

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070918\
 Data File : BM015843.D
 Acq On : 09 Jul 2018 15:07
 Operator : SJ/JU
 Sample : PB110790BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB110790BL

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-BM070518.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	5.322	341	346	363	rVB	212959	332999	3.90%	0.844%
2	5.804	422	428	449	rBV	2181610	3289828	38.53%	8.338%
3	7.451	701	708	725	rBV	2128521	3461552	40.54%	8.773%
4	7.828	765	772	789	rBV	2217705	3627114	42.48%	9.193%
5	8.298	846	852	861	rBV	346802	593003	6.95%	1.503%
6	8.628	900	908	923	rBV	1559520	2584992	30.28%	6.552%
7	9.486	1047	1054	1077	rBV	1250698	2146044	25.13%	5.439%
8	11.127	1326	1333	1345	rBV	453776	782581	9.17%	1.983%
9	13.545	1738	1744	1759	rBV	3310149	4748169	55.61%	12.034%
10	14.921	1972	1978	1990	rBV2	666576	1016374	11.90%	2.576%
11	16.398	2223	2229	2246	rBV	3065204	4271526	50.03%	10.826%
12	17.645	2436	2441	2451	rBV	898490	1275677	14.94%	3.233%
13	20.209	2871	2877	2882	rBV	7788935	8538117	100.00%	21.639%
14	21.756	3135	3140	3148	rBV	1097501	1482200	17.36%	3.757%
15	24.150	3541	3547	3561	rBV2	639287	1306090	15.30%	3.310%

