

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM101918\
 Data File : BM017230.D
 Acq On : 19 Oct 2018 15:14
 Operator : MJ/SJ
 Sample : PB113600BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB113600BL

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:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM101718.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	4.840	310	315	325	rBV	171526	233570	2.17%	0.322%
2	5.281	384	390	404	rBV	4257578	5936234	55.28%	8.193%
3	6.846	648	656	670	rBV	4027996	6329846	58.94%	8.736%
4	7.204	710	717	728	rBV	4232618	6457607	60.13%	8.912%
5	7.669	789	796	803	rBV	1181135	1879875	17.51%	2.594%
6	7.981	842	849	861	rBV	3104183	4985411	46.42%	6.880%
7	8.822	983	992	1002	rBV	2361171	3866162	36.00%	5.336%
8	10.439	1259	1267	1282	rBV	1489956	2530644	23.57%	3.493%
9	12.922	1682	1689	1704	rBV	6386683	9413352	87.66%	12.991%
10	14.310	1918	1925	1935	rBV2	2103250	3389453	31.56%	4.678%
11	15.804	2172	2179	2196	rBV	4683178	6752532	62.88%	9.319%
12	17.057	2386	2392	2406	rBV2	2733396	3901129	36.33%	5.384%
13	19.698	2835	2841	2849	rBV	8289999	10738943	100.00%	14.821%
14	21.250	3100	3105	3112	rBV	2752743	3382319	31.50%	4.668%
15	23.462	3475	3481	3491	rVB	1499815	2661946	24.79%	3.674%

