

Data Path : Z:\HPCHEM1\BNA M\DATA\BM052417\
 Data File : BM010173.D
 Acq On : 24 May 2017 10:39
 Operator : SJ/MA
 Sample : PB99196BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB99196BL

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_M\METHODS\8270-BM051917.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	5.075	334	338	346	rBV	175205	253113	2.58%	0.558%
2	5.534	410	416	431	rVB	2599446	3655396	37.25%	8.055%
3	7.140	682	689	702	rBV	2275779	3506185	35.73%	7.726%
4	7.493	742	749	761	rBV	2475564	3810066	38.82%	8.396%
5	7.951	821	827	835	rBV	504761	793554	8.09%	1.749%
6	8.275	874	882	892	rBV	1924988	3082593	31.41%	6.793%
7	9.140	1022	1029	1040	rBV	1287293	2079376	21.19%	4.582%
8	10.757	1297	1304	1313	rBV	555791	930589	9.48%	2.051%
9	13.204	1713	1720	1733	rBV	3606603	5259585	53.59%	11.590%
10	14.586	1948	1955	1963	rBV2	791608	1210978	12.34%	2.669%
11	16.074	2202	2208	2218	rBV	2974530	4077000	41.54%	8.984%
12	17.327	2415	2421	2428	rBV	1076453	1573080	16.03%	3.466%
