

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060618\
 Data File : BM015506.D
 Acq On : 07 Jun 2018 02:13
 Operator : SJ/JU
 Sample : PB109923BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB109923BL

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-BM052518.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	5.363	347	353	371	rVB	197518	301258	4.34%	0.681%
2	5.845	429	435	447	rBV	2662614	4010117	57.73%	9.065%
3	7.498	708	716	733	rBV	2644675	4244619	61.11%	9.595%
4	7.875	772	780	792	rBV	2772097	4429785	63.77%	10.013%
5	8.351	854	861	871	rBV	629819	1021166	14.70%	2.308%
6	8.681	909	917	927	rBV	2081025	3368644	48.50%	7.615%
7	9.539	1055	1063	1079	rBV	1540422	2578677	37.12%	5.829%
8	11.186	1335	1343	1355	rBV	757726	1298851	18.70%	2.936%
9	13.598	1746	1753	1764	rBV	3716867	5315445	76.52%	12.015%
10	14.968	1980	1986	1997	rVB2	1010065	1540822	22.18%	3.483%
11	16.445	2231	2237	2248	rBV	2680479	3735521	53.78%	8.444%
12	17.692	2443	2449	2456	rBV	1208290	1710431	24.62%	3.866%
13	20.245	2878	2883	2889	rBV	5518397	6946152	100.00%	15.701%
14	21.791	3141	3146	3156	rBV	1440816	1902903	27.40%	4.301%
15	24.215	3551	3558	3566	rVB	911306	1835296	26.42%	4.149%

