

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG032018\
 Data File : BG033277.D
 Acq On : 20 Mar 2018 22:19
 Operator : SJ/JU
 Sample : PB107498BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB107498BL

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-BG031918.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	5.790	422	426	434	rBV	28672	49301	1.28%	0.287%
2	6.307	507	514	535	rBV	607881	1254432	32.60%	7.311%
3	7.994	794	801	824	rBV	628229	1416723	36.82%	8.257%
4	8.399	863	870	890	rBV	603335	1300133	33.79%	7.577%
5	8.893	946	954	968	rBV2	90472	220216	5.72%	1.283%
6	9.228	1003	1011	1025	rBV	459612	939107	24.41%	5.473%
7	10.091	1151	1158	1172	rBV	362774	815778	21.20%	4.754%
8	11.778	1437	1445	1461	rBV2	146767	329579	8.57%	1.921%
9	14.139	1840	1847	1869	rBV	1260016	2149496	55.87%	12.528%
10	15.508	2072	2080	2094	rBV2	301752	539712	14.03%	3.146%
11	16.983	2324	2331	2347	rBV	1087585	1851897	48.13%	10.793%
12	18.241	2538	2545	2559	rBV	441526	756217	19.65%	4.407%
13	20.761	2967	2974	2986	rBV	2821117	3847553	100.00%	22.424%
14	22.594	3279	3286	3298	rBV2	451600	866323	22.52%	5.049%
15	26.390	3921	3932	3947	rVB3	241668	821571	21.35%	4.788%

