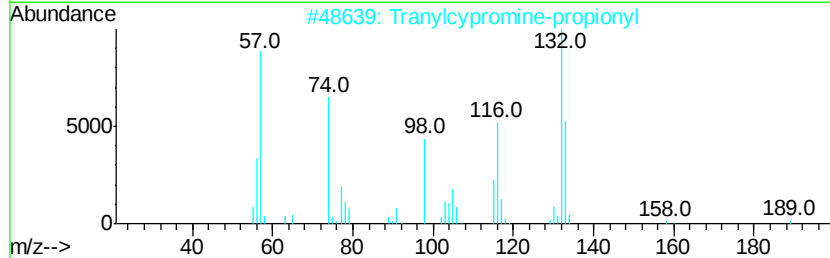
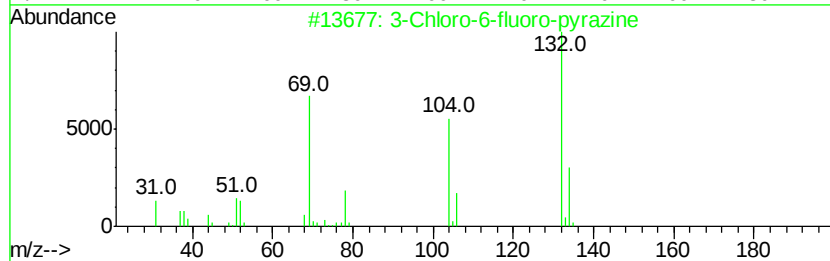
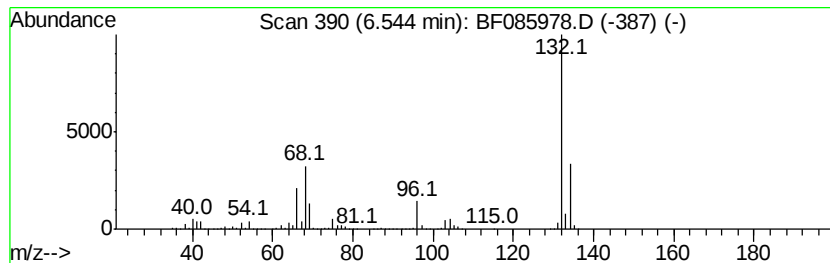
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Peak Number 2 unknown6.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.54	86.45 ng	2624690	1,4-Dichlorobenzene-d4	6.77

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		Xenon	132	Xe	007440-63-3	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.97	2.2	ng	67686	1	6.77	607239	20.0
unknown6.54	6.54	86.5	ng	2624690	1	6.77	607239	20.0