

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050517\
 Data File : BM009918.D
 Acq On : 06 May 2017 00:41
 Operator : SJ/MA
 Sample : I2893-15
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 3074-CC5

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_M\METHODS\8270-BM050517.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	4.904	305	309	316	rBV	153548	232996	1.90%	0.474%
2	5.363	382	387	398	rBV	1113748	1700205	13.87%	3.458%
3	6.957	651	658	670	rBV	708135	1161408	9.47%	2.362%
4	7.316	711	719	728	rBV	2086930	3371968	27.50%	6.859%
5	7.775	789	797	805	rBV	451800	750506	6.12%	1.527%
6	8.098	844	852	860	rBV	1864794	3041678	24.81%	6.187%
7	8.951	987	997	1007	rBV	1849637	3048761	24.86%	6.202%
8	10.575	1265	1273	1287	rVB	575064	981490	8.00%	1.997%
9	13.045	1683	1693	1702	rBV	4506159	6301121	51.39%	12.817%
10	14.433	1923	1929	1937	rBV2	852965	1321829	10.78%	2.689%
11	15.933	2177	2184	2199	rBV	4063993	5652736	46.10%	11.499%
12	17.186	2391	2397	2403	rBV2	1198650	1711332	13.96%	3.481%
13	19.815	2837	2844	2849	rBV	10591464	12261741	100.00%	24.942%
14	21.056	3052	3055	3061	rBV3	173661	201060	1.64%	0.409%
15	21.262	3086	3090	3095	rBV	3178999	3226679	26.32%	6.564%
16	21.374	3104	3109	3114	rBV	1830794	2225201	18.15%	4.526%
17	23.680	3495	3501	3508	rVB2	1019054	1969617	16.06%	4.007%