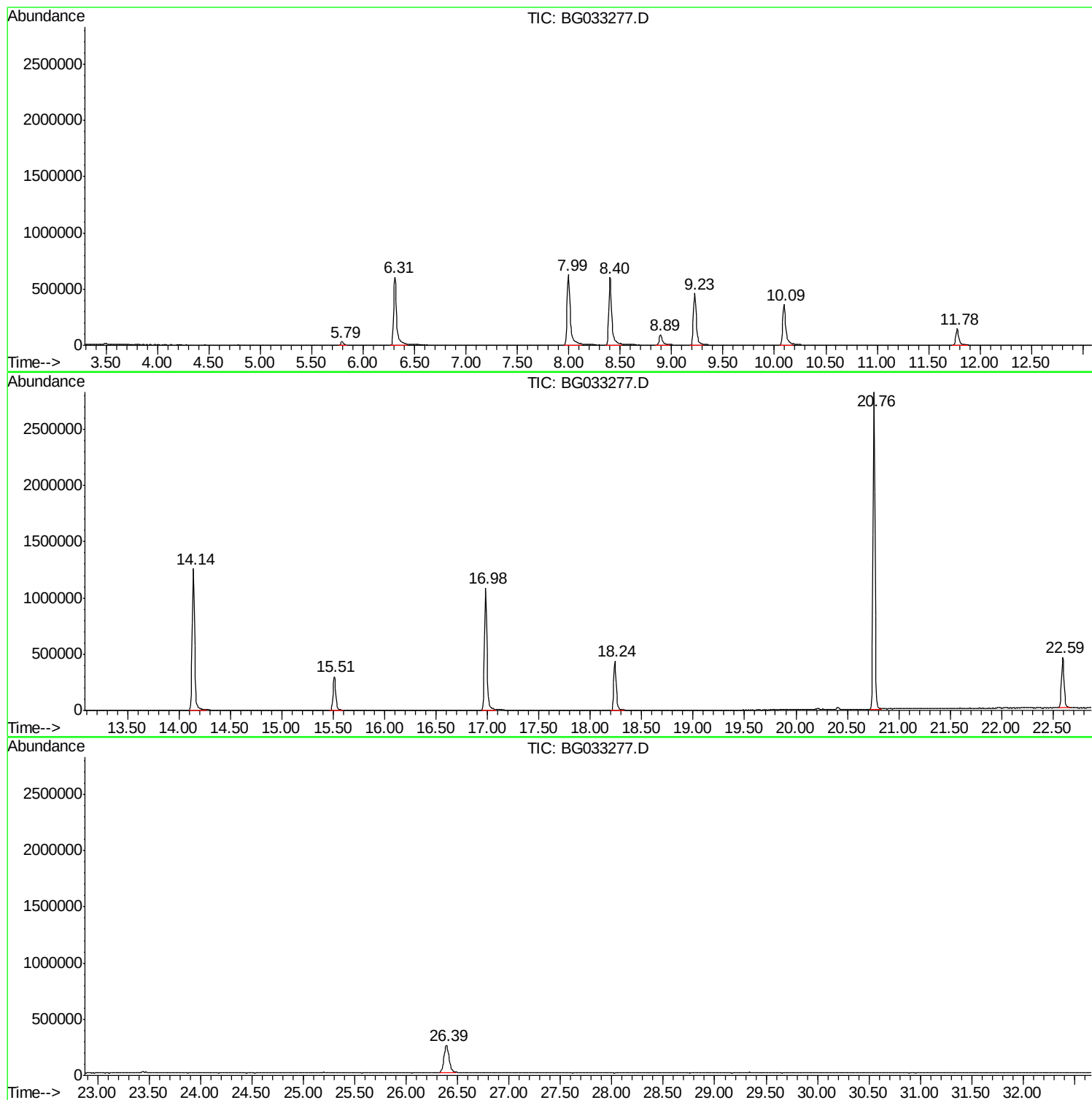
Sum of corrected areas: 17158038

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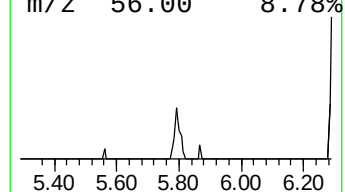
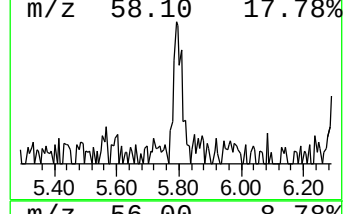
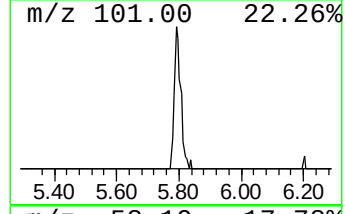
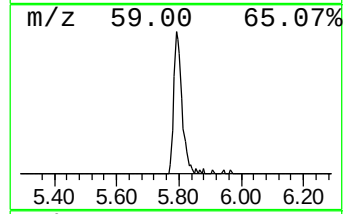
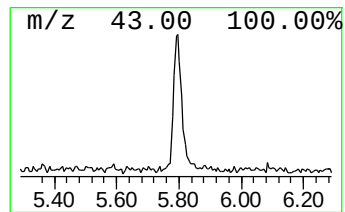
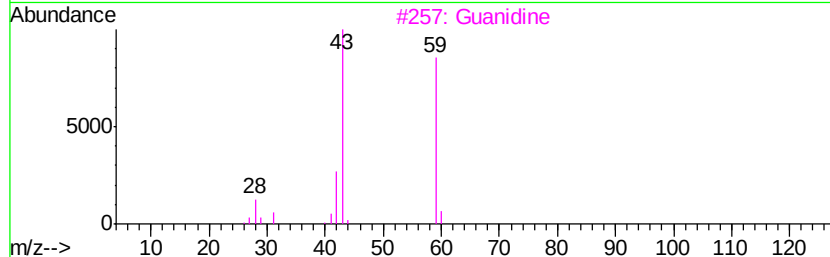
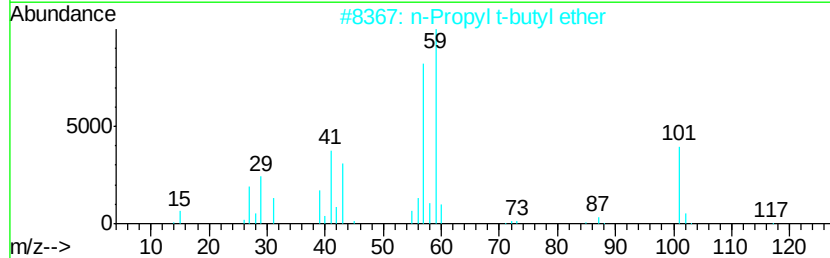
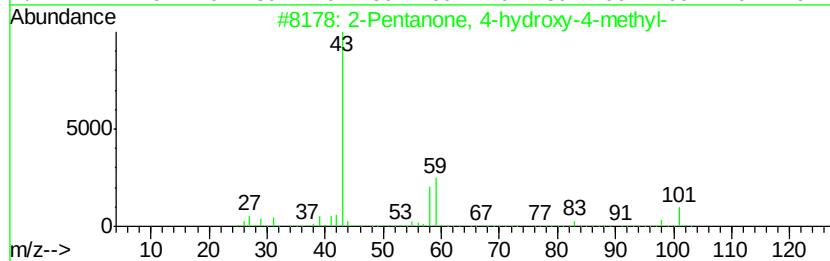
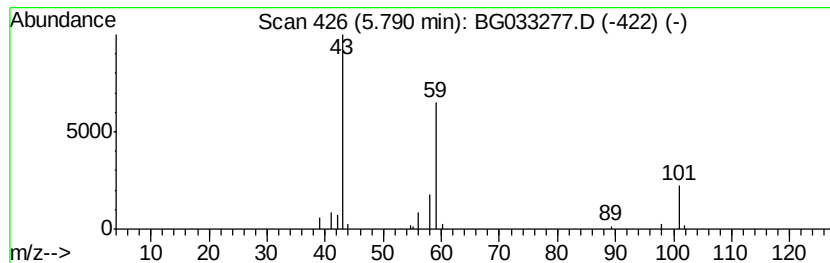
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.79	4.48 ng	49301	1,4-Dichlorobenzene-d4	8.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	33
3			Guanidine	59	CH5N3	000113-00-8	9
4			Acetone	58	C3H6O	000067-64-1	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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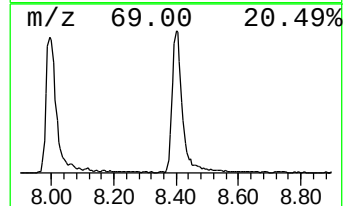