13	19.933	2858	2864	2870	rBV	8067446	9814025	100.00%	21.626%
14	21.486	3124	3128	3134	rBV	1995301	2501564	25.49%	5.512%
15	23.850	3524	3530	3540	rVB	1471680	2833296	28.87%	6.243%

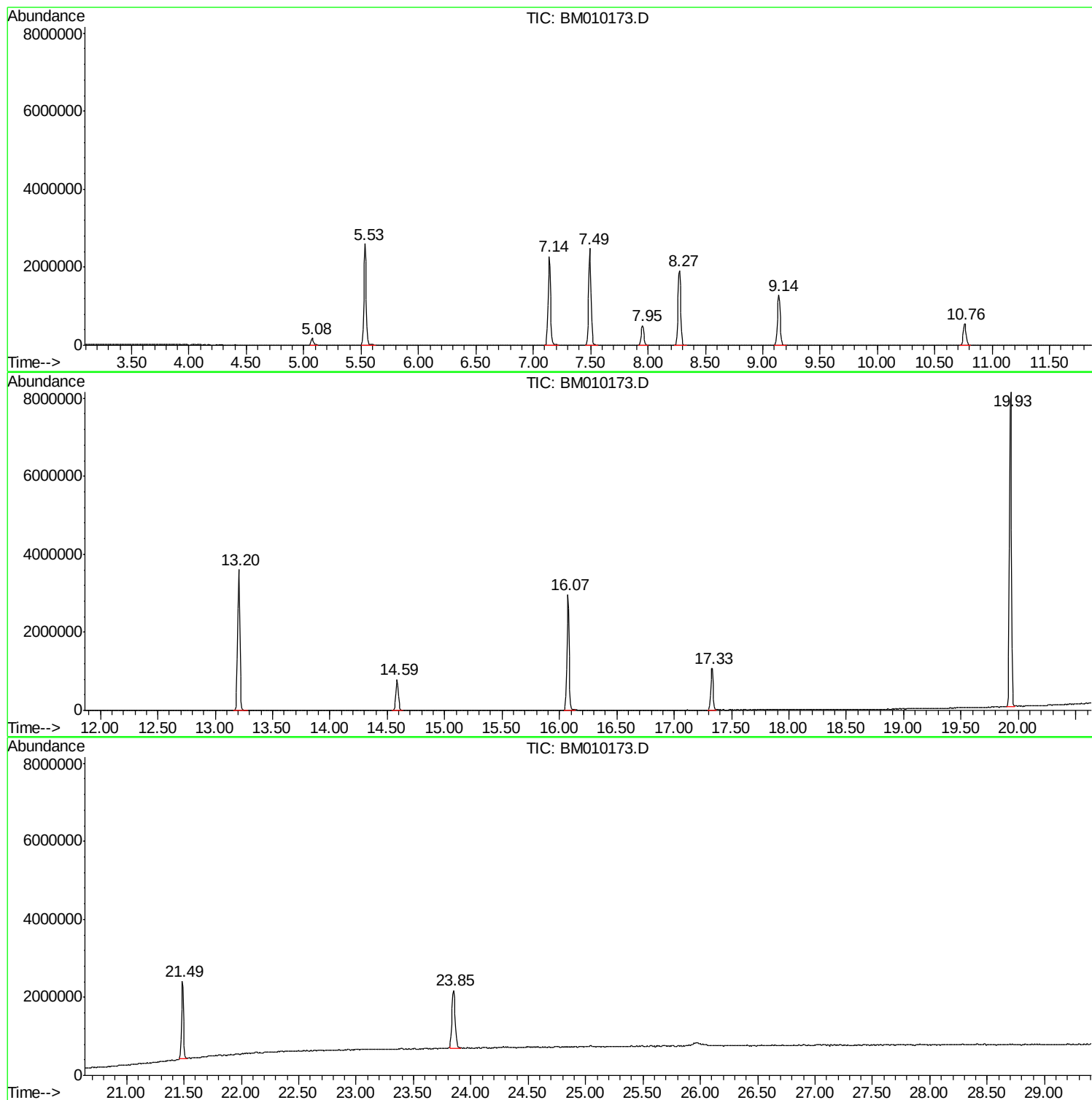
Sum of corrected areas: 45380400

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Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051917.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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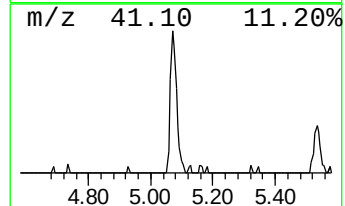
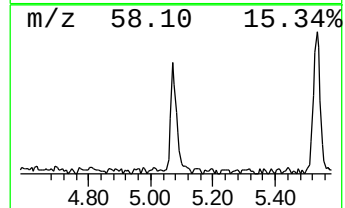
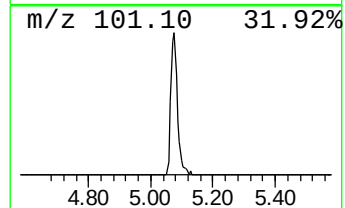
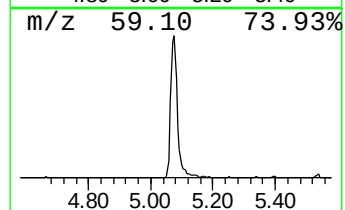
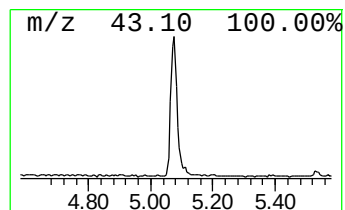
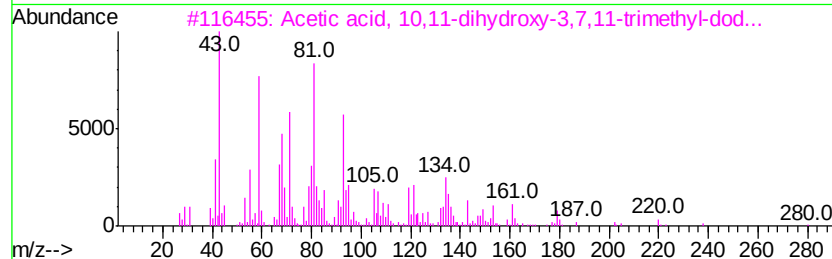
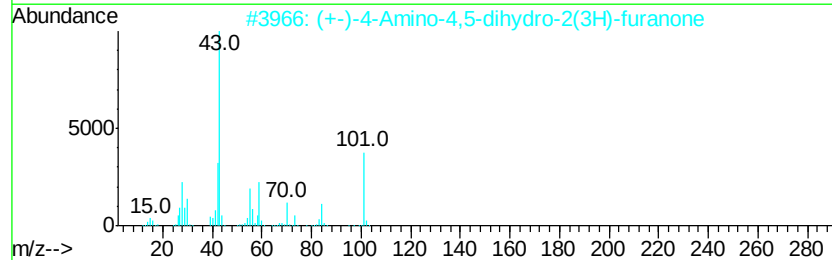
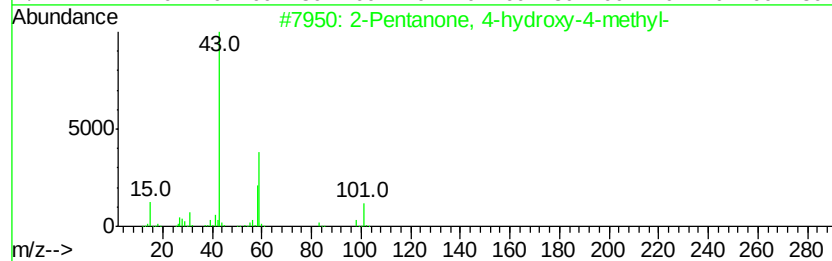
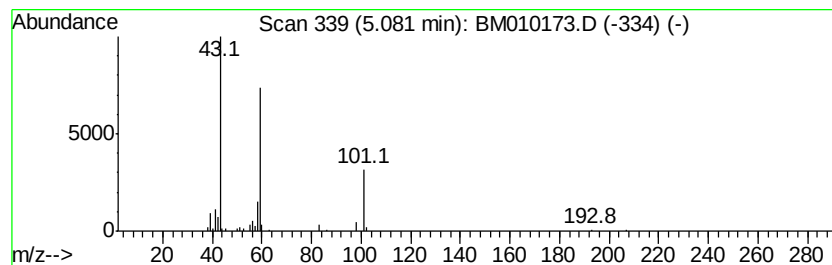
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	6.38 ng	253113	1,4-Dichlorobenzene-d4	7.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	28
