

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM101918\
 Data File : BM017237.D
 Acq On : 19 Oct 2018 19:28
 Operator : MJ/SJ
 Sample : PB113945BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB113945BL

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	172418	235168	1.99%	0.309%
2	5.281	383	390	401	rBV	4379885	6092122	51.52%	8.017%
3	6.845	649	656	667	rBV	4140775	6467760	54.70%	8.511%
4	7.204	709	717	729	rBV	4324159	6603639	55.85%	8.690%
5	7.669	789	796	803	rBV	1193720	1909379	16.15%	2.513%
6	7.981	841	849	860	rBV	3245478	5123011	43.33%	6.741%
7	8.822	983	992	1006	rBV	2413870	3986206	33.71%	5.245%
8	10.439	1259	1267	1279	rBV	1521243	2576268	21.79%	3.390%
9	12.921	1682	1689	1706	rBV	6753281	9741006	82.38%	12.818%
10	14.310	1918	1925	1940	rBV2	2239963	3517677	29.75%	4.629%
11	15.798	2172	2178	2199	rBV	4854493	7172232	60.66%	9.438%
12	17.056	2385	2392	2405	rBV	2910476	4121460	34.86%	5.423%
13	19.692	2835	2840	2850	rBV	9274171	11824400	100.00%	15.560%
14	21.250	3100	3105	3111	rBV	2925001	3665277	31.00%	4.823%
15	23.462	3475	3481	3495	rVB	1562406	2957886	25.02%	3.892%