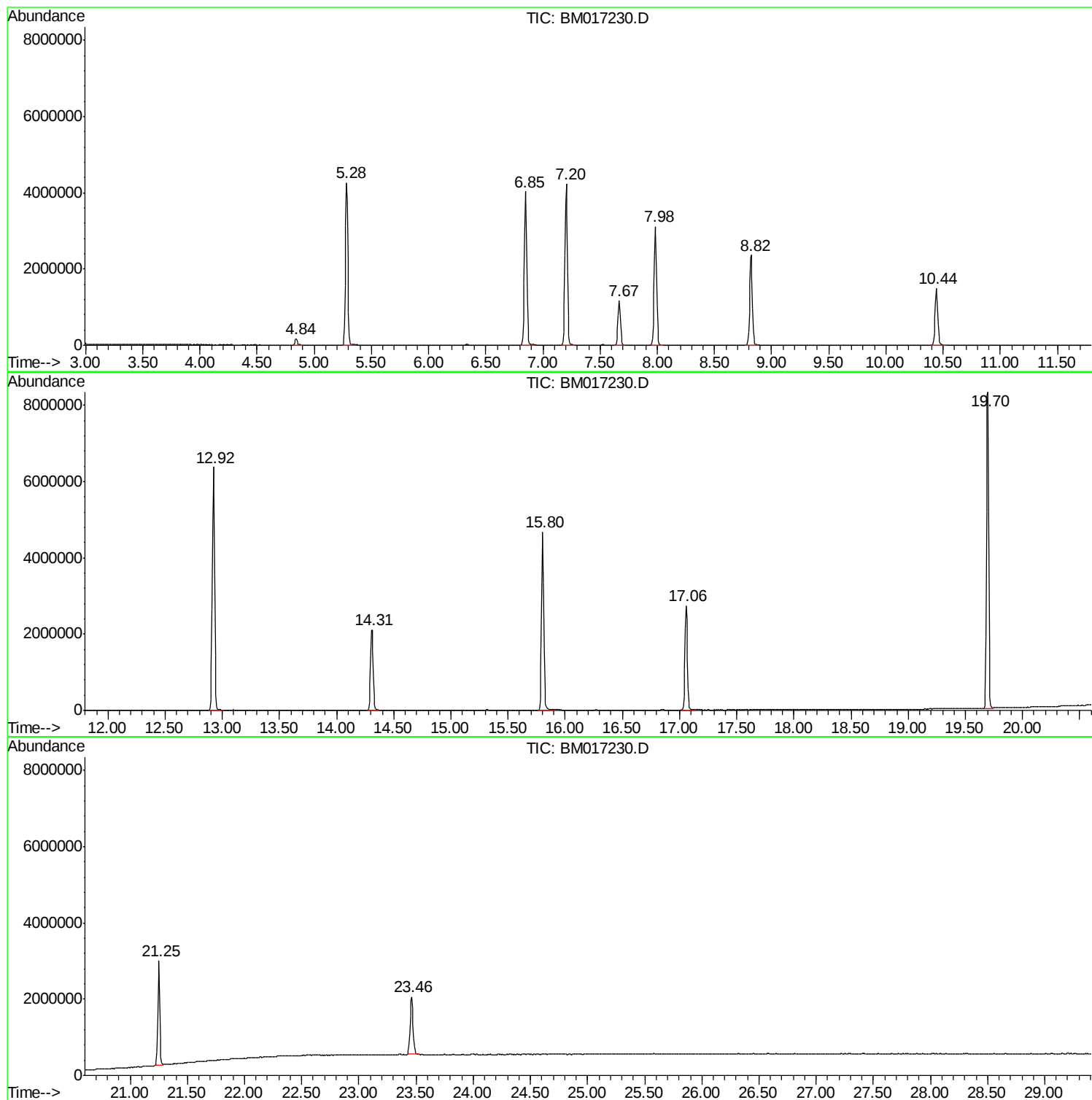
Sum of corrected areas: 72459023

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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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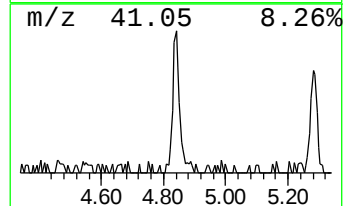
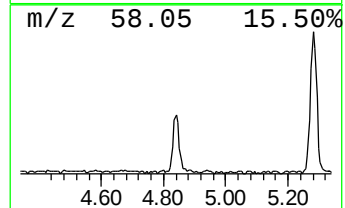
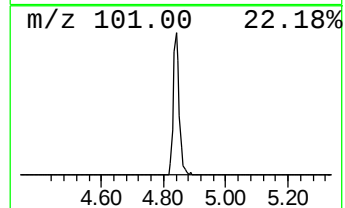
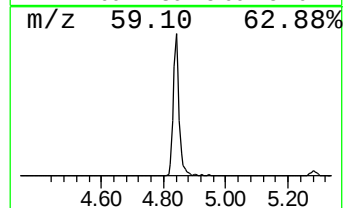
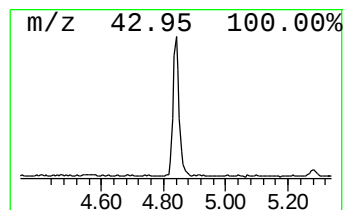
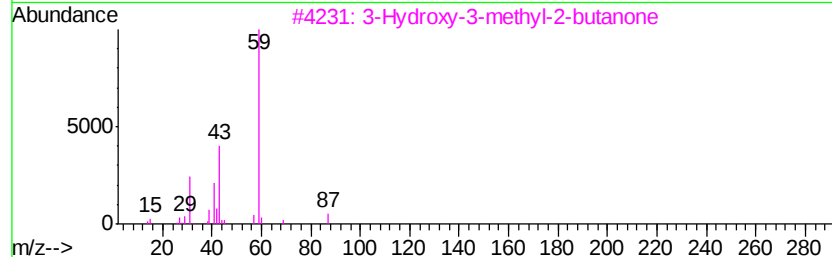
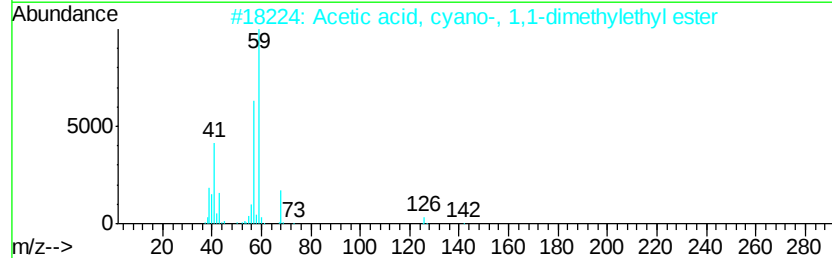
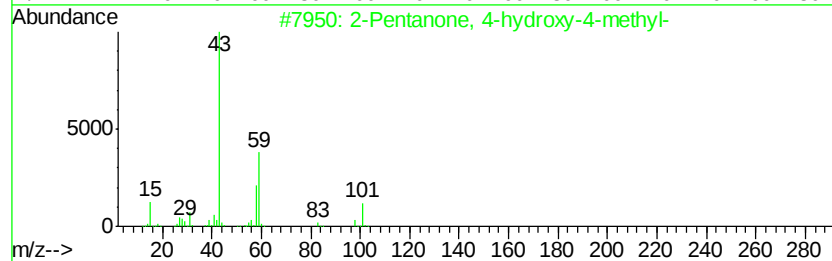
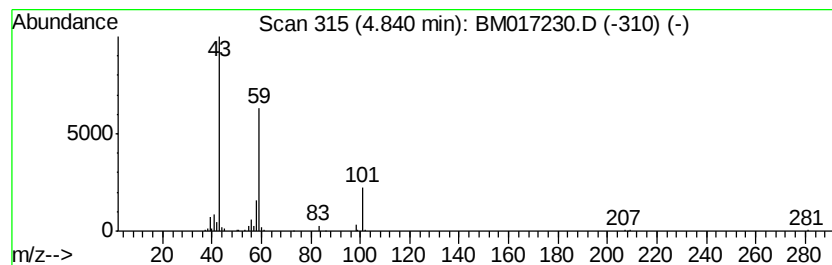
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.84	2.48 ng	233570	1,4-Dichlorobenzene-d4	7.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
3			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17
4			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	17
5			t-Butyl ethyl ether	102	C6H14O	1000118-18-4	17



