

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070317\
 Data File : BG027839.D
 Acq On : 3 Jul 2017 19:33
 Operator : SJ/JU
 Sample : PB100333BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB100333BL

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_G\METHODS\8270-BG062817.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	5.577	330	338	349	rVB	70370	112252	3.67%	0.899%
2	6.030	408	415	425	rBV	484007	843282	27.56%	6.753%
3	7.598	673	682	694	rBV	483527	897243	29.32%	7.185%
4	7.986	737	748	763	rBV	482356	885117	28.93%	7.088%
5	8.450	821	827	835	rVB	92199	160905	5.26%	1.288%
6	8.779	875	883	892	rVB	353304	636165	20.79%	5.094%
7	9.619	1017	1026	1037	rBV	270994	528995	17.29%	4.236%
8	11.270	1299	1307	1317	rBV	134593	258363	8.44%	2.069%
9	13.662	1707	1714	1725	rBV	875221	1435154	46.90%	11.492%
10	15.042	1942	1949	1960	rVB2	218113	379318	12.40%	3.037%
11	16.529	2194	2202	2213	rBV	850695	1354201	44.26%	10.844%
12	17.792	2410	2417	2426	rBV2	366695	576305	18.83%	4.615%
13	20.372	2849	2856	2863	rBV	2314859	3059892	100.00%	24.503%
14	22.140	3149	3157	3169	rBV	361790	691684	22.60%	5.539%
15	25.730	3754	3768	3784	rBV2	205838	669120	21.87%	5.358%