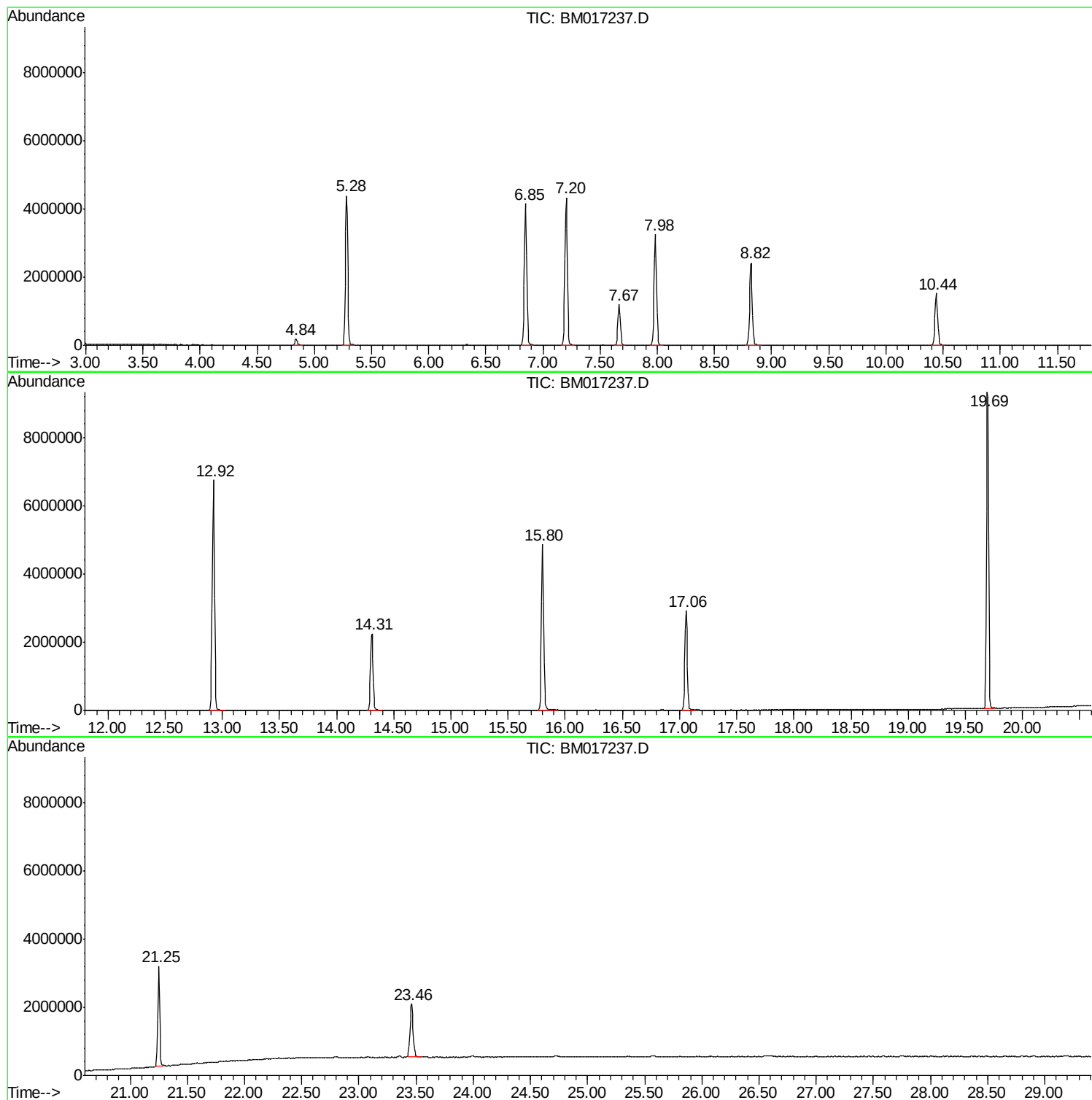
Sum of corrected areas: 75993491

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



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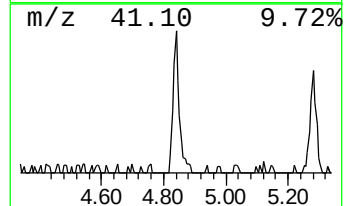
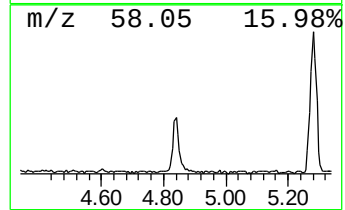
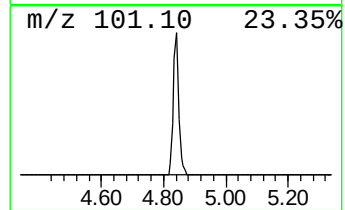
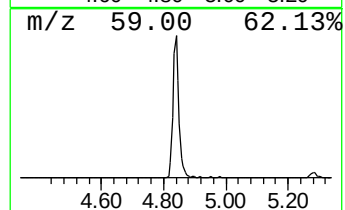
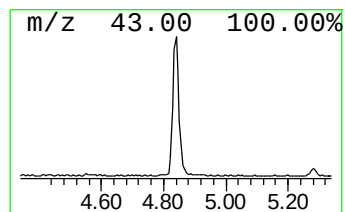
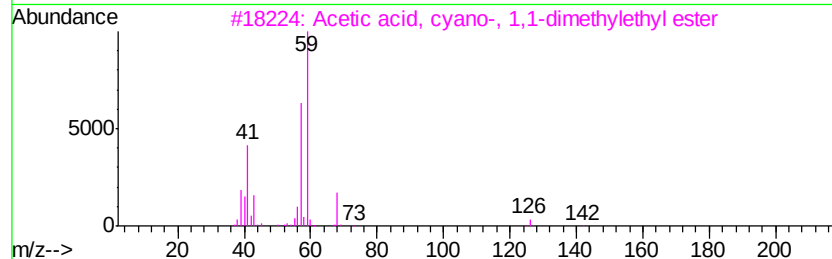
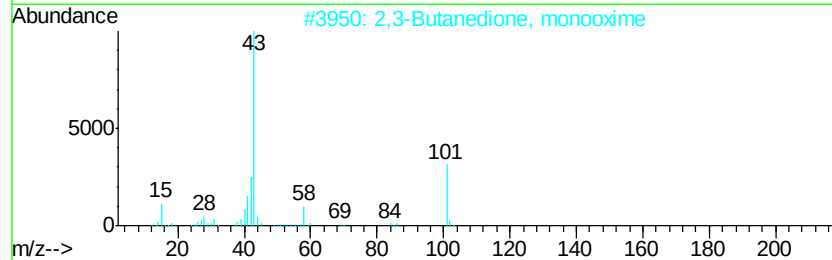
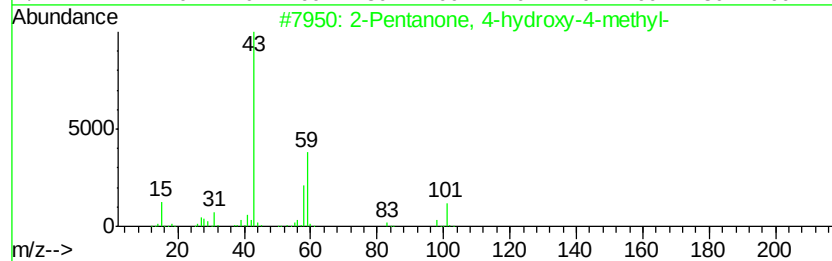
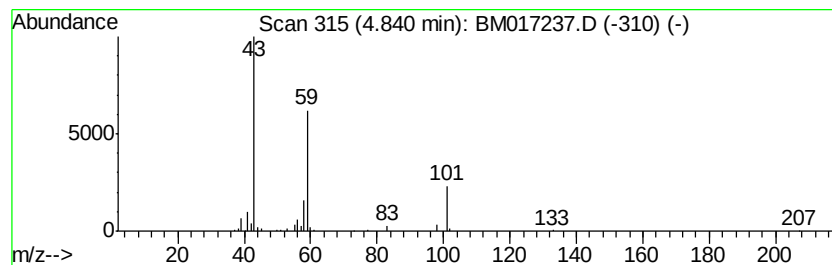
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.84	2.46 ng	235168	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	72
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	9



