

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM101918\
 Data File : BM017232.D
 Acq On : 19 Oct 2018 16:27
 Operator : MJ/SJ
 Sample : PB113946BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB113946BL

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	rVB	175287	236756	1.67%	0.271%
2	5.281	383	390	399	rBV	5280218	7447927	52.48%	8.520%
3	6.851	646	657	666	rBV	5058040	7916483	55.78%	9.056%
4	7.204	710	717	730	rBV	5300656	8078911	56.92%	9.242%
5	7.669	787	796	806	rVB	1175160	1891373	13.33%	2.164%
6	7.981	842	849	860	rBV	3865365	6213219	43.78%	7.108%
7	8.822	984	992	1008	rBV	2947157	4862408	34.26%	5.562%
8	10.439	1260	1267	1286	rBV	1492195	2523628	17.78%	2.887%
9	12.922	1682	1689	1698	rBV	7901137	11734912	82.68%	13.424%
10	14.304	1918	1924	1937	rBV2	2209999	3408025	24.01%	3.899%
11	15.804	2172	2179	2195	rBV	5963992	8606164	60.64%	9.845%
12	17.057	2385	2392	2405	rBV	2864052	3951729	27.84%	4.521%
13	19.698	2835	2841	2849	rBV	11431280	14192770	100.00%	16.236%
14	21.250	3100	3105	3112	rBV	2966375	3617191	25.49%	4.138%
15	23.462	3475	3481	3488	rVB	1482455	2734174	19.26%	3.128%