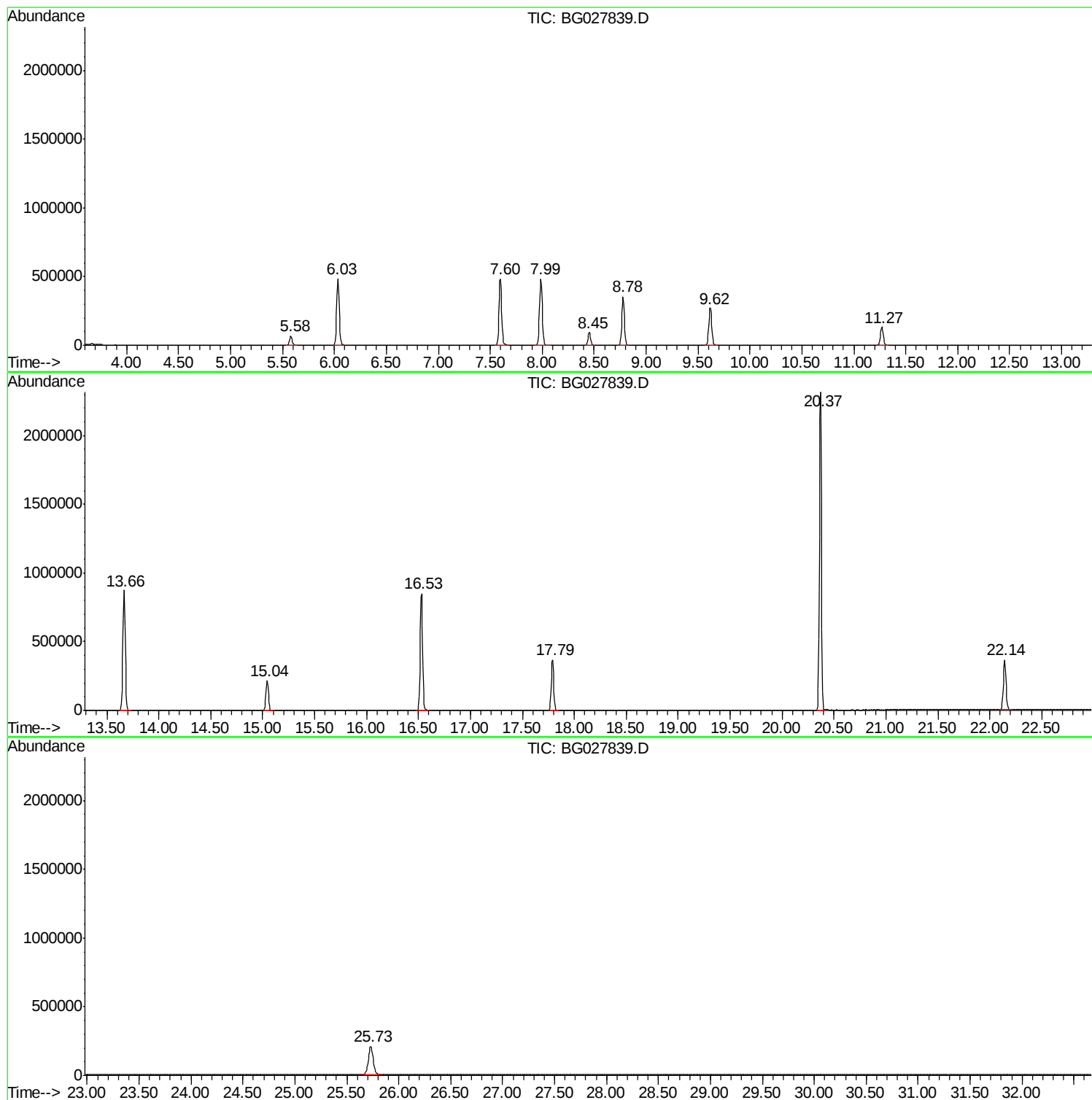
Sum of corrected areas: 12487996

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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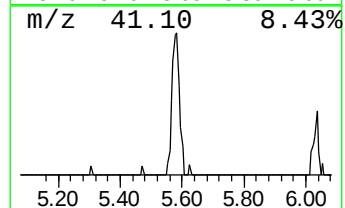
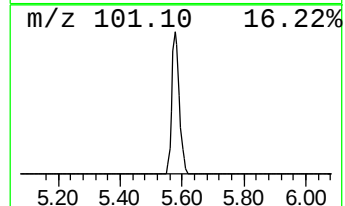
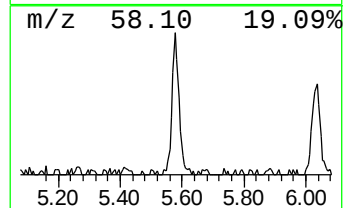
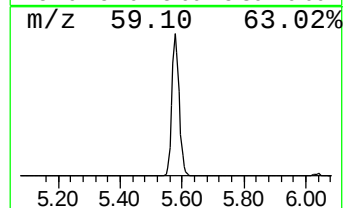
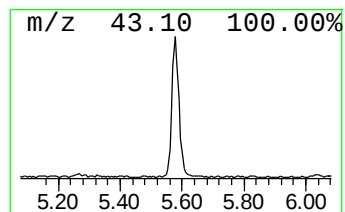
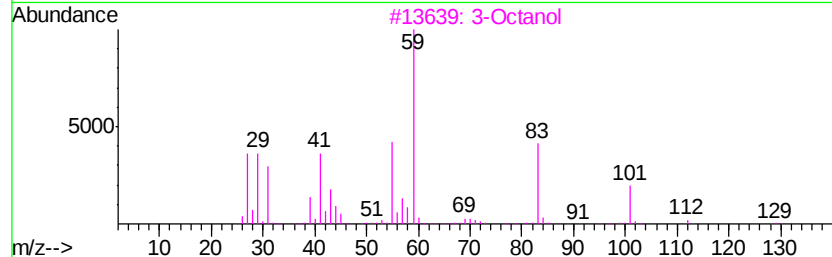
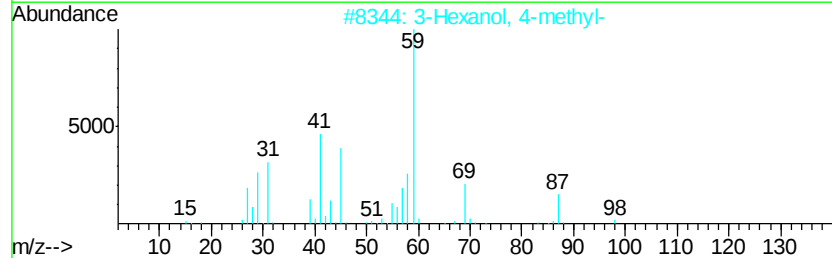
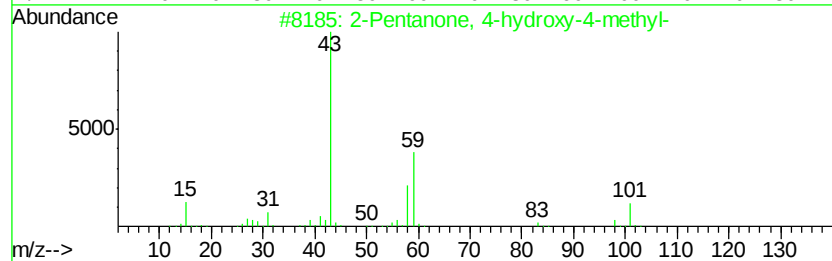
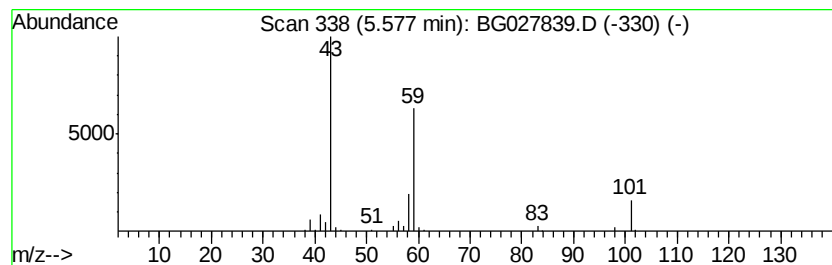
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.58	13.95 ng	112252	1,4-Dichlorobenzene-d4	8.46