3			Acetic acid, 10,11-dihydroxy-3,7,11-trimethyl-dodecanoic	298	C17H30O4	1000194-28-5	23
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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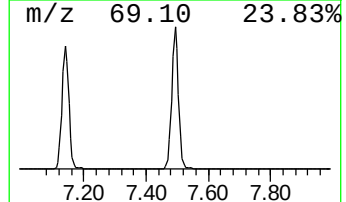
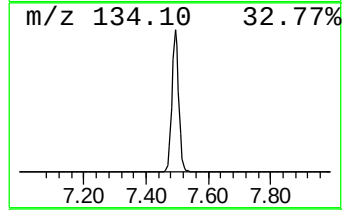
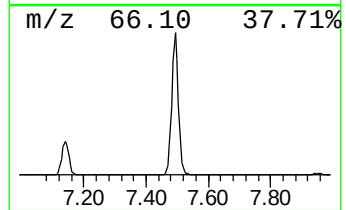
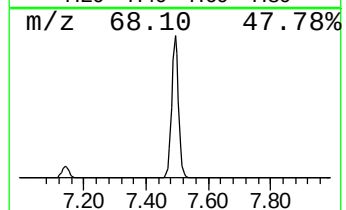
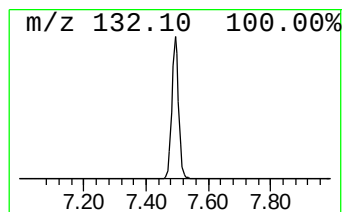
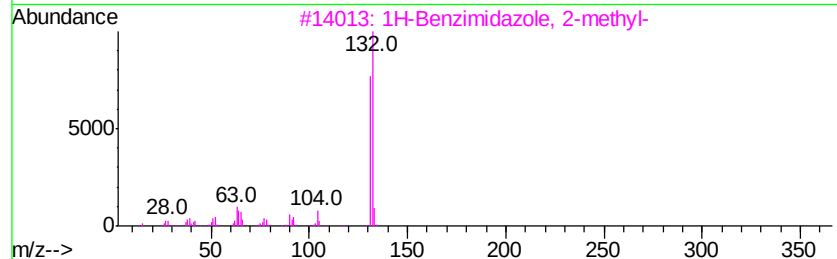
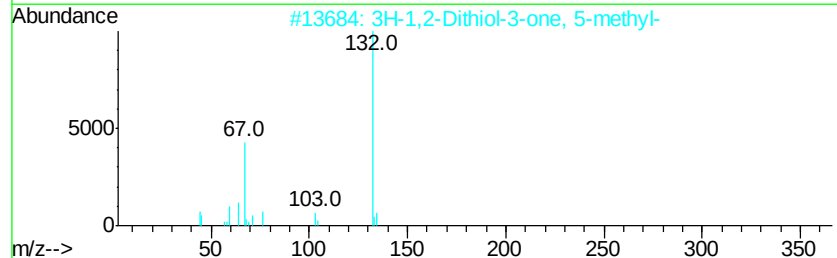
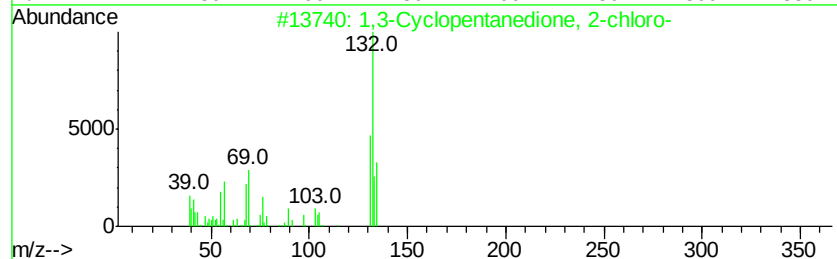
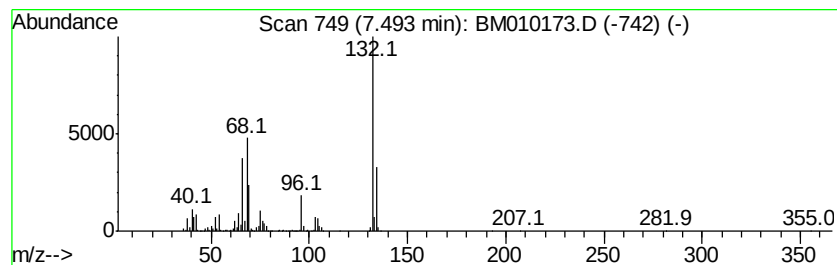
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Peak Number 2 unknown7.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.49	96.03 ng	3810070	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	43
2		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	27
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	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...	5.08	6.4	ng	253113	1	7.95	793554	20.0
unknown7.49	7.49	96.0	ng	3810070	1	7.95	793554	20.0