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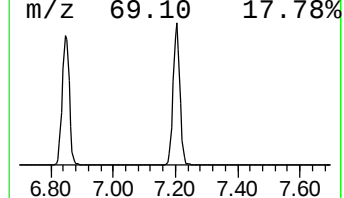
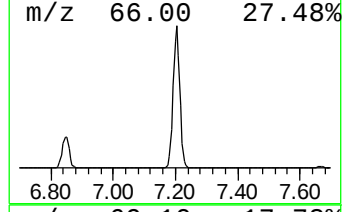
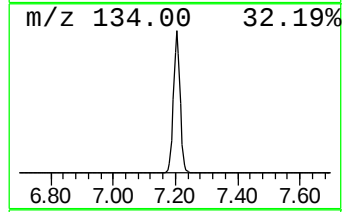
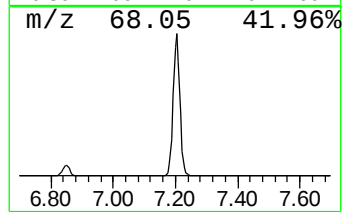
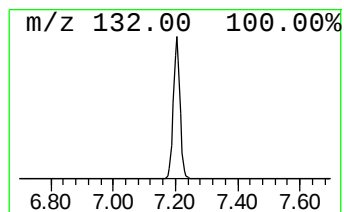
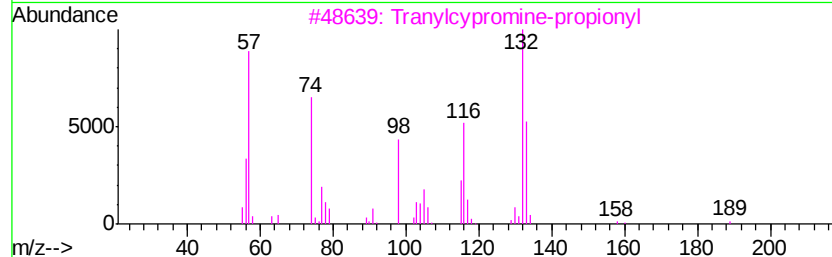
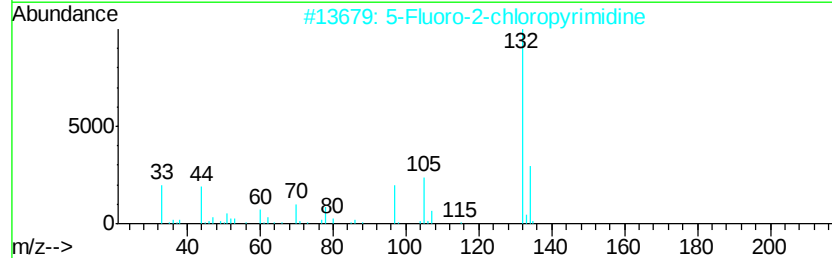
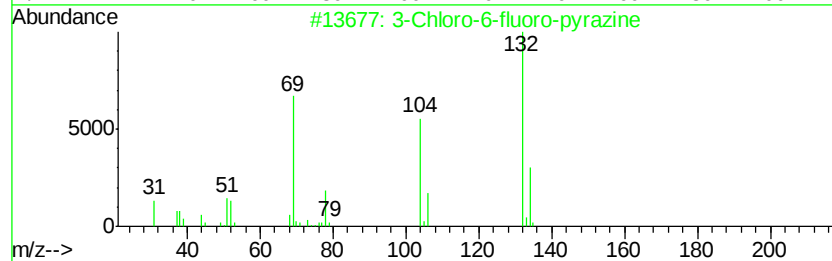
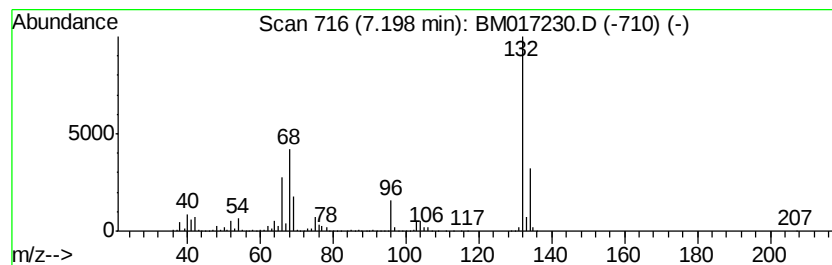
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Peak Number 2 unknown7.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	68.70 ng	6457610	1,4-Dichlorobenzene-d4	7.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.84	2.5	ng	233570	1	7.67	1879880	20.0
unknown7.20	7.20	68.7	ng	6457610	1	7.67	1879880	20.0