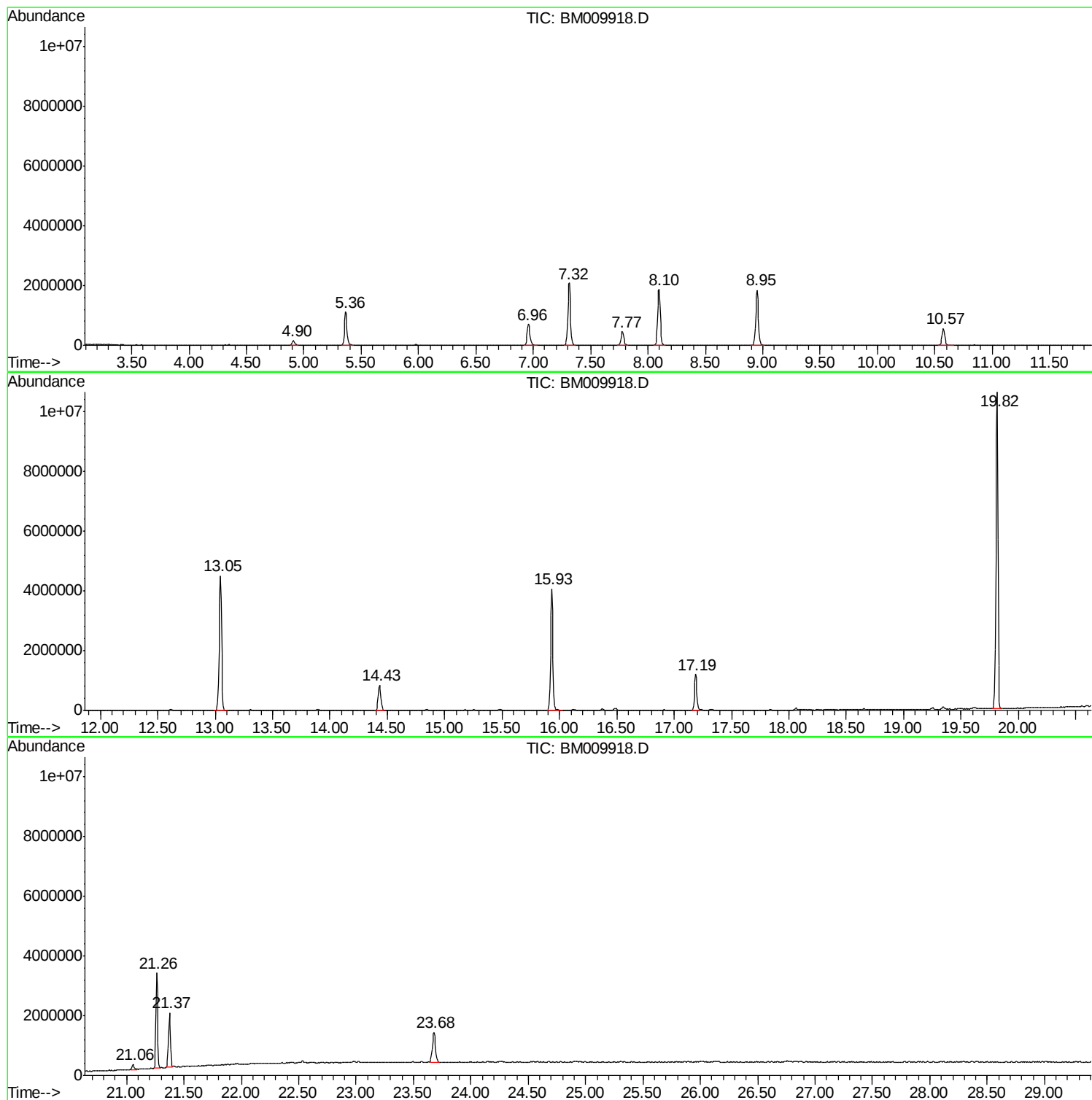
Sum of corrected areas: 49160328

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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM050517.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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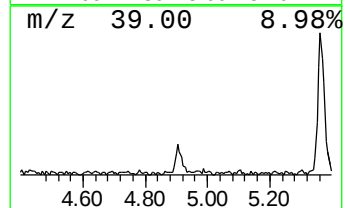
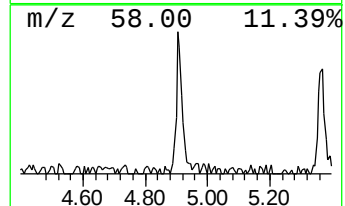
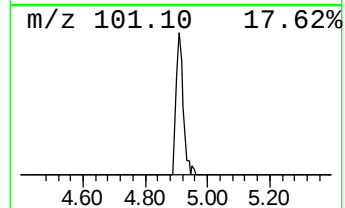
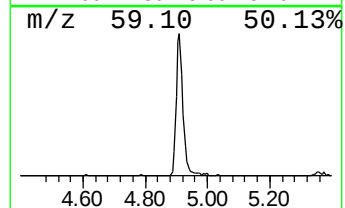
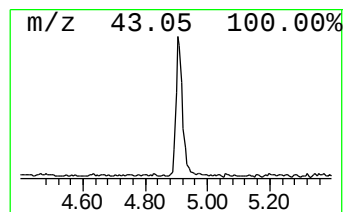
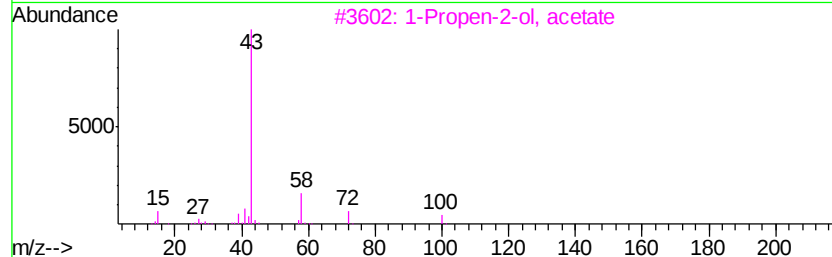
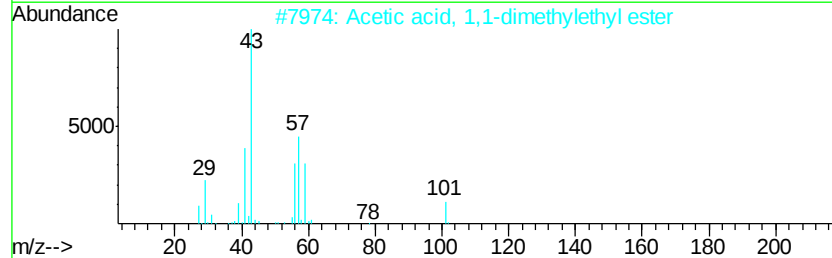
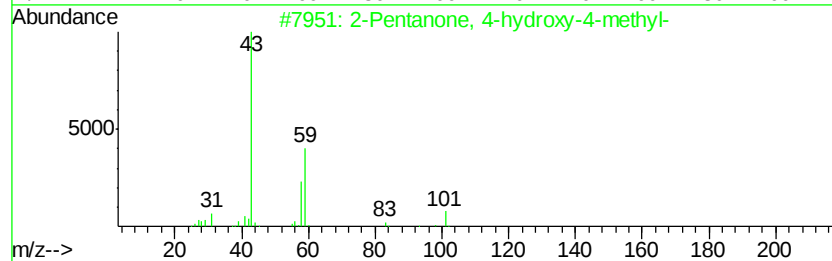
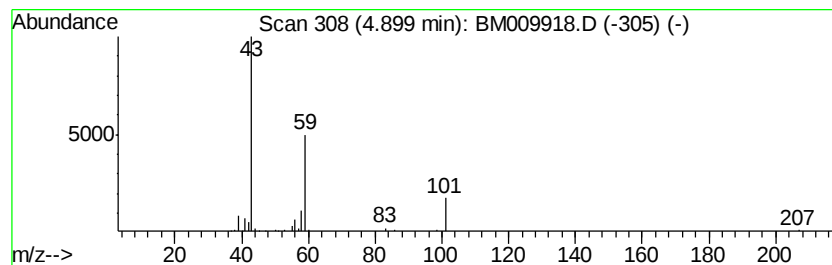
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	6.21 ng	232996	1,4-Dichlorobenzene-d4	7.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Acetone	58	C3H6O	000067-64-1	9