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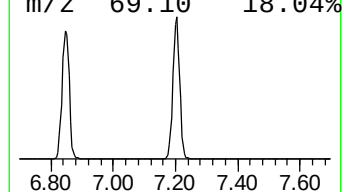
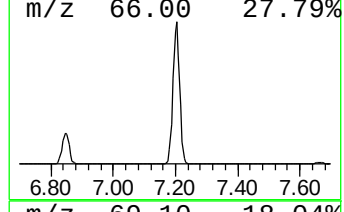
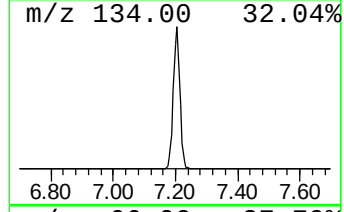
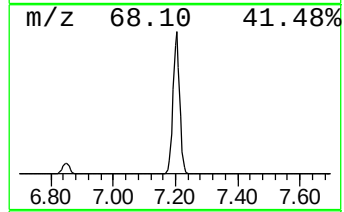
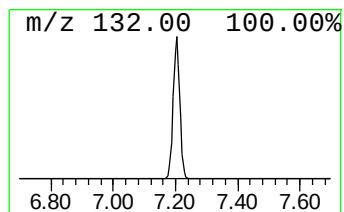
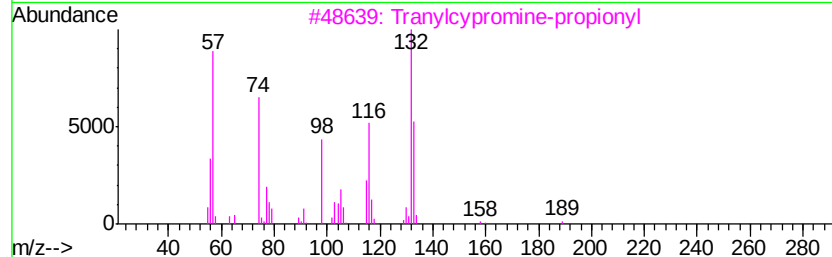
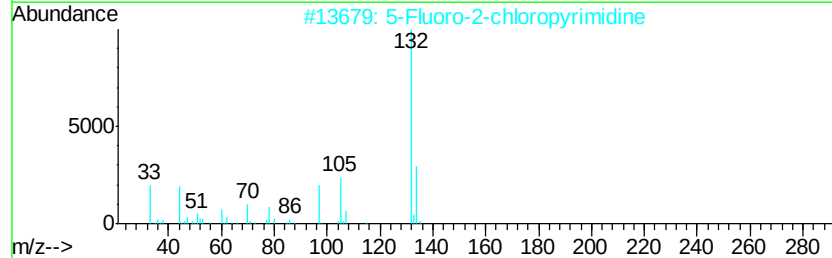
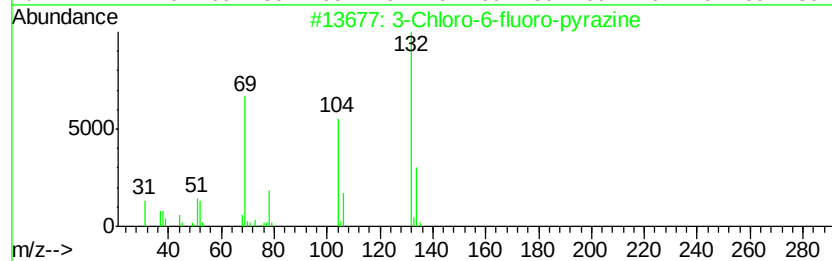
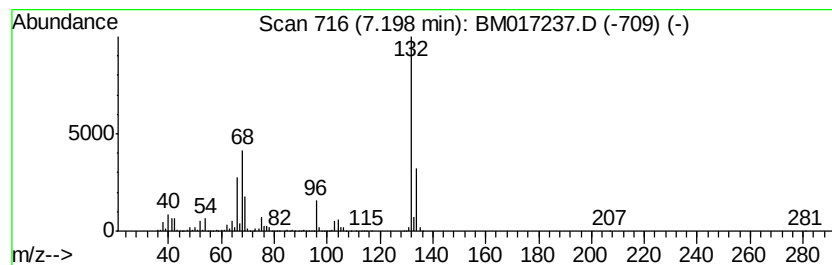
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Peak Number 2 unknown7.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	69.17 ng	6603640	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	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	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	235168	1	7.67	1909380	20.0
unknown7.20	7.20	69.2	ng	6603640	1	7.67	1909380	20.0