

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062617\
 Data File : BG027677.D
 Acq On : 26 Jun 2017 16:32
 Operator : SJ/JU
 Sample : PB100146BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB100146BL

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-BG062117.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	5.627	290	296	306	rVB	41310	66102	3.22%	0.840%
2	6.080	365	373	385	rVB	316851	542899	26.42%	6.897%
3	7.637	630	638	649	rBV	324616	589708	28.69%	7.492%
4	8.030	697	705	715	rBV	307501	563693	27.43%	7.161%
5	8.494	775	784	793	rBB	49747	87099	4.24%	1.107%
6	8.817	831	839	847	rBB	212103	389120	18.93%	4.943%
7	9.658	972	982	991	rBV	180741	336749	16.39%	4.278%
8	11.309	1255	1263	1270	rBV	71673	131116	6.38%	1.666%
9	13.700	1663	1670	1681	rBB	570143	873134	42.48%	11.093%
10	15.075	1897	1904	1911	rVB2	120711	202610	9.86%	2.574%
11	16.561	2150	2157	2169	rBV	565059	899911	43.79%	11.433%
12	17.825	2364	2372	2379	rBV	220652	334312	16.27%	4.247%
13	20.404	2804	2811	2818	rBV	1544547	2055207	100.00%	26.110%
14	22.190	3108	3115	3122	rBV	219195	408141	19.86%	5.185%
15	25.809	3717	3731	3746	rBV	125388	391584	19.05%	4.975%

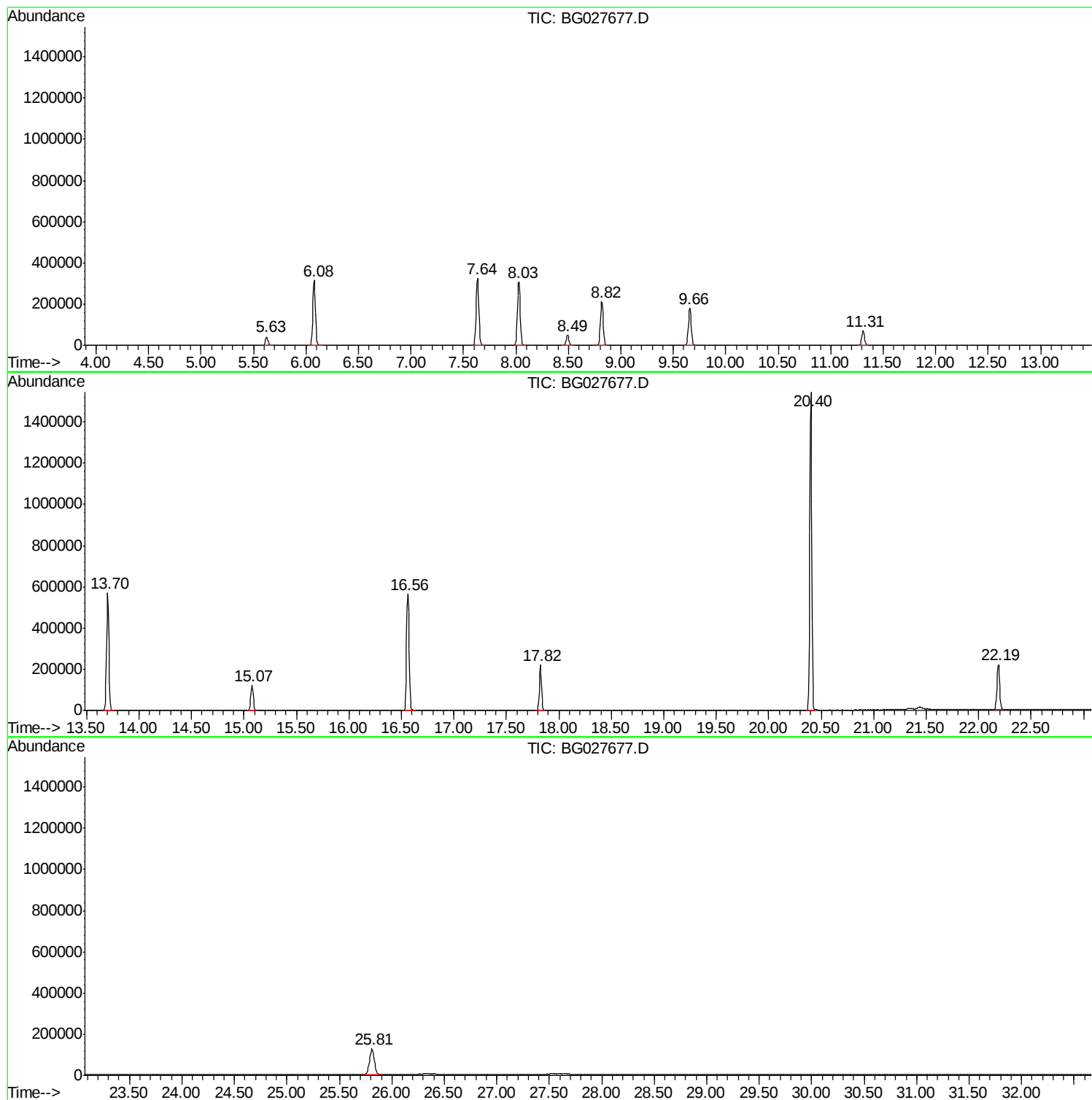
Sum of corrected areas: 7871385

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062117.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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Instrument :
BNA_G
ClientSampled :
PB100146BL

