

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082917\
 Data File : BF098110.D
 Acq On : 29 Aug 2017 14:55
 Operator : SJ/JU
 Sample : PB101845BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB101845BL

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:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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	4.945	481	486	496	rBV	194474	298644	7.34%	0.929%
2	5.351	549	555	558	rBV	3145140	3371762	82.90%	10.490%
3	6.375	723	729	732	rBV	3177257	3407362	83.78%	10.601%
4	6.504	746	751	754	rBV	3313795	3324944	81.75%	10.344%
5	6.734	786	790	793	rBV	863874	764635	18.80%	2.379%
6	6.887	812	816	820	rBV	2660738	2605459	64.06%	8.106%
7	7.298	880	886	889	rBV	2035684	2214854	54.46%	6.891%
8	8.016	1003	1008	1016	rBV	1013467	1031229	25.36%	3.208%
9	9.092	1185	1191	1194	rBV	3350831	3876985	95.32%	12.062%
10	9.769	1301	1306	1317	rVB	1241068	1163185	28.60%	3.619%
11	10.557	1434	1440	1443	rBV	2204648	2379819	58.51%	7.404%
12	11.251	1552	1558	1561	rBV	1501339	1284226	31.58%	3.995%
13	12.845	1822	1829	1832	rBV	3681379	4067127	100.00%	12.653%
14	13.886	2001	2006	2009	rBV	1332130	1246143	30.64%	3.877%
15	15.304	2241	2247	2251	rBV	746161	910604	22.39%	2.833%
16	15.745	2317	2322	2328	rBV2	36083	52532	1.29%	0.163%
17	16.221	2398	2403	2408	rVB	34030	53851	1.32%	0.168%
18	16.768	2492	2496	2504	rVB4	26193	46479	1.14%	0.145%
19	17.421	2602	2607	2616	rVB6	18877	43033	1.06%	0.134%

