

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM101818\
 Data File : BM017208.D
 Acq On : 19 Oct 2018 00:21
 Operator : MJ/SJ
 Sample : PB113966BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB113966BL

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	323	rBV	149470	203212	2.06%	0.303%
2	5.287	383	391	405	rBV	3661424	5333274	53.96%	7.965%
3	6.851	648	657	670	rBV	3643861	5618632	56.85%	8.391%
4	7.204	710	717	727	rBV	3698619	5767483	58.36%	8.613%
5	7.669	789	796	807	rVB	1244471	1911040	19.34%	2.854%
6	7.987	842	850	858	rBV	2808487	4465574	45.18%	6.669%
7	8.822	984	992	1004	rBV	2082694	3424188	34.65%	5.114%
8	10.445	1260	1268	1280	rBV	1502707	2542807	25.73%	3.797%
9	12.927	1682	1690	1702	rBV	5543821	8119632	82.16%	12.126%
10	14.310	1918	1925	1935	rBV2	2192869	3314916	33.54%	4.951%
11	15.804	2172	2179	2192	rBV	4068142	5764923	58.33%	8.609%
12	17.062	2386	2393	2403	rBV	2658637	3864703	39.10%	5.772%
13	19.698	2835	2841	2851	rBV	7970471	9883072	100.00%	14.759%
14	21.250	3100	3105	3115	rBV	2946792	3697784	37.42%	5.522%
15	23.462	3475	3481	3490	rVB2	1692686	3049953	30.86%	4.555%

