

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111018\
 Data File : BM017596.D
 Acq On : 09 Nov 2018 18:00
 Operator : MJ/SJ
 Sample : PB114669BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB114669BL

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	4.799	303	308	314	rBV	163625	231540	1.69%	0.301%
2	5.246	377	384	398	rBV	4515368	6442755	47.04%	8.382%
3	6.810	641	650	666	rBV	4325646	6868500	50.15%	8.936%
4	7.157	702	709	719	rBV	4553892	6961223	50.83%	9.056%
5	7.616	781	787	795	rBV	923726	1492942	10.90%	1.942%
6	7.934	832	841	854	rBV	3265569	5240435	38.26%	6.818%
7	8.775	975	984	1011	rBV	2429106	4220759	30.82%	5.491%
8	10.392	1251	1259	1279	rBV	1022772	1991649	14.54%	2.591%
9	12.880	1674	1682	1705	rBV	5997512	9600253	70.10%	12.490%
10	14.263	1910	1917	1932	rBV2	1563250	2605653	19.03%	3.390%
11	15.763	2165	2172	2198	rBV	5086090	7629618	55.71%	9.926%
12	17.021	2379	2386	2408	rBV	1826568	3166712	23.12%	4.120%
13	19.662	2829	2835	2847	rBV	10434967	13695257	100.00%	17.817%
14	21.215	3094	3099	3113	rBV2	2484048	3687308	26.92%	4.797%
15	23.409	3466	3472	3486	rVB	1498187	3030804	22.13%	3.943%