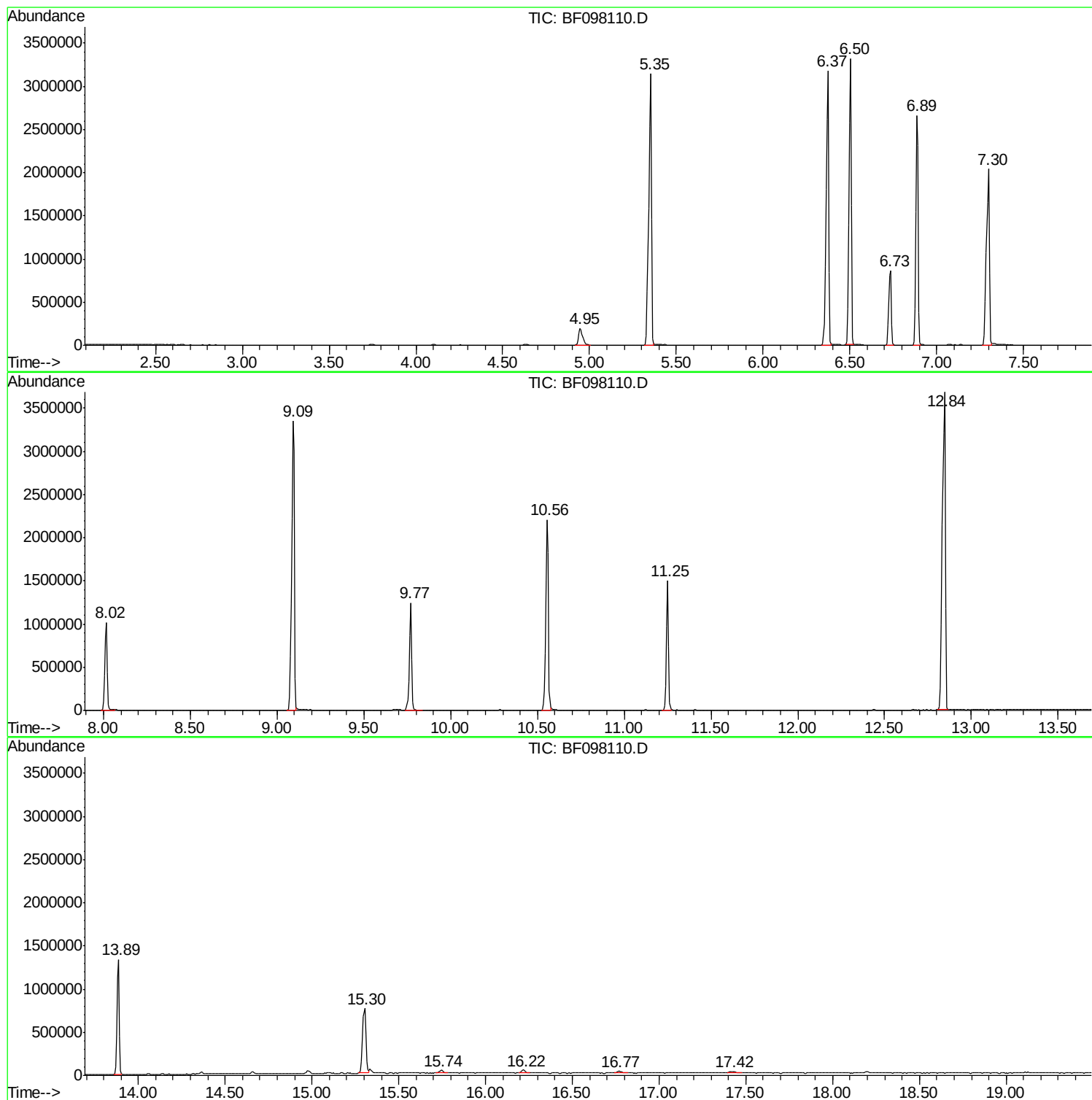
Sum of corrected areas: 32142873

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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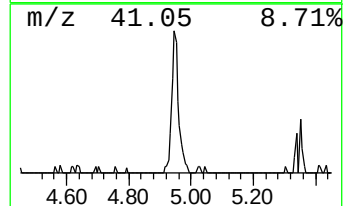
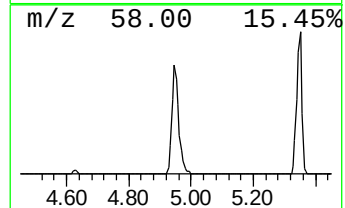
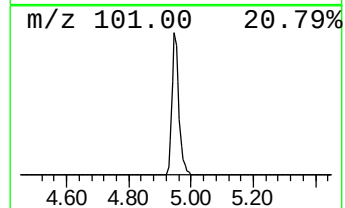
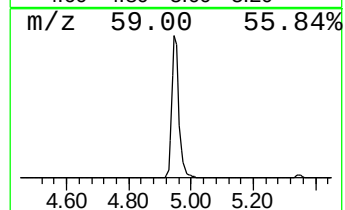
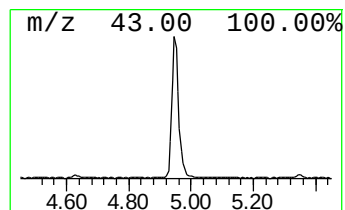
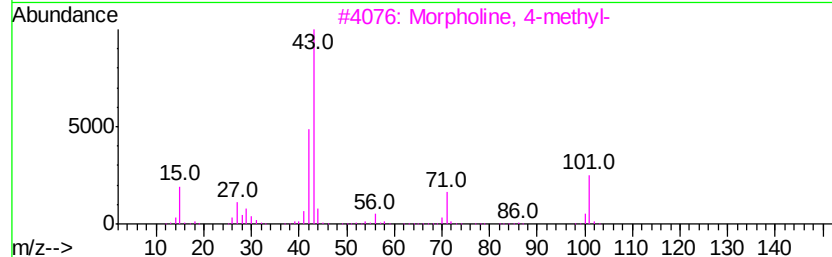
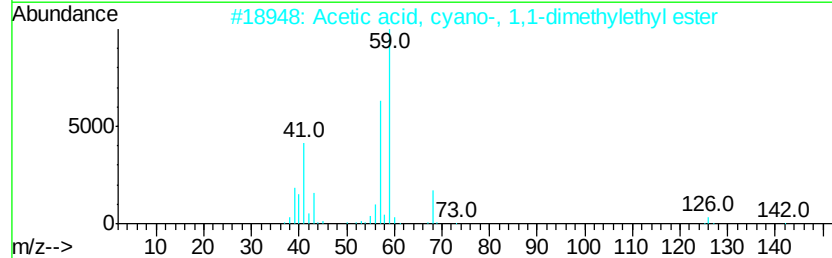
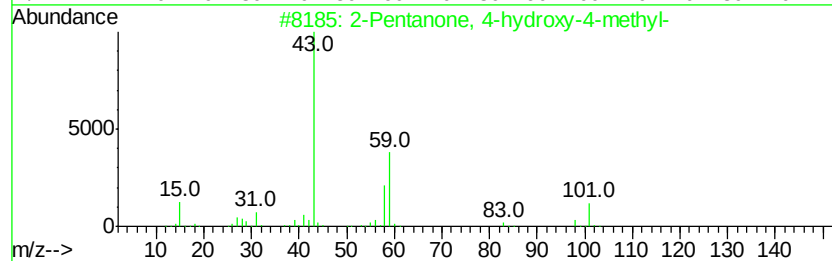
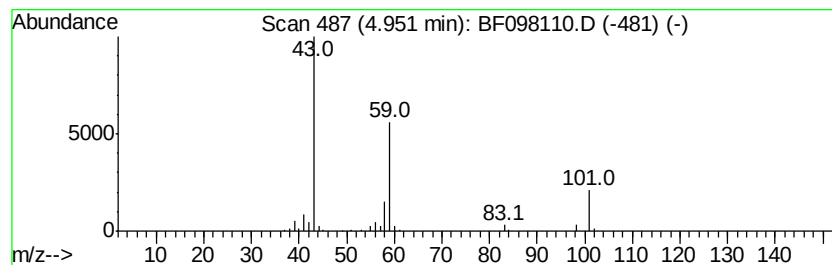
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	7.81 ng	298644	1,4-Dichlorobenzene-d4	6.73

