

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070918\
 Data File : BM015845.D
 Acq On : 09 Jul 2018 16:22
 Operator : SJ/JU
 Sample : PB110833BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB110833BL

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	356	rBV	185821	284610	4.16%	0.837%
2	5.804	422	428	450	rBV	1768122	2697837	39.47%	7.929%
3	7.451	701	708	724	rBV	1834454	2887263	42.24%	8.486%
4	7.822	765	771	792	rBV	1837678	3040707	44.49%	8.937%
5	8.298	846	852	867	rVB	391195	654907	9.58%	1.925%
6	8.628	900	908	922	rBV	1325910	2217909	32.45%	6.519%
7	9.486	1047	1054	1072	rBV	1048510	1756115	25.69%	5.161%
8	11.127	1323	1333	1346	rBV	501297	849044	12.42%	2.495%
9	13.545	1738	1744	1755	rBV	2661682	3864451	56.54%	11.358%
10	14.921	1972	1978	1998	rVB2	752178	1126752	16.49%	3.312%
11	16.398	2223	2229	2246	rBV	2540276	3483928	50.97%	10.240%
12	17.645	2435	2441	2451	rBV	978919	1423612	20.83%	4.184%
13	20.209	2871	2877	2884	rBV	5897817	6834713	100.00%	20.088%
14	21.756	3135	3140	3148	rBV	1195654	1568538	22.95%	4.610%
15	24.156	3541	3548	3556	rVB	684838	1333349	19.51%	3.919%