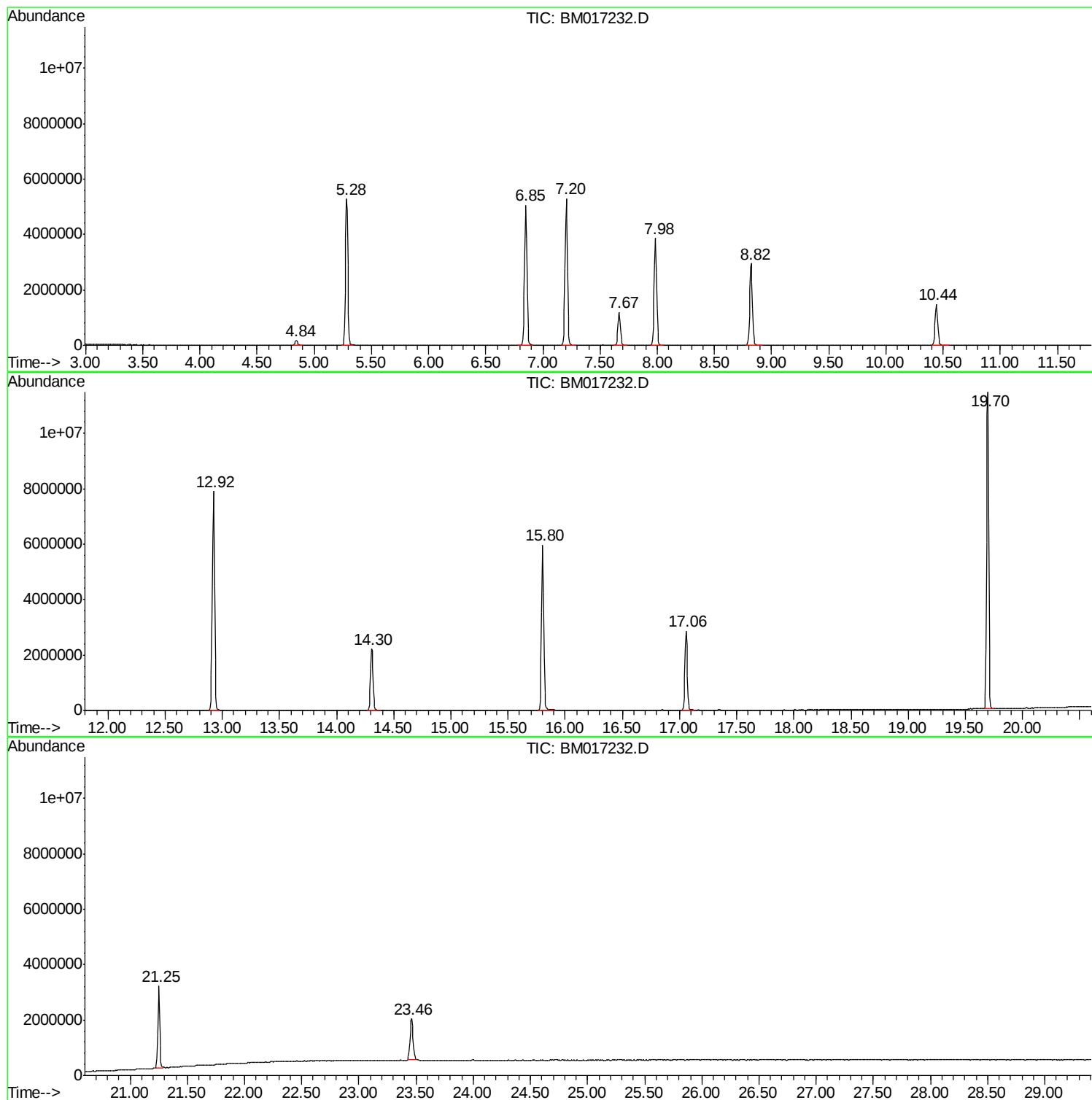
Sum of corrected areas: 87415670

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



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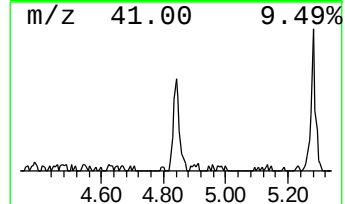
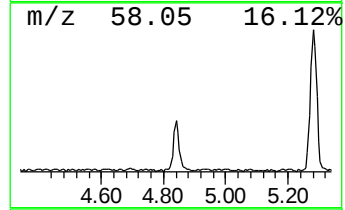
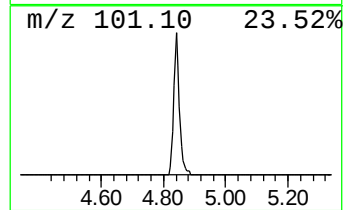
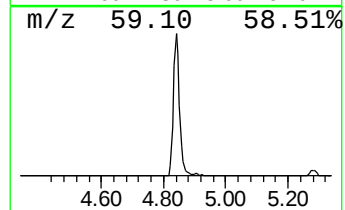
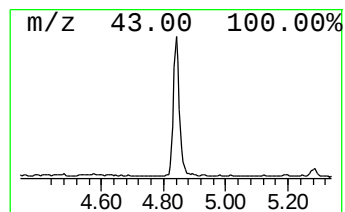
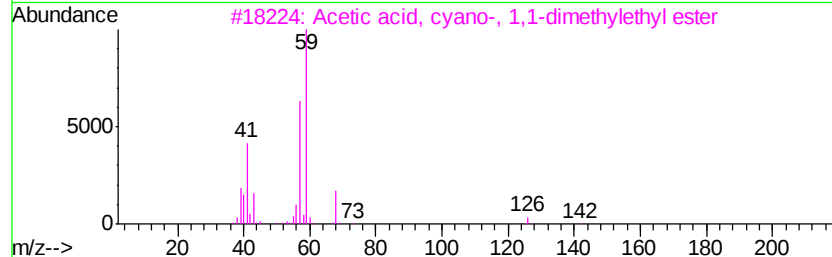
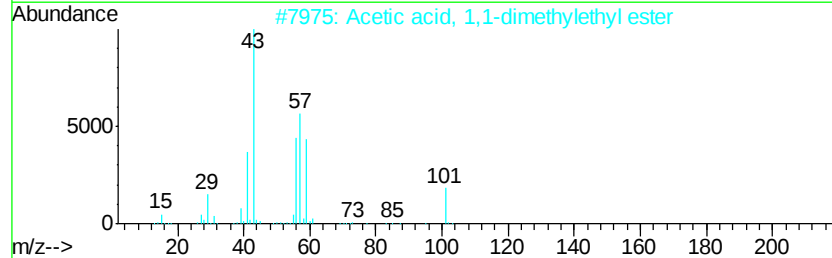
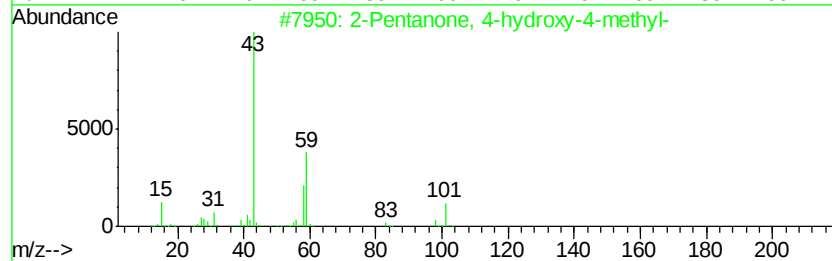
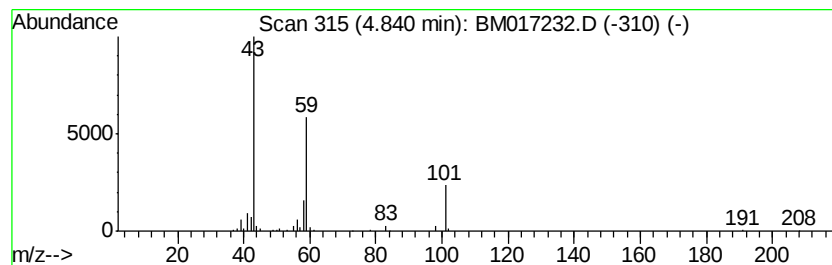
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.84	2.50 ng	236756	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	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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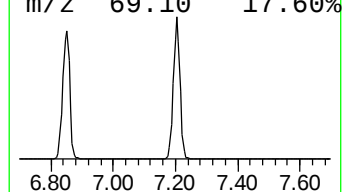