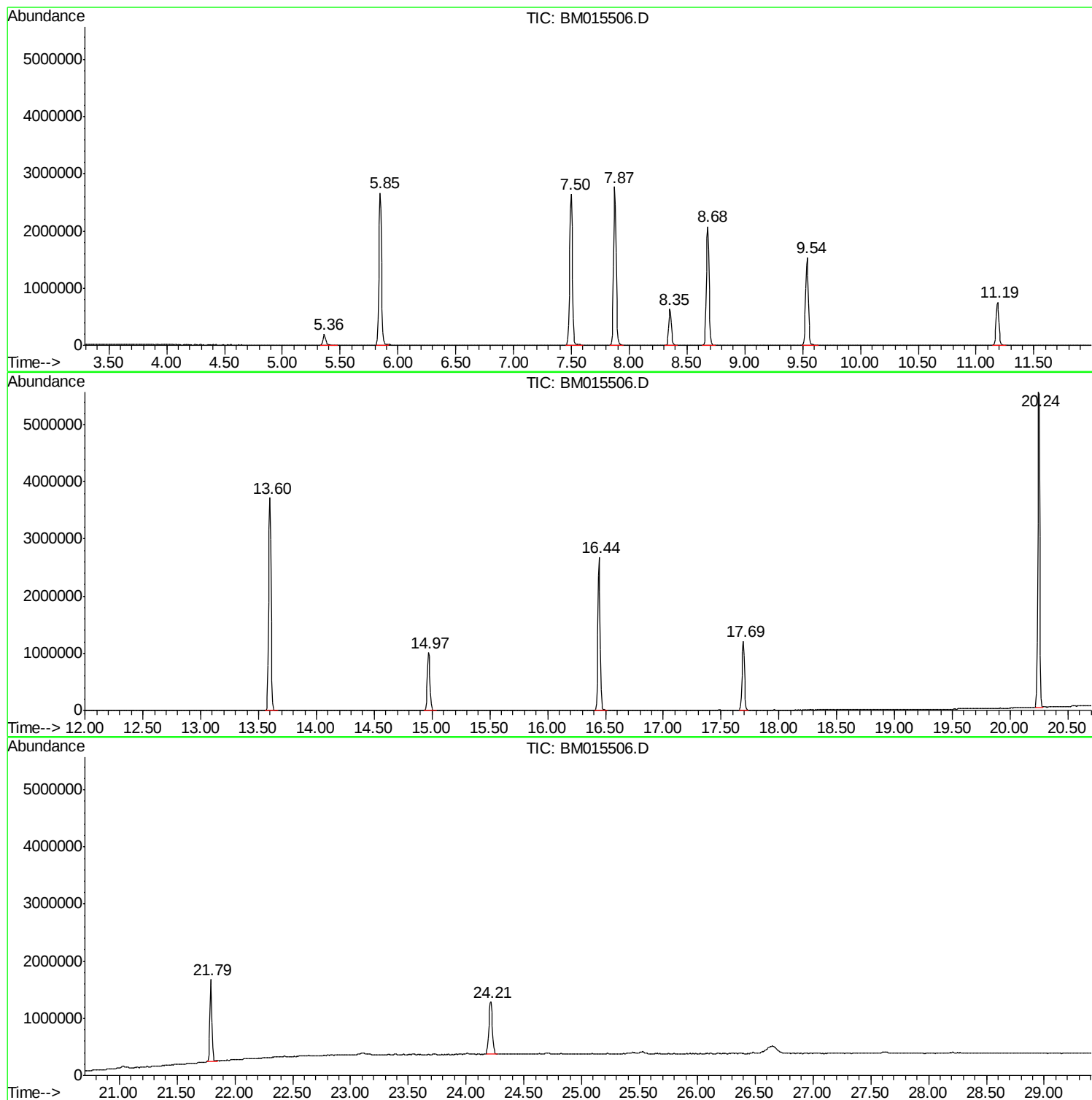
Sum of corrected areas: 44239687

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



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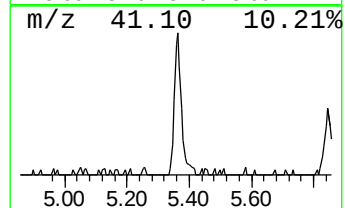
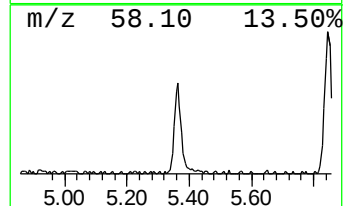
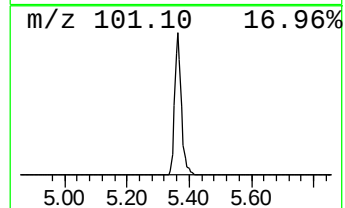
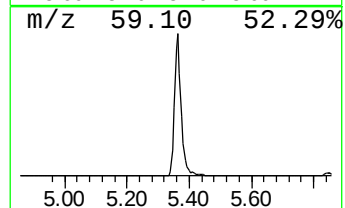
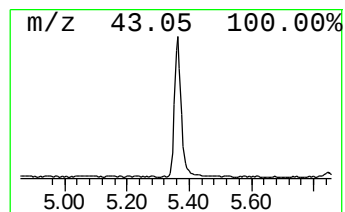
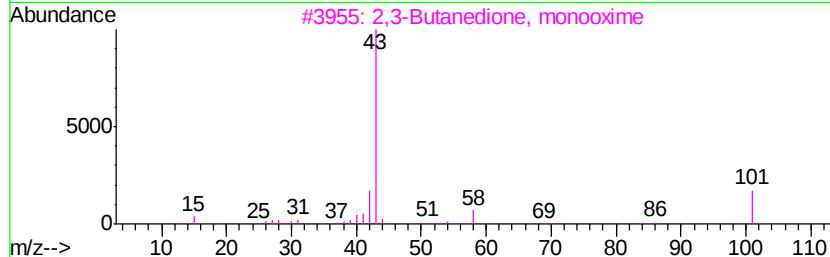
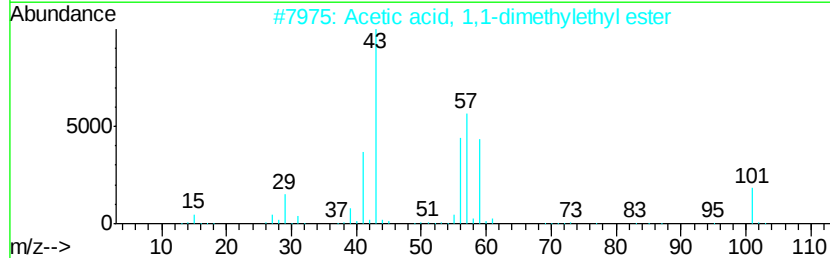
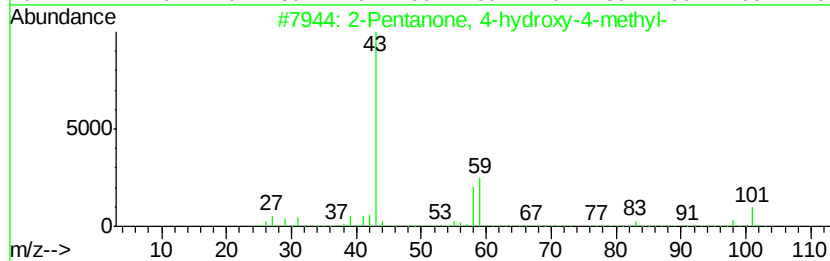
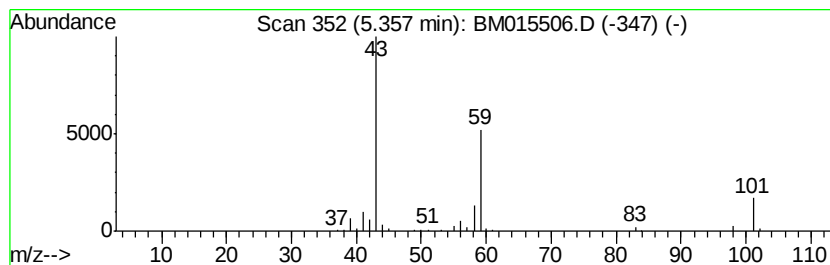
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.36	5.90 ng	301258	1,4-Dichlorobenzene-d4	8.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
4			1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	23