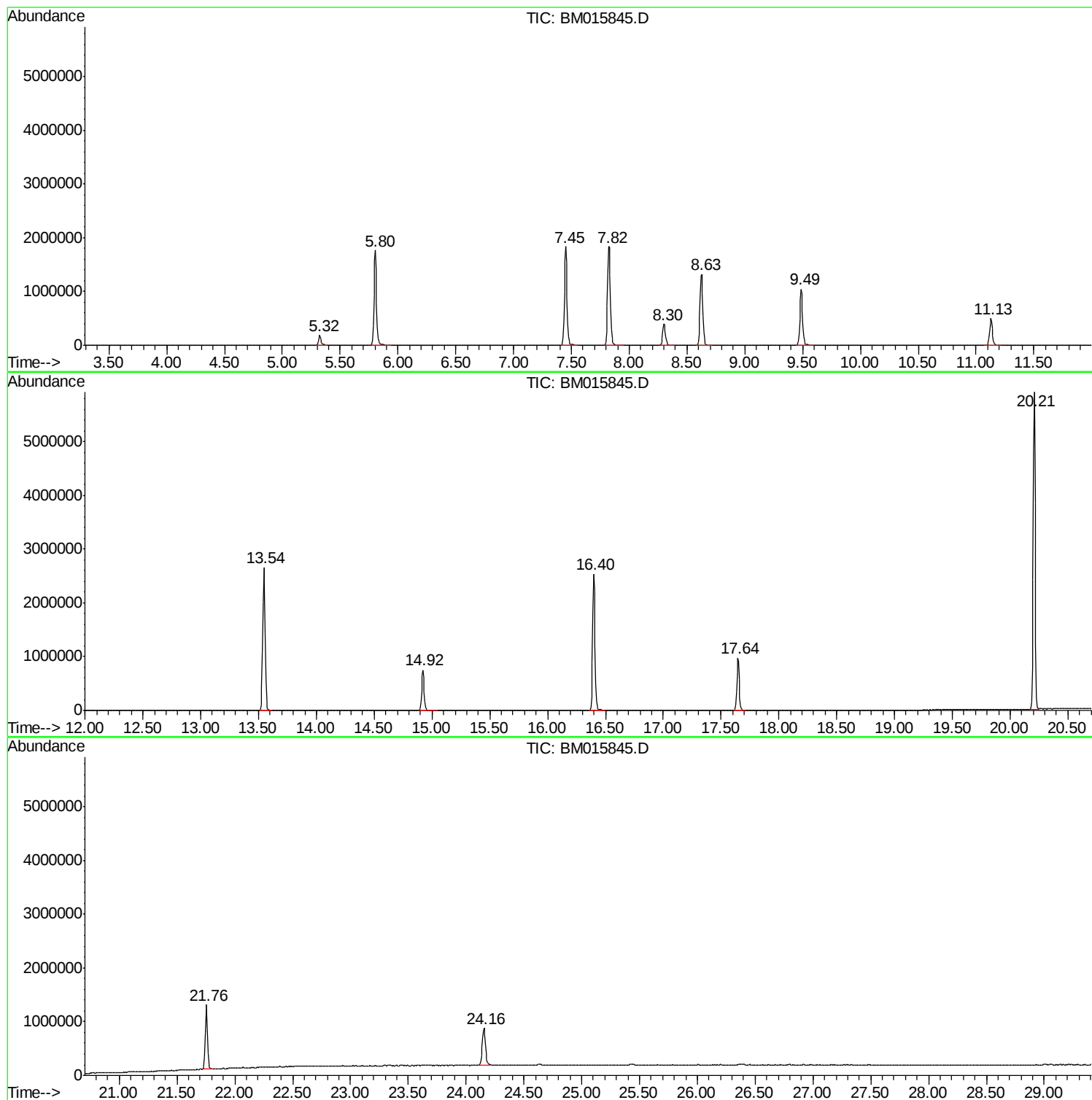
Sum of corrected areas: 34023735

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



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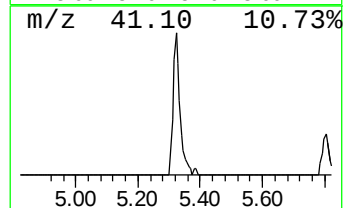
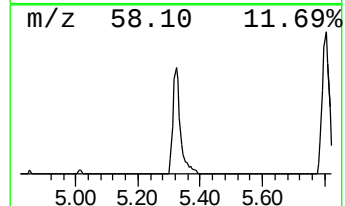
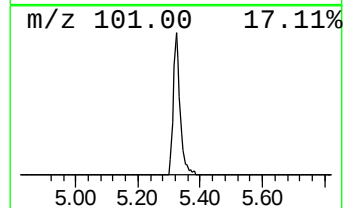
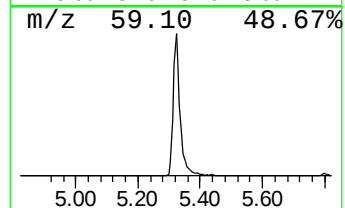
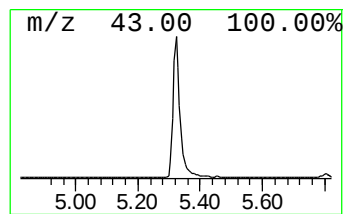
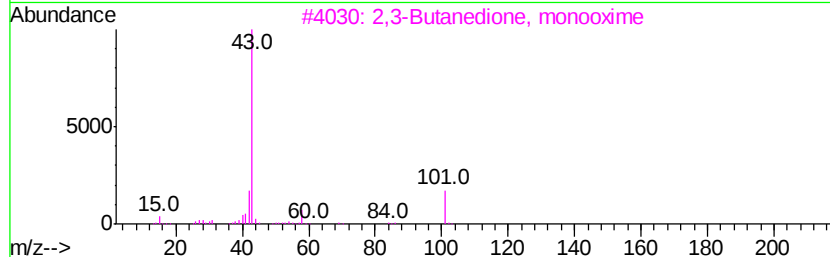
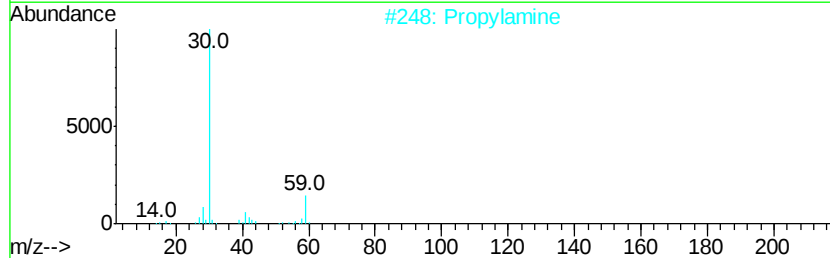
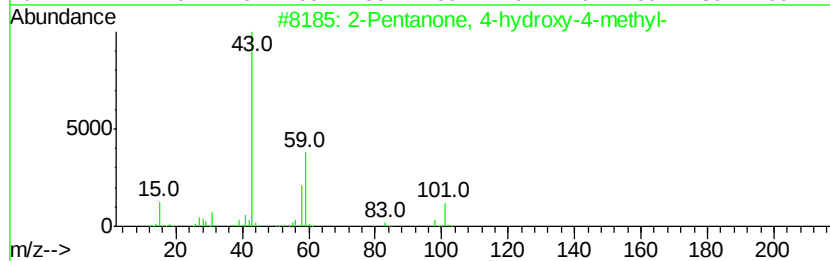
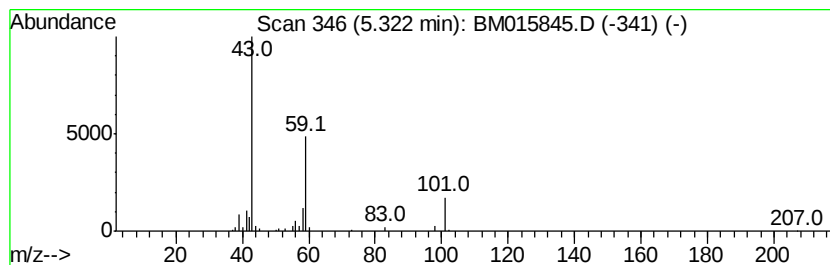
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.32	8.69 ng	284610	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			Propylamine	59	C3H9N	000107-10-8	40
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			Allyl acetate	100	C5H8O2	000591-87-7	10