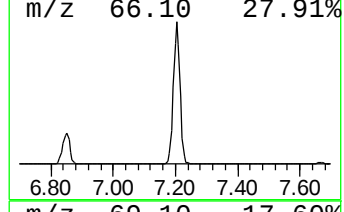
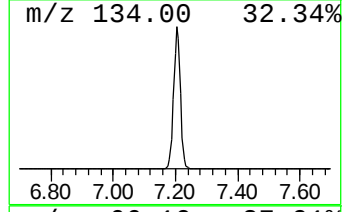
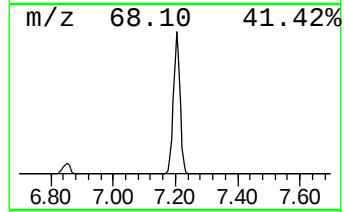
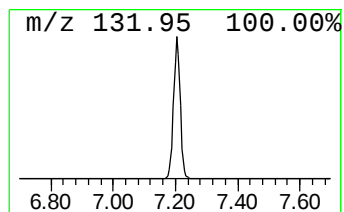
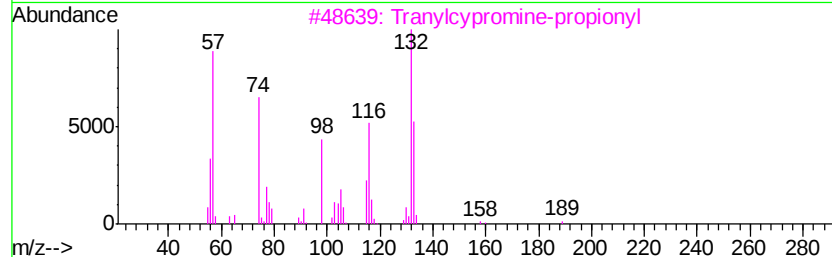
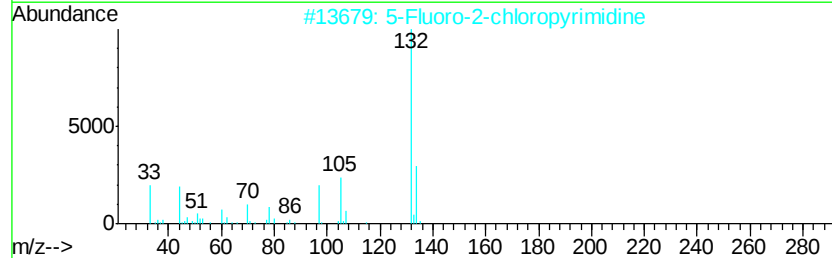
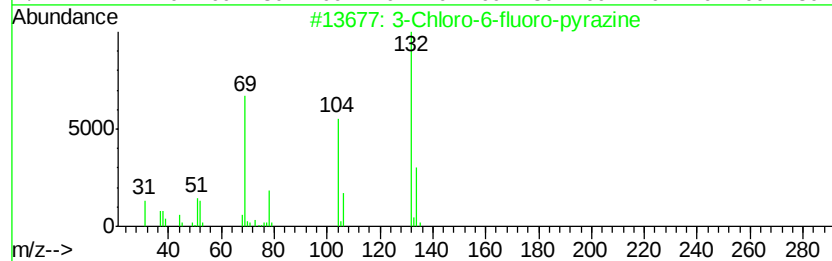
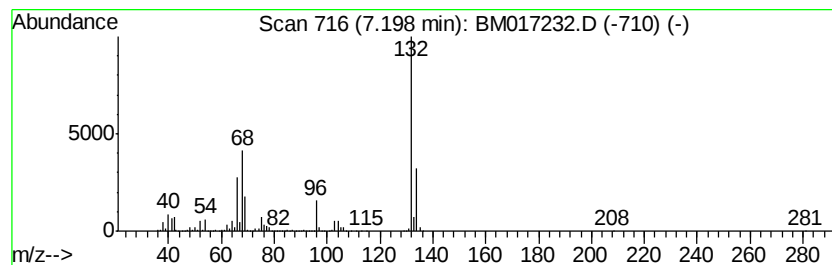
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Peak Number 2 unknown7.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	85.43 ng	8078910	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	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	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	236756	1	7.67	1891370	20.0
unknown7.20	7.20	85.4	ng	8078910	1	7.67	1891370	20.0