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, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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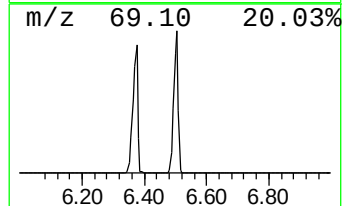
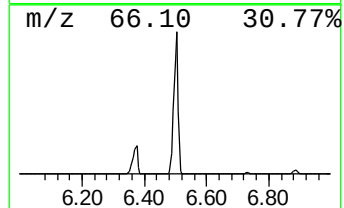
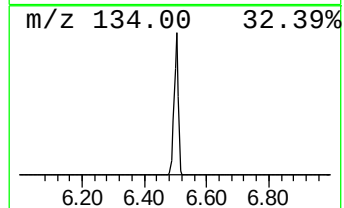
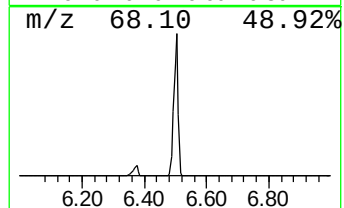
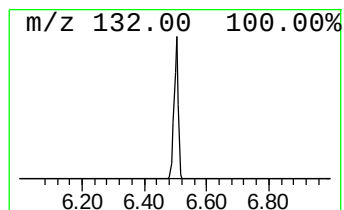
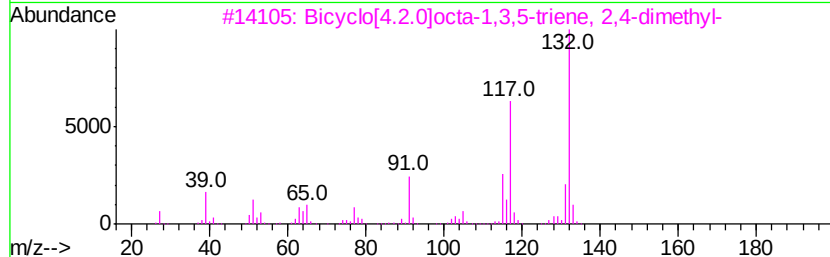
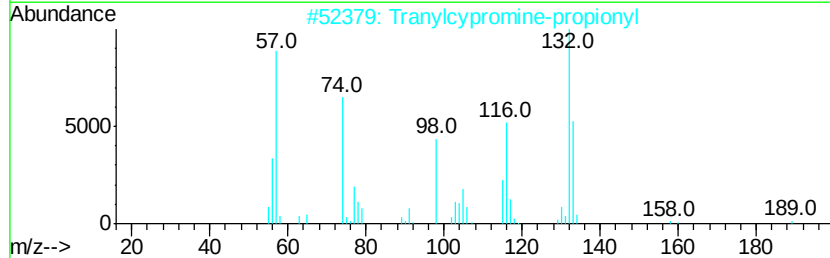
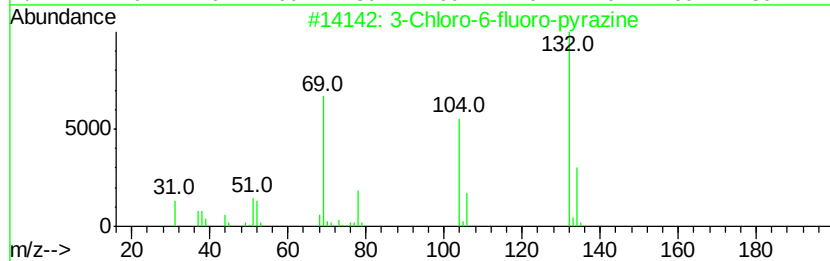
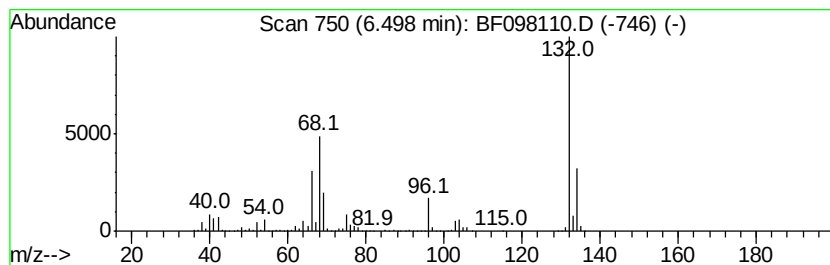
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Peak Number 2 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	86.97 ng	3324940	1,4-Dichlorobenzene-d4	6.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.95	7.8	ng	298644	1	6.73	764635	20.0
unknown6.50	6.50	87.0	ng	3324940	1	6.73	764635	20.0