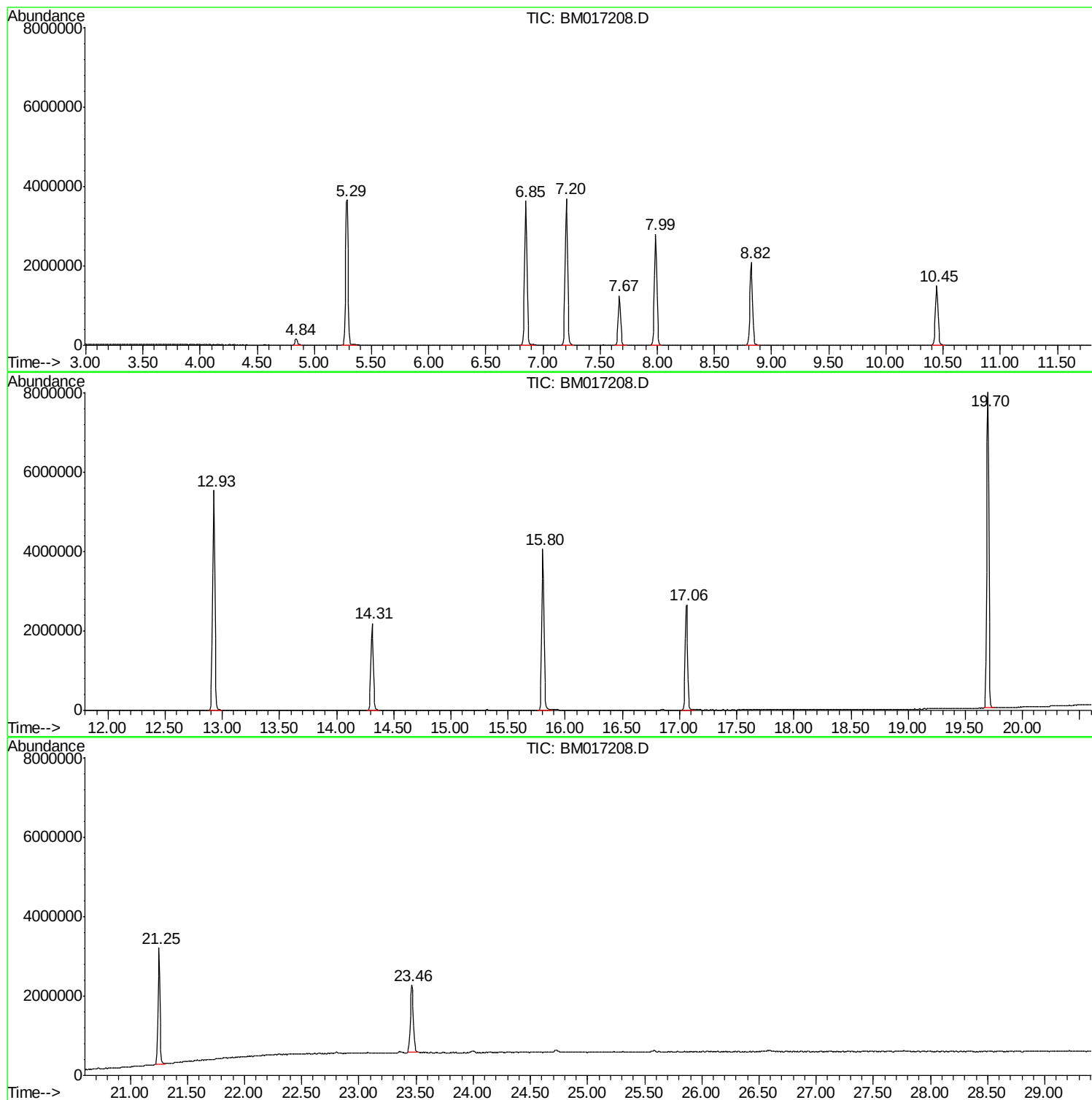
Sum of corrected areas: 66961193

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



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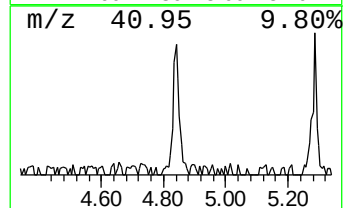
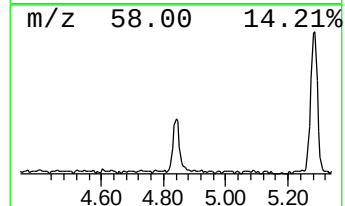
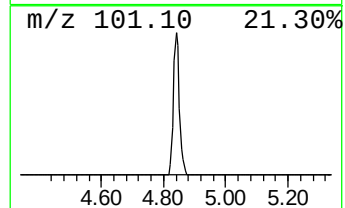
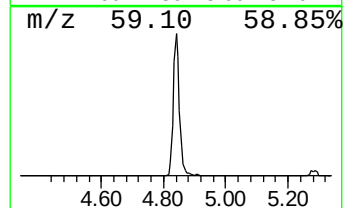
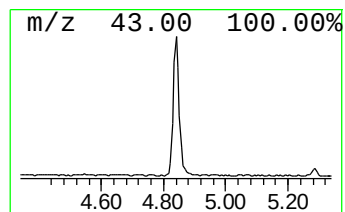
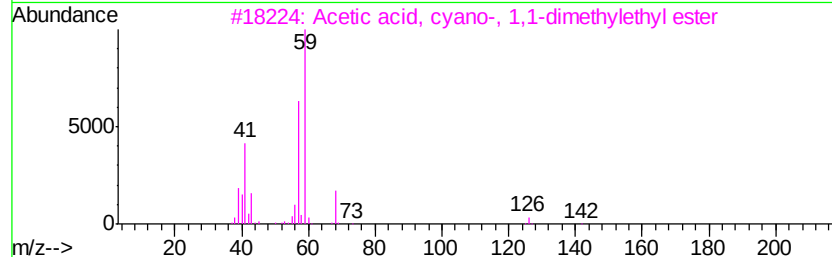
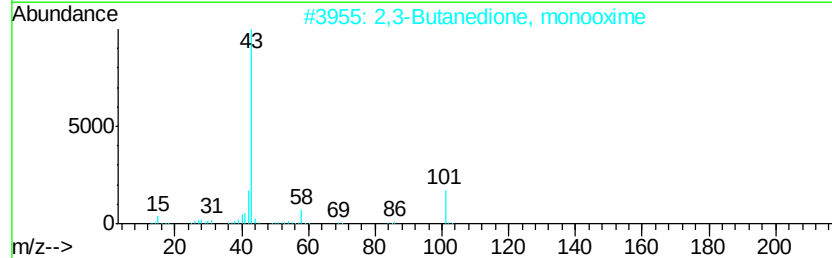
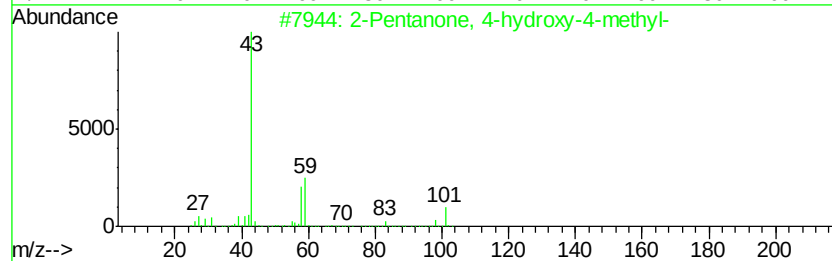
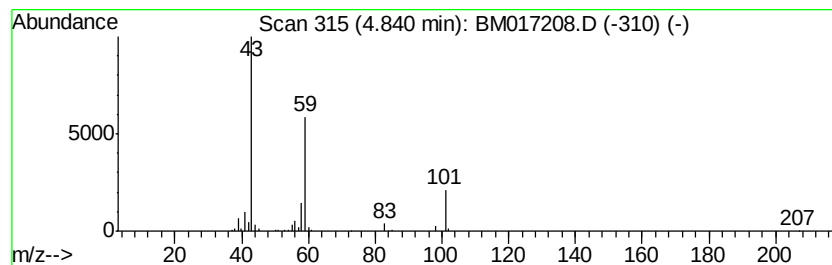
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.84	2.13 ng	203212	1,4-Dichlorobenzene-d4	7.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane	58	C4H10	000106-97-8	9