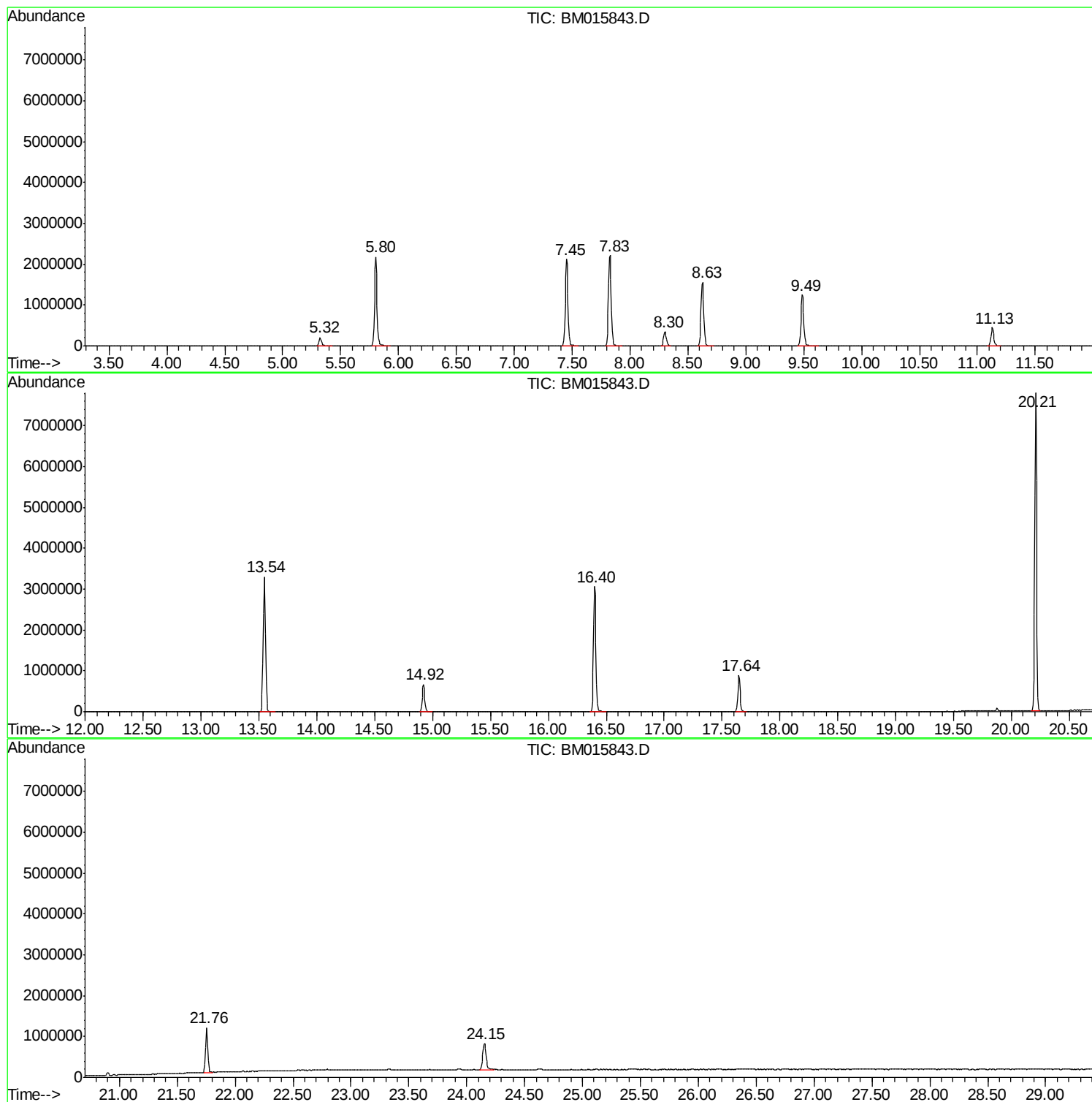
Sum of corrected areas: 39456266

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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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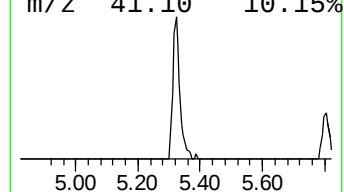
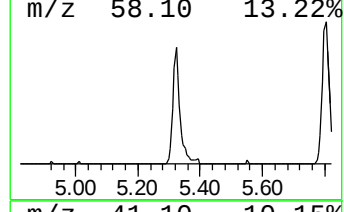
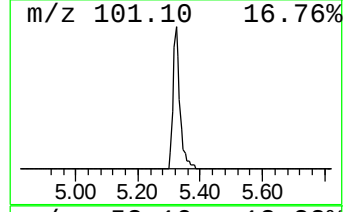
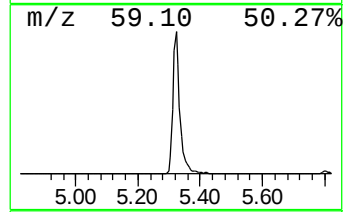
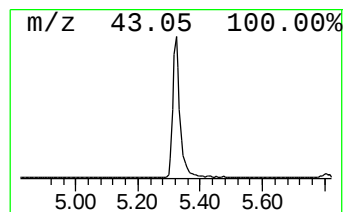
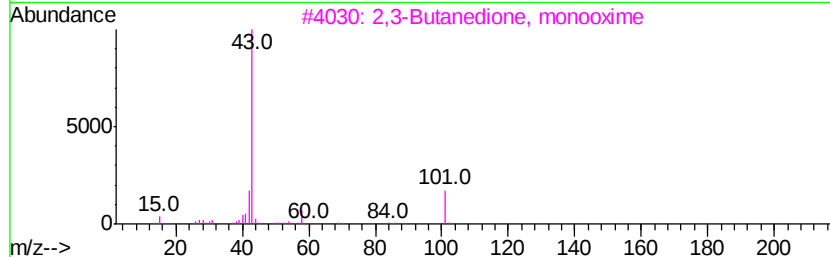
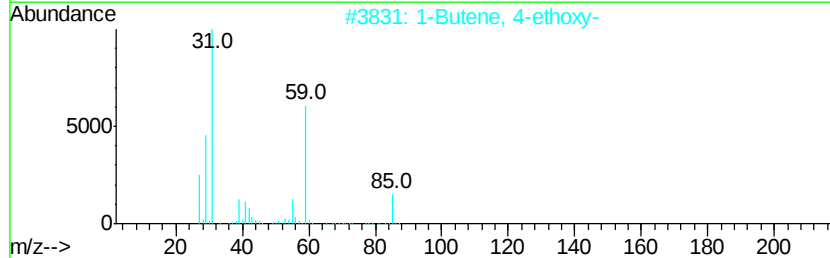
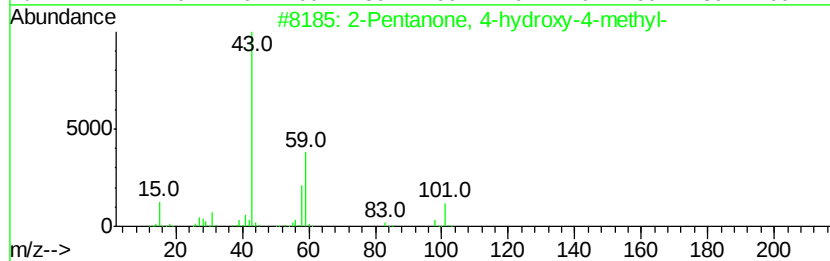
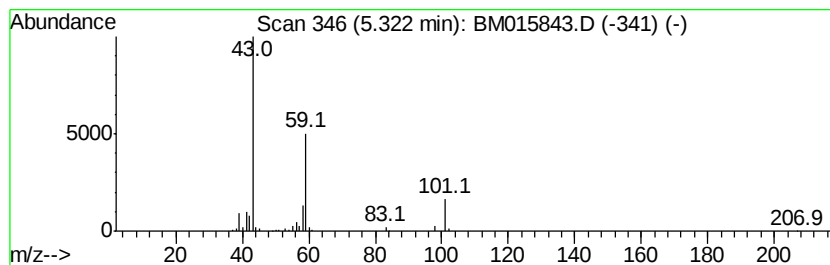
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.32	11.23 ng	332999	1,4-Dichlorobenzene-d4	8.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	23
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			Butane	58	C4H10	000106-97-8	9
5			Guanidine	59	CH5N3	000113-00-8	9



