

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102817\
 Data File : BG030532.D
 Acq On : 28 Oct 2017 18:02
 Operator : SJ/JU
 Sample : PB103686TB
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103686TB

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-BG101617.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	5.239	326	332	343	rBV	199901	308933	5.83%	1.203%
2	5.720	406	414	427	rVB	1145611	1854928	35.03%	7.223%
3	7.318	678	686	696	rBV	1132996	1941807	36.67%	7.561%
4	7.694	739	750	763	rBV	1117747	1936400	36.56%	7.540%
5	8.164	822	830	843	rBV	243738	409346	7.73%	1.594%
6	8.494	877	886	894	rBV	844156	1510068	28.51%	5.880%
7	9.328	1020	1028	1038	rBV	629413	1152653	21.77%	4.488%
8	10.985	1300	1310	1319	rVB	318279	579635	10.95%	2.257%
9	13.411	1715	1723	1733	rBV	1913629	3091477	58.38%	12.038%
10	14.780	1949	1956	1966	rVB2	534163	866965	16.37%	3.376%
11	16.267	2202	2209	2228	rBV	1610954	2644586	49.94%	10.298%
12	17.518	2415	2422	2434	rBV	780071	1212870	22.90%	4.723%
13	20.109	2857	2863	2872	rBV	4058344	5295854	100.00%	20.621%
14	21.790	3142	3149	3157	rBV	860975	1460673	27.58%	5.688%
15	25.103	3701	3713	3726	rBV	475592	1415602	26.73%	5.512%