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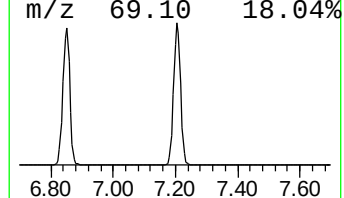
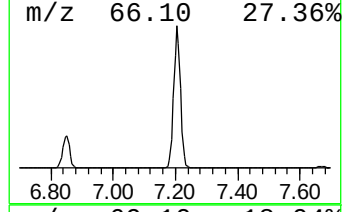
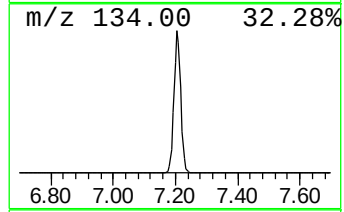
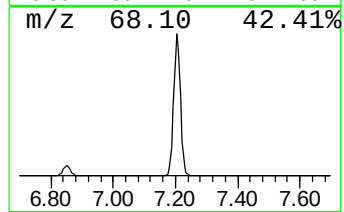
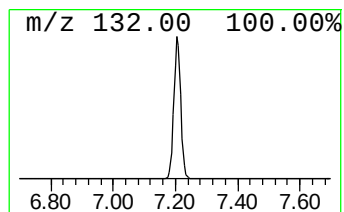
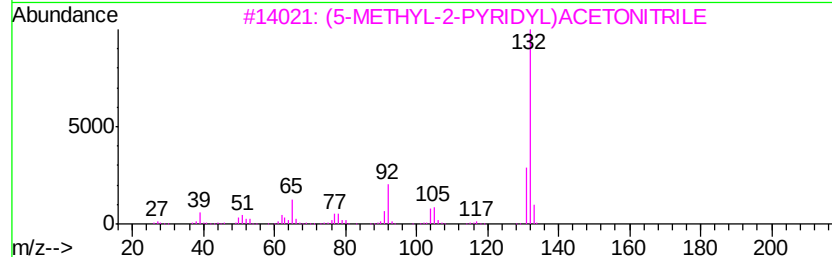
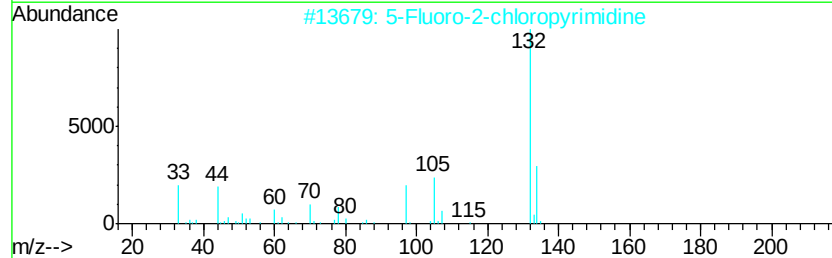
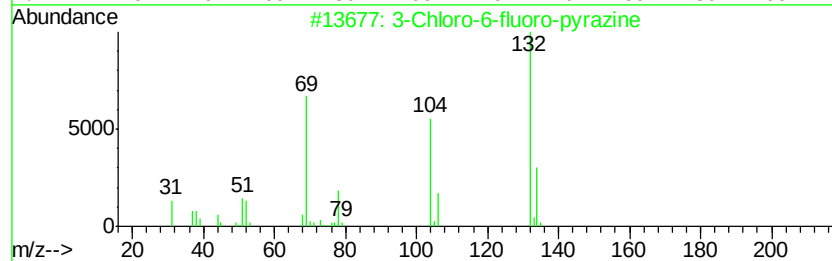
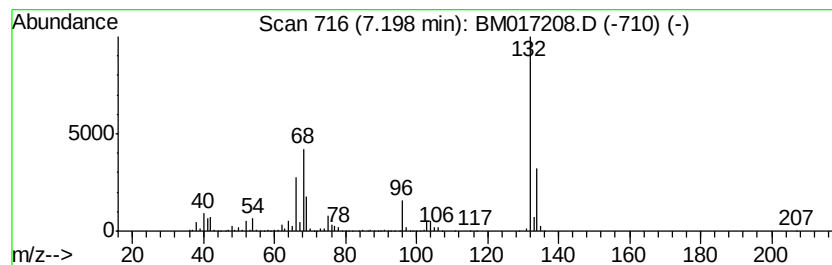
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Peak Number 2 unknown7.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	60.36 ng	5767480	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		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.84	2.1	ng	203212	1	7.67	1911040	20.0
unknown7.20	7.20	60.4	ng	5767480	1	7.67	1911040	20.0