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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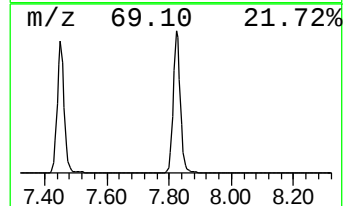
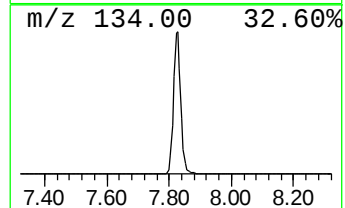
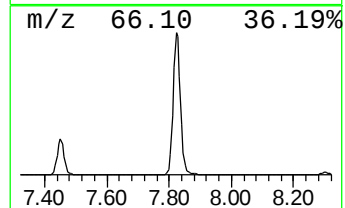
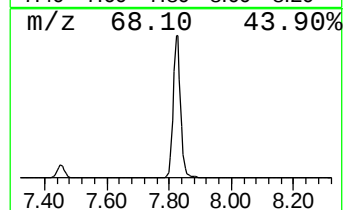
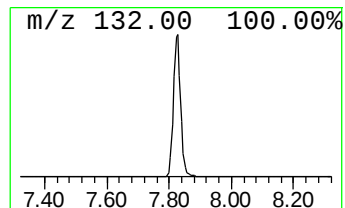
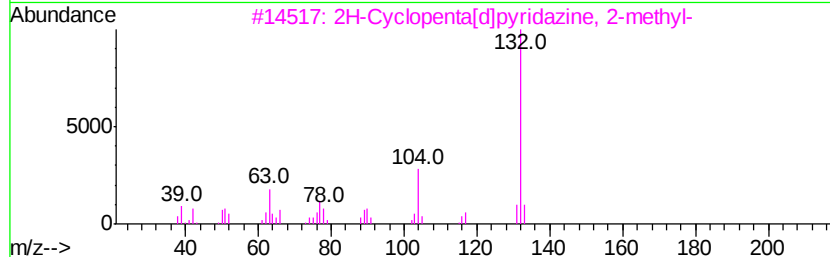
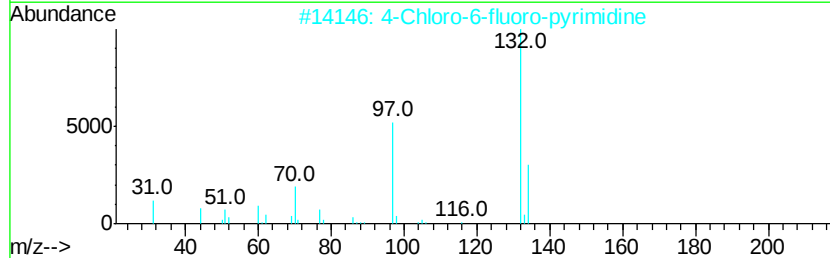
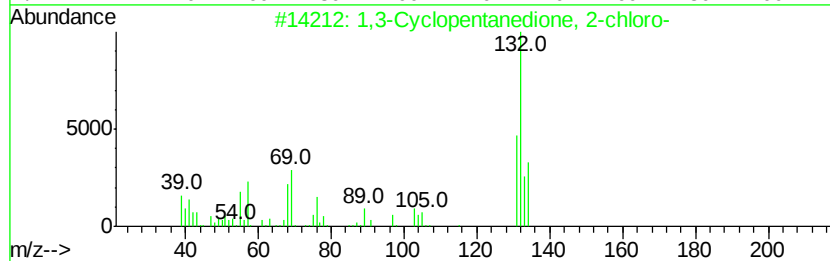
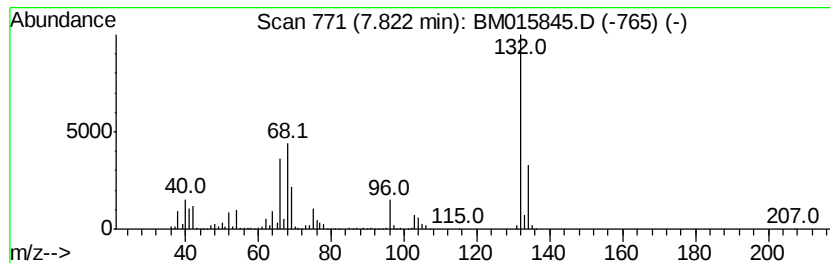
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Peak Number 2 unknown7.82 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.82	92.86 ng	3040710	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	35
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	27
4		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	22
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.32	8.7	ng	284610	1	8.30	654907	20.0
unknown7.82	7.82	92.9	ng	3040710	1	8.30	654907	20.0