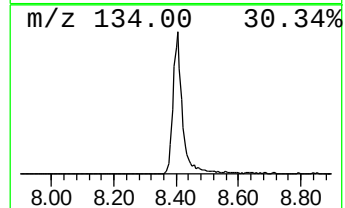
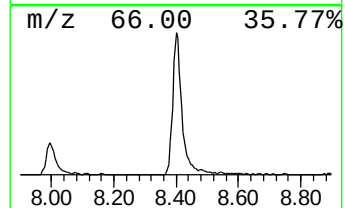
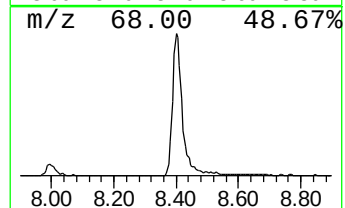
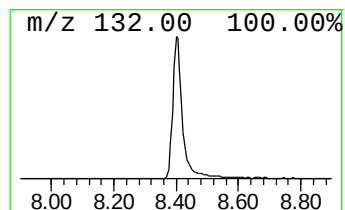
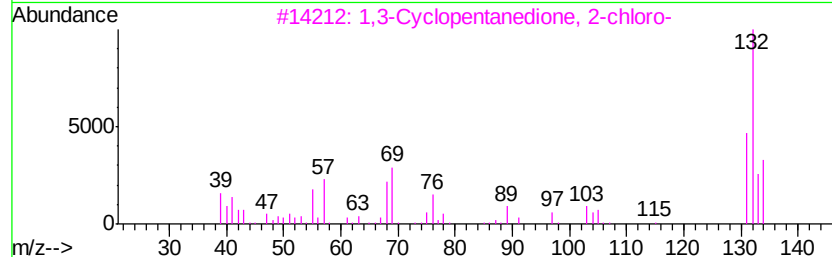
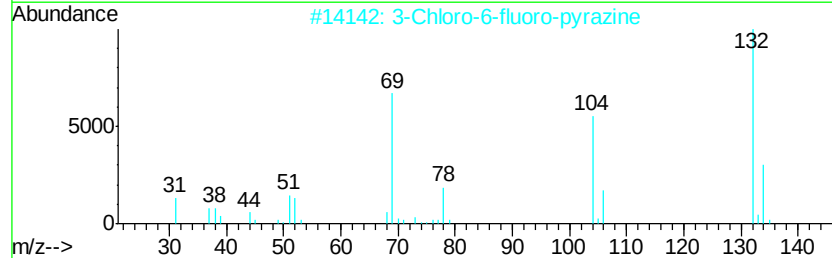
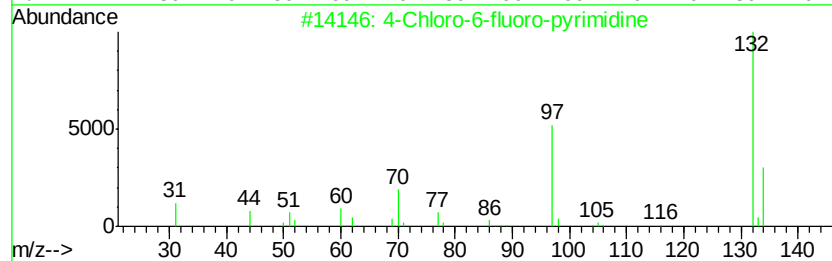
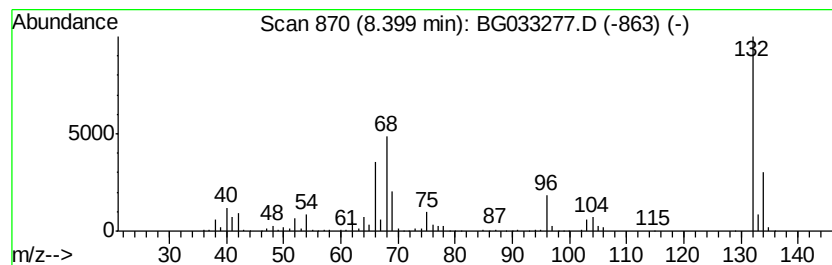
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Peak Number 2 unknown8.40 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.40	118.08 ng	1300130	1,4-Dichlorobenzene-d4	8.89

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
4		1H-Tetrazaborole, 5-chloro-4,5-d...	132	C2H6BClN4	021960-49-6	16
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.79	4.5	ng	49301	1	8.89	220216	20.0
unknown8.40	8.40	118.1	ng	1300130	1	8.89	220216	20.0