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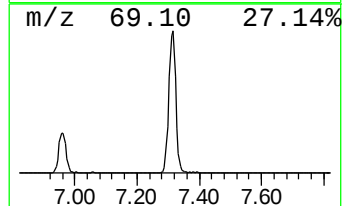
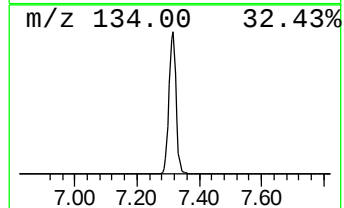
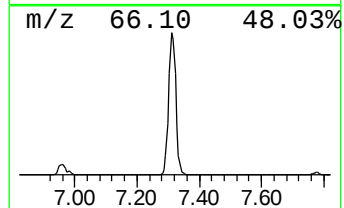
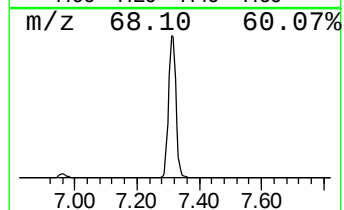
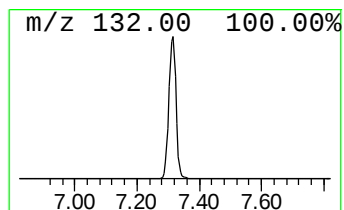
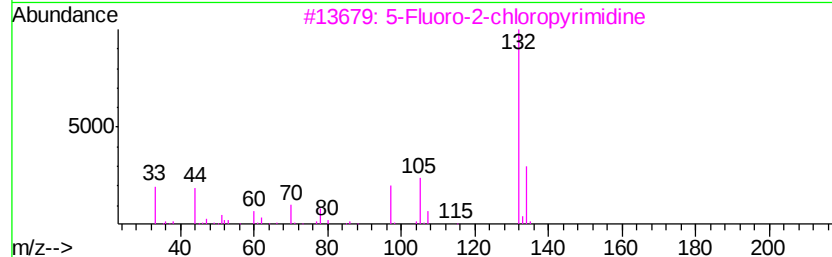
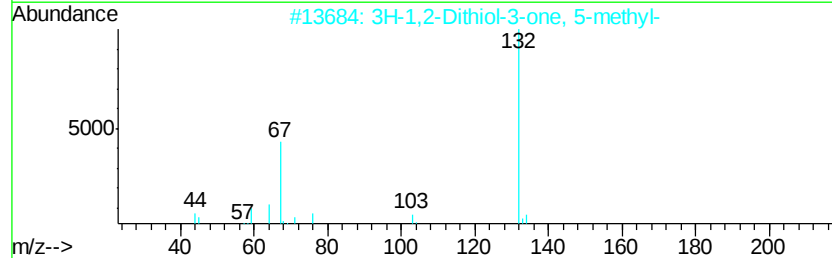
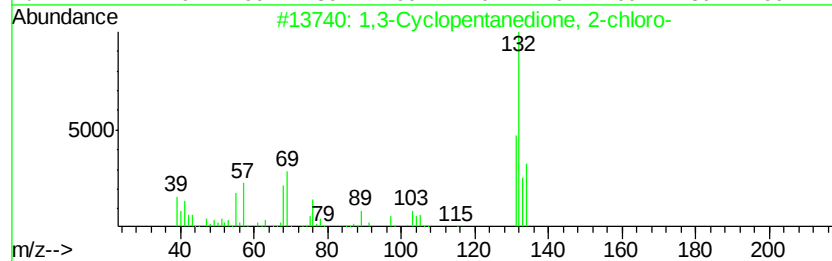
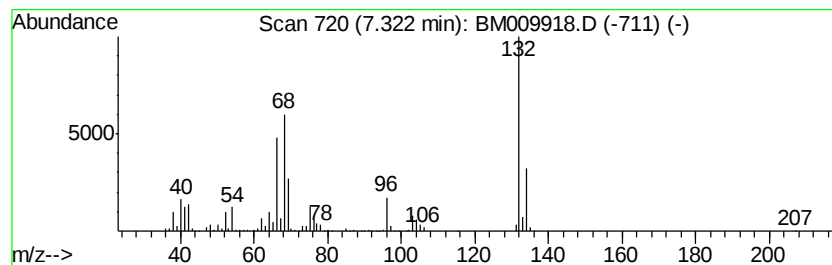
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Peak Number 2 unknown7.32 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.32	89.86 ng	3371970	1,4-Dichlorobenzene-d4	7.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	16
2		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	14
4		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	14
5		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.90	6.2	ng	232996	1	7.78	750506	20.0
unknown7.32	7.32	89.9	ng	3371970	1	7.78	750506	20.0