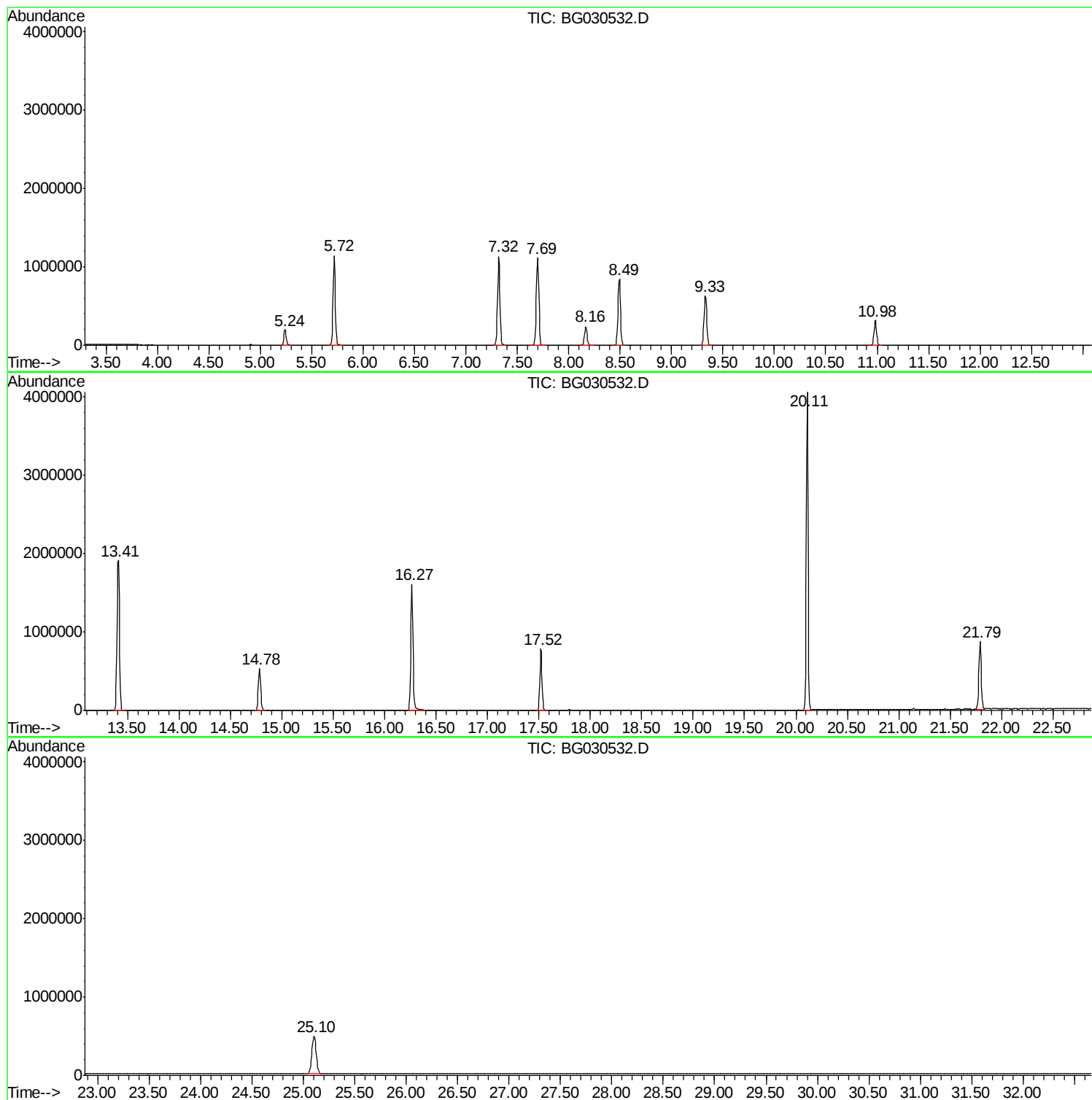
Sum of corrected areas: 25681797

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.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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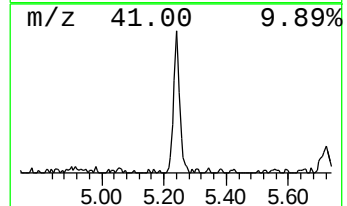
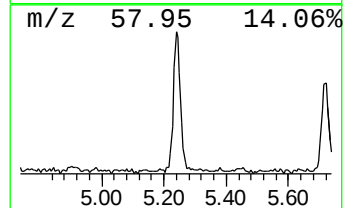
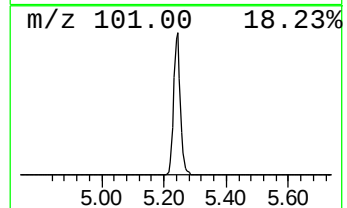
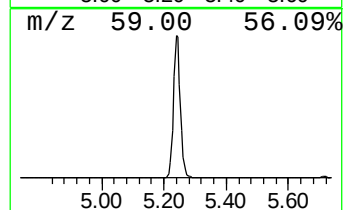
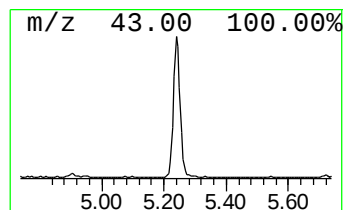
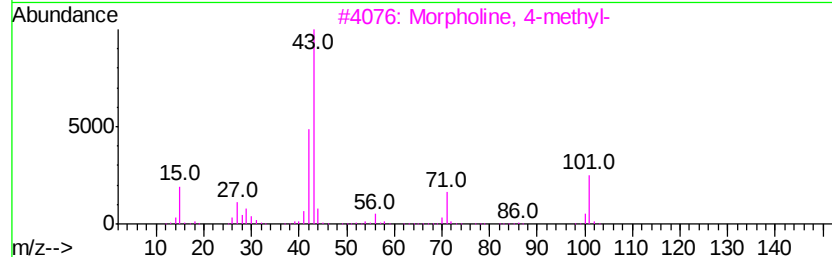
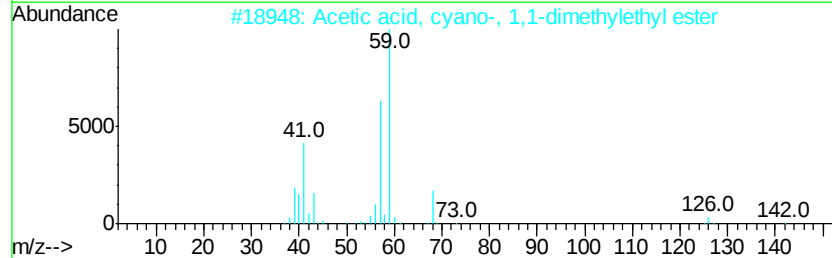
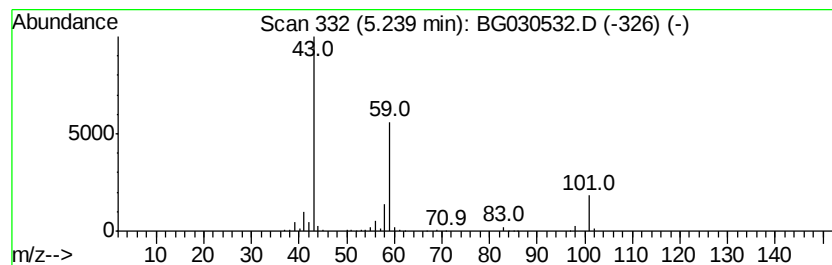
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	15.09 ng	308933	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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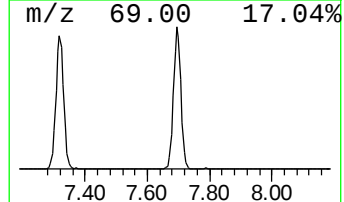
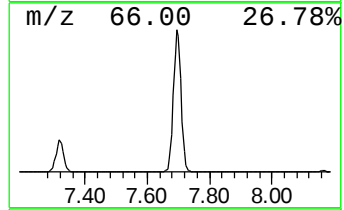
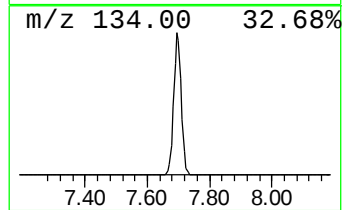
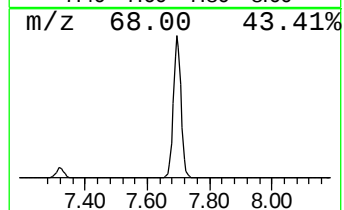
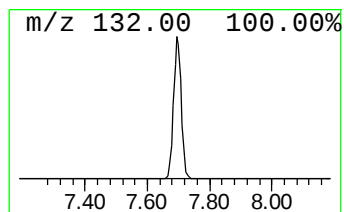
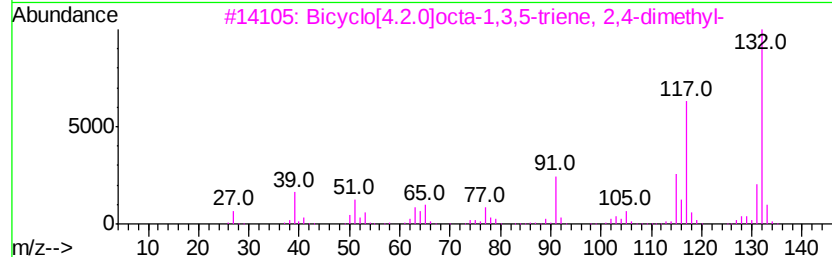