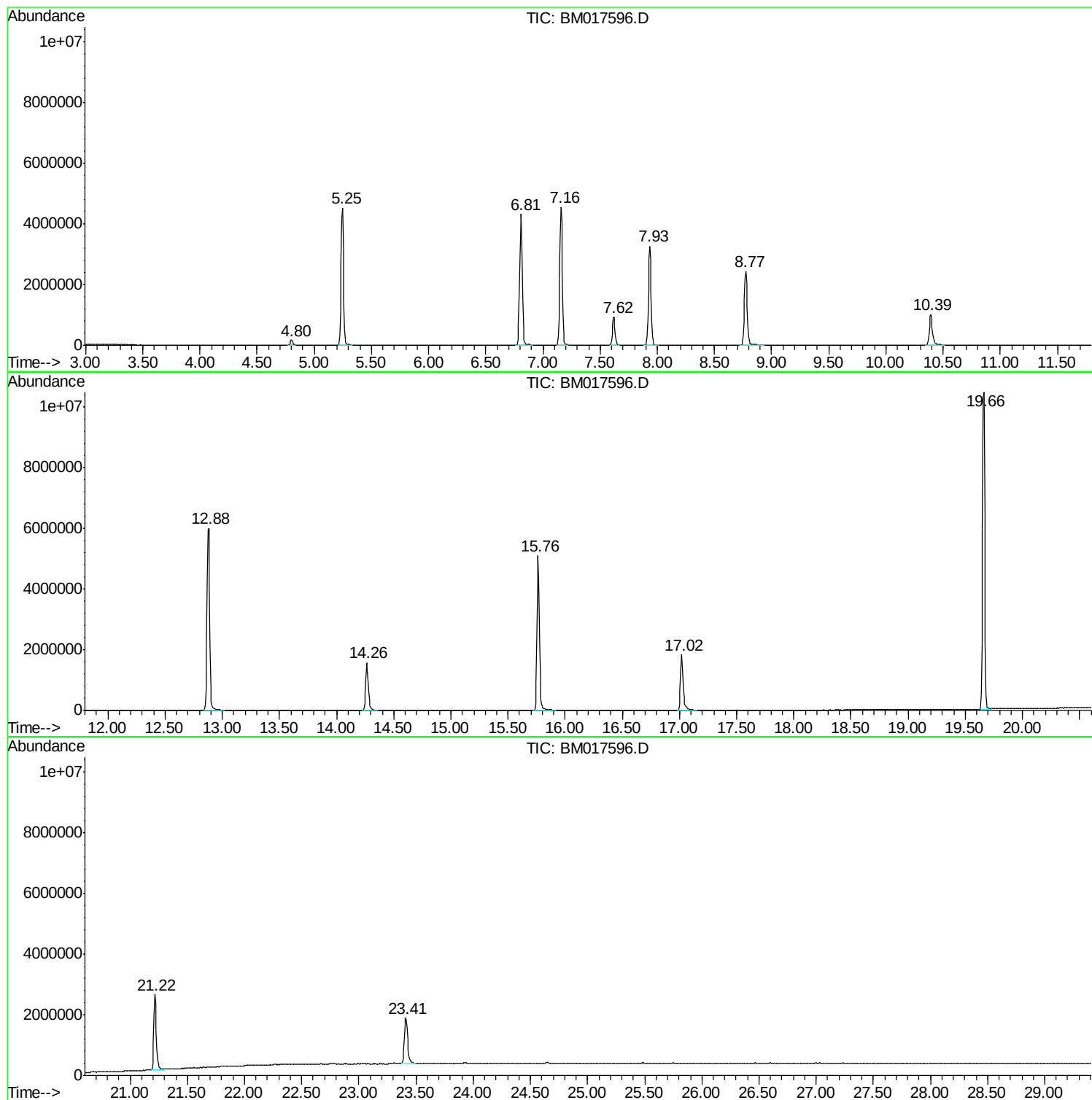
Sum of corrected areas: 76865408

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



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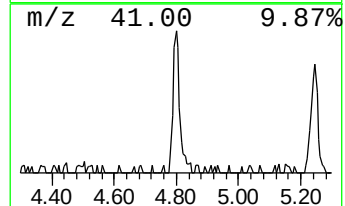
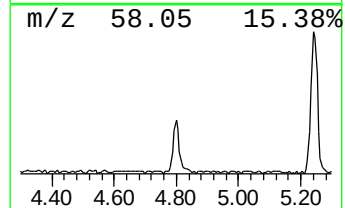
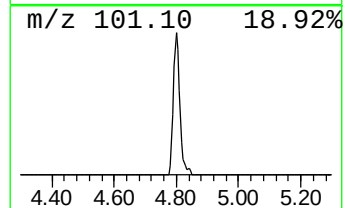
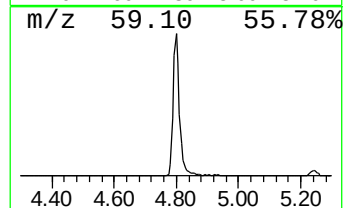
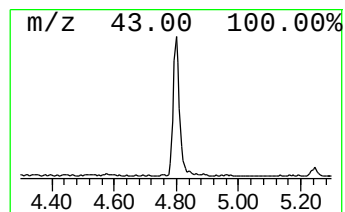
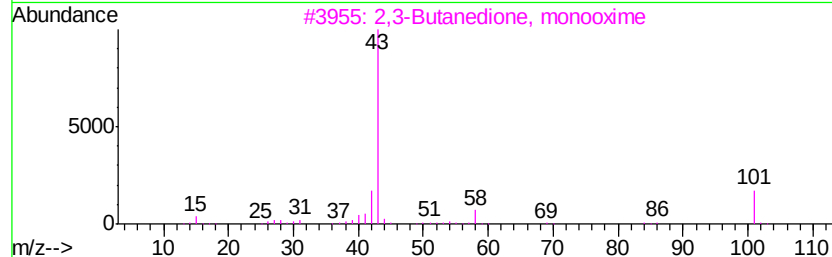
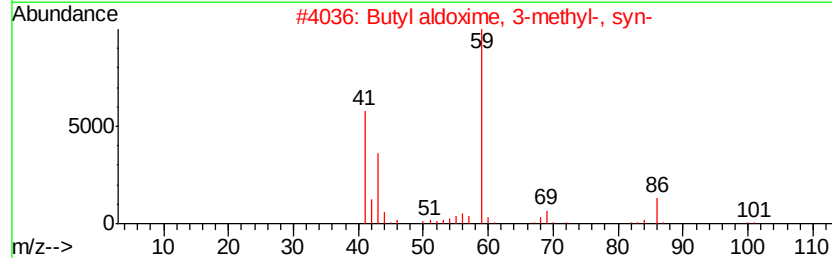
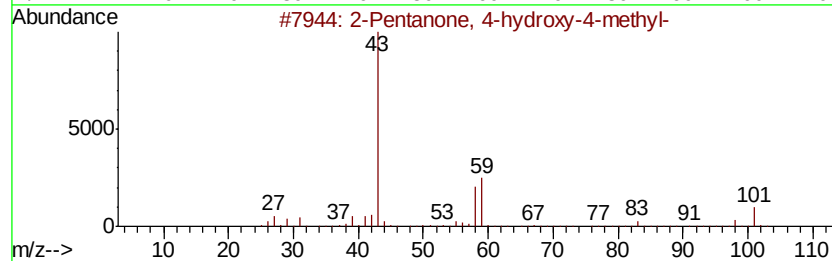
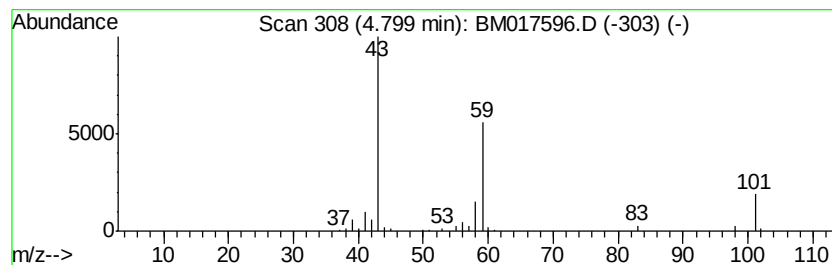
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	3.10 ng	231540	1,4-Dichlorobenzene-d4	7.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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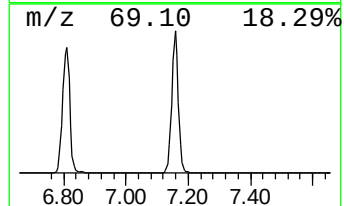