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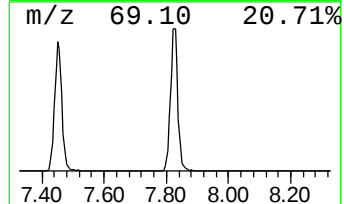
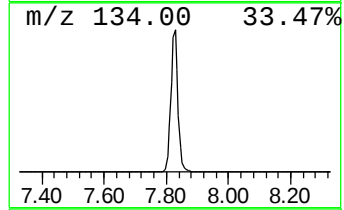
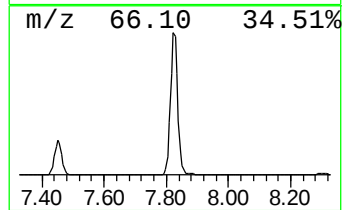
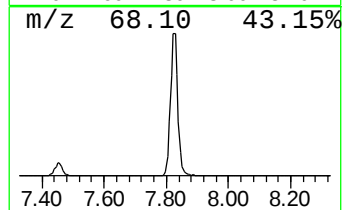
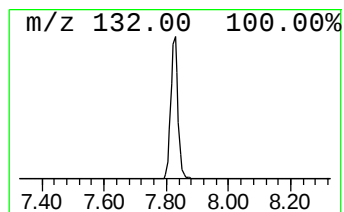
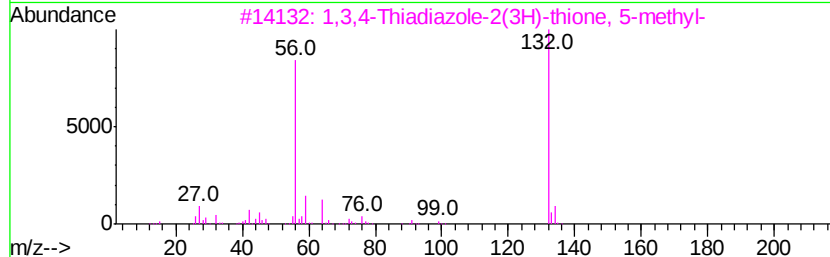
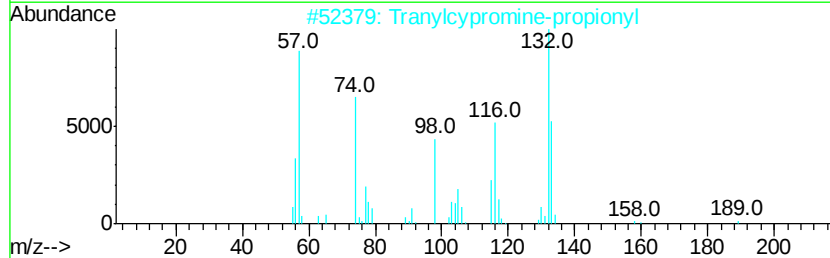
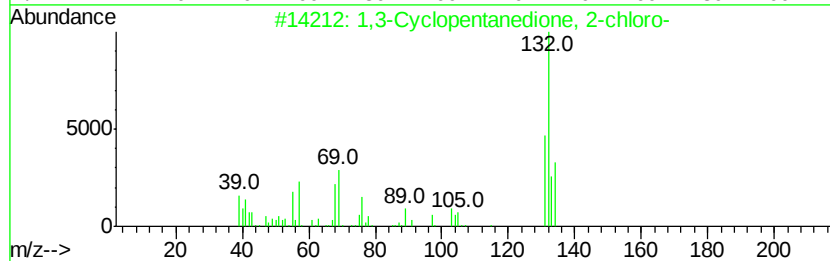
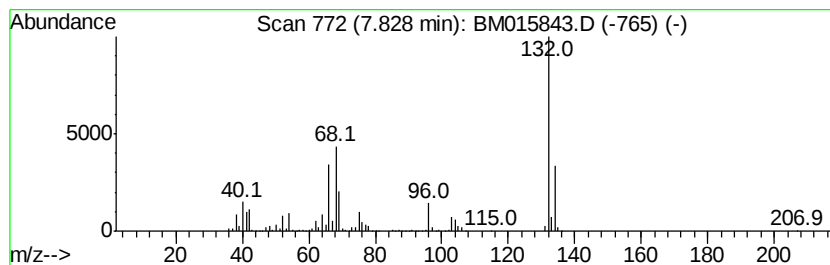
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Peak Number 2 unknown7.83 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.83	122.33 ng	3627110	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
3		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.32	11.2	ng	332999	1	8.30	593003	20.0
unknown7.83	7.83	122.3	ng	3627110	1	8.30	593003	20.0