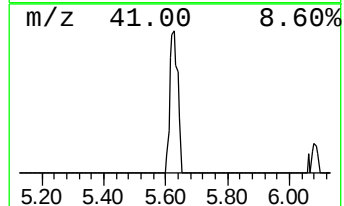
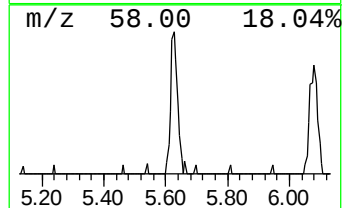
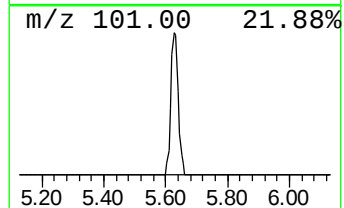
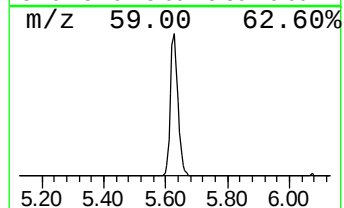
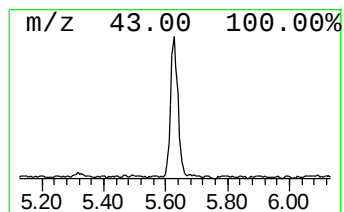
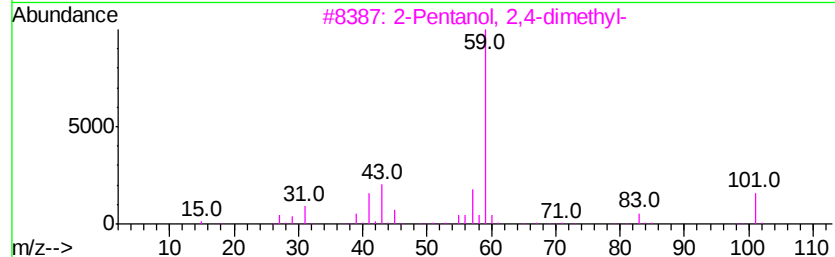
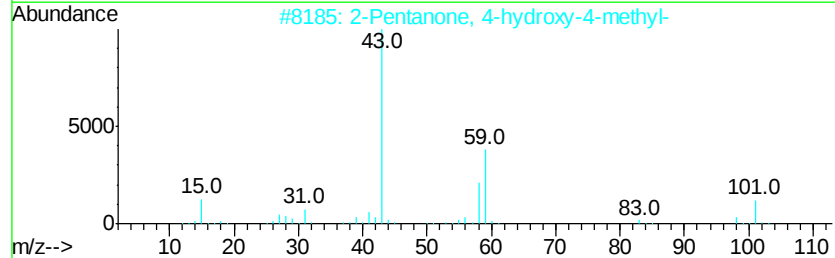
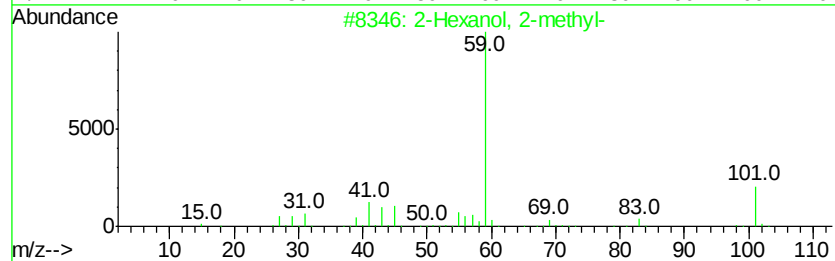
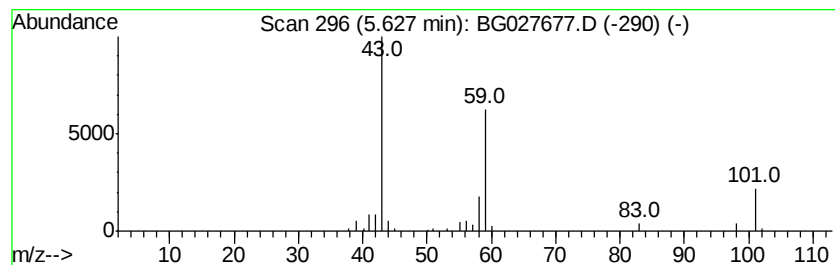
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Hexanol, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	15.18 ng	66102	1,4-Dichlorobenzene-d4	8.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	53
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	42
4			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25