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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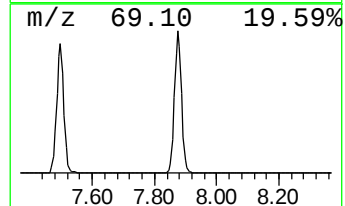
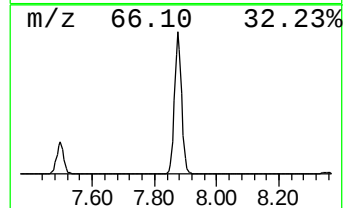
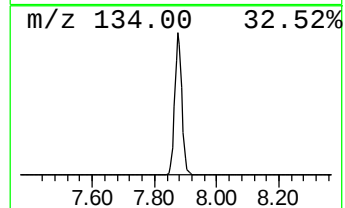
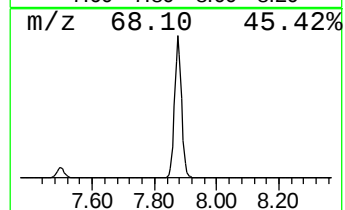
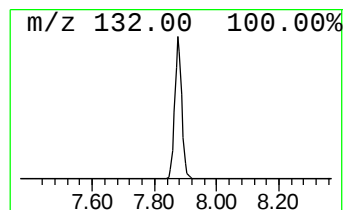
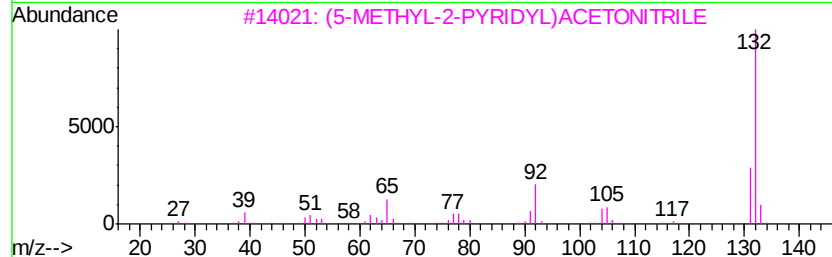
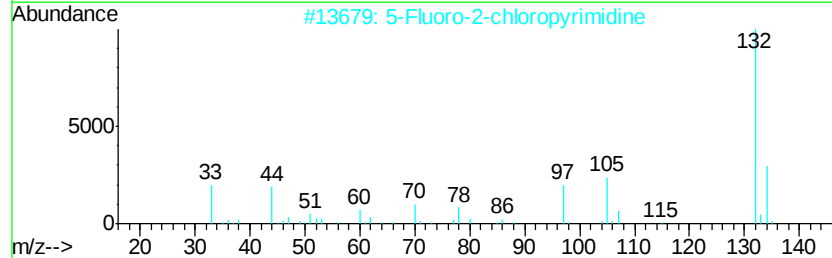
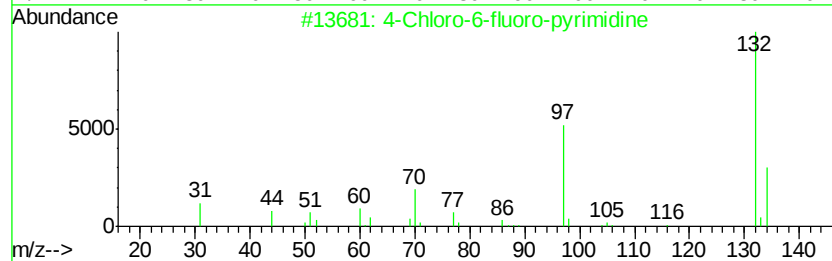
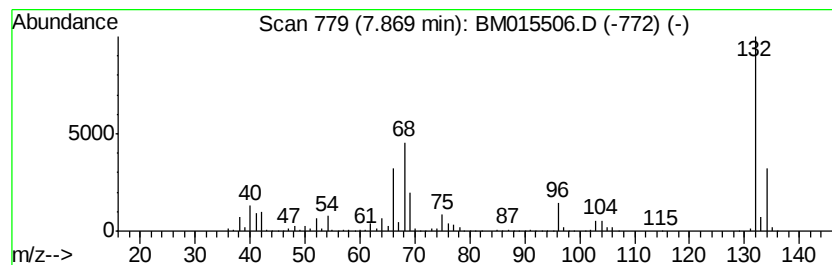
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Peak Number 2 unknown7.87 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.87	86.76 ng	4429790	1,4-Dichlorobenzene-d4	8.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	10
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...	5.36	5.9	ng	301258	1	8.35	1021170	20.0
unknown7.87	7.87	86.8	ng	4429790	1	8.35	1021170	20.0