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		3-Octanol	130	C8H18O	000589-98-0	28
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12



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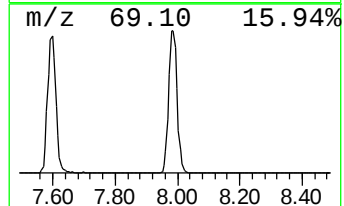
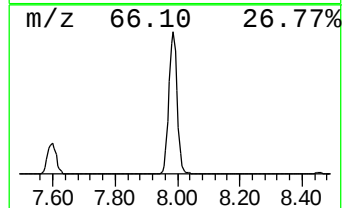
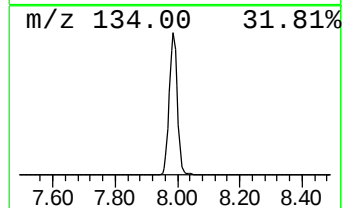
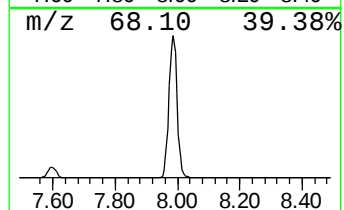
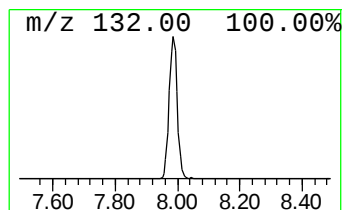
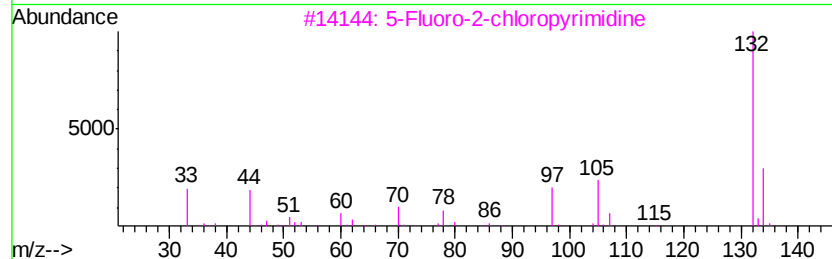
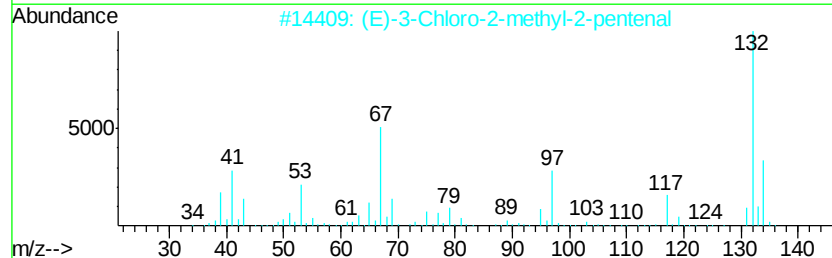
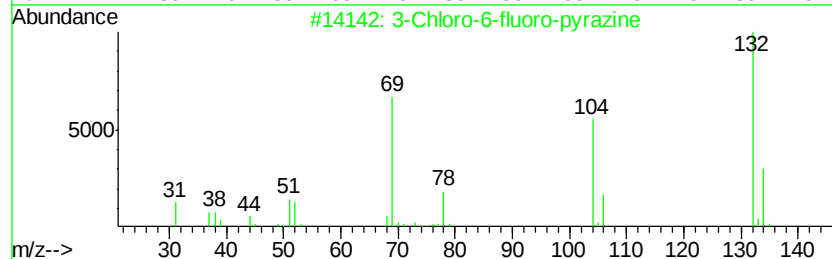
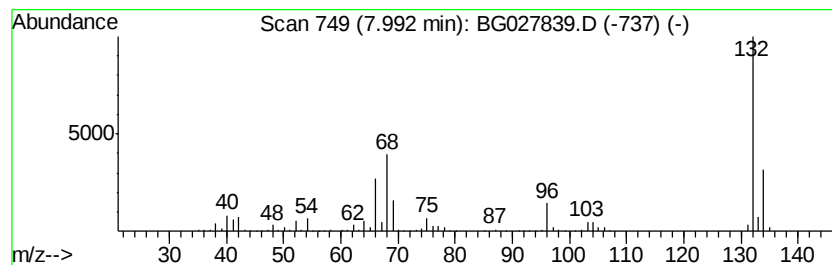
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Peak Number 2 unknown7.99 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.99	110.02 ng	885117	1,4-Dichlorobenzene-d4	8.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.58	13.9	ng	112252	1	8.46	160905	20.0
unknown7.99	7.99	110.0	ng	885117	1	8.46	160905	20.0