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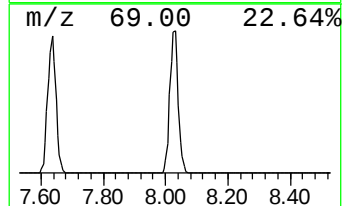
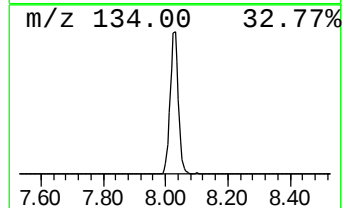
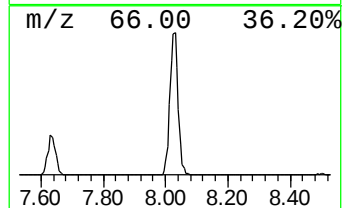
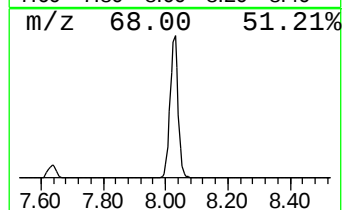
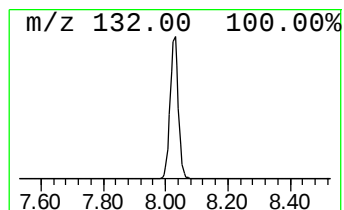
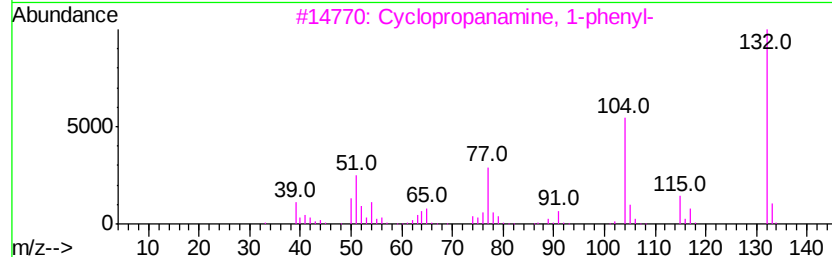
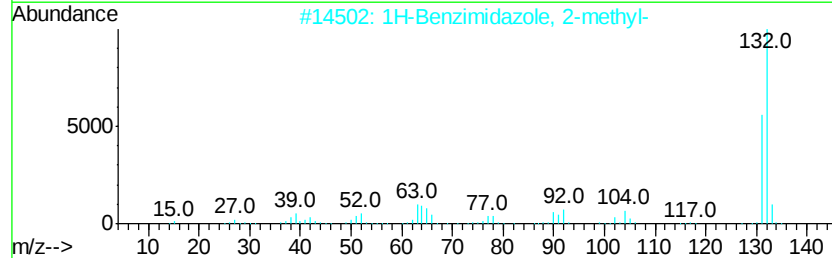
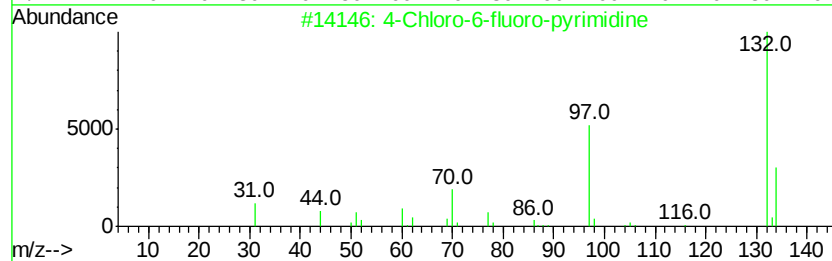
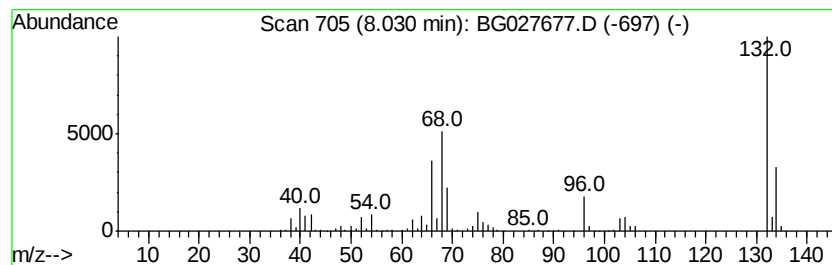
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Peak Number 2 unknown8.03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.03	129.44 ng	563693	1,4-Dichlorobenzene-d4	8.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Hexanol, 2-methyl-	5.63	15.2	ng	66102	1	8.49	87099	20.0
unknown8.03	8.03	129.4	ng	563693	1	8.49	87099	20.0