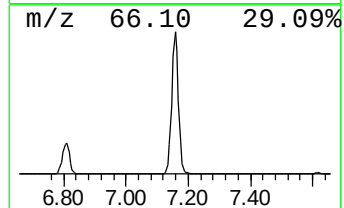
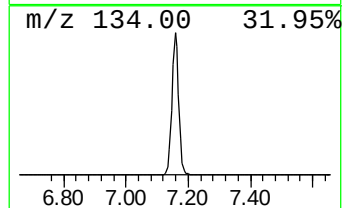
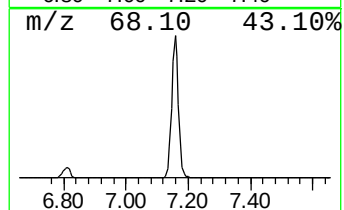
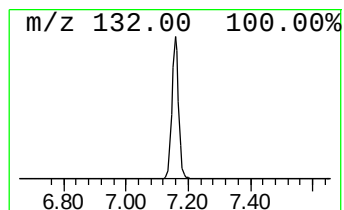
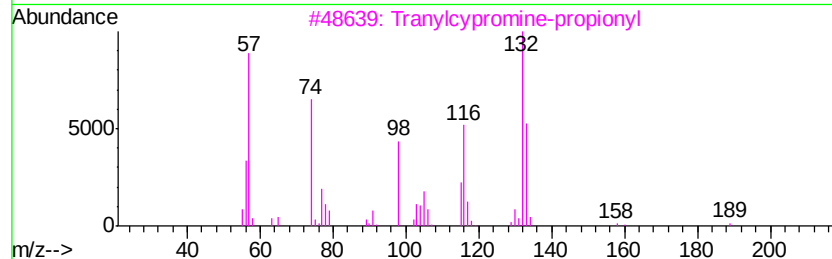
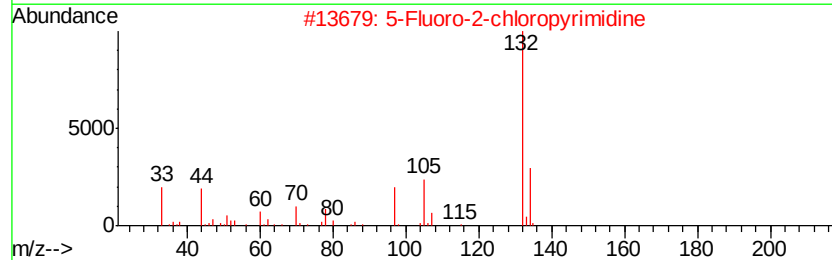
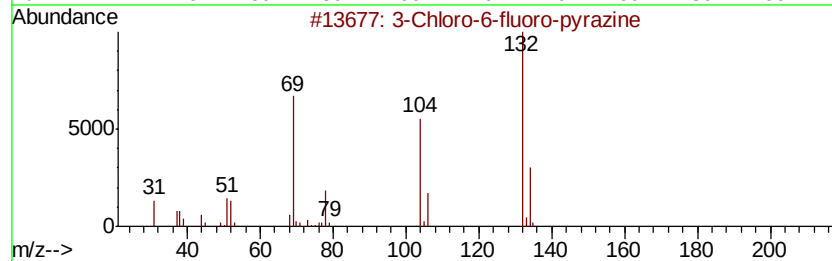
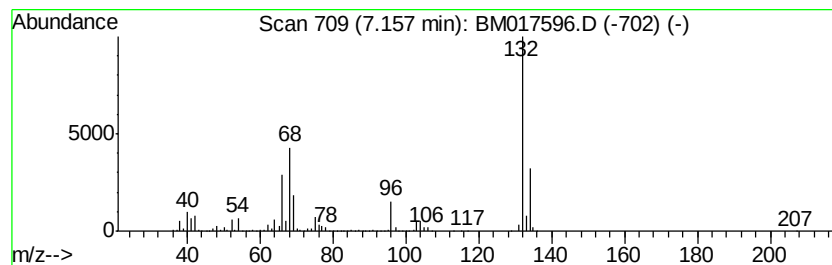
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Peak Number 2 unknown7.16 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	93.26 ng	6961220	1,4-Dichlorobenzene-d4	7.62

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		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.80	3.1	ng	231540	1	7.62	1492940	20.0
unknown7.16	7.16	93.3	ng	6961220	1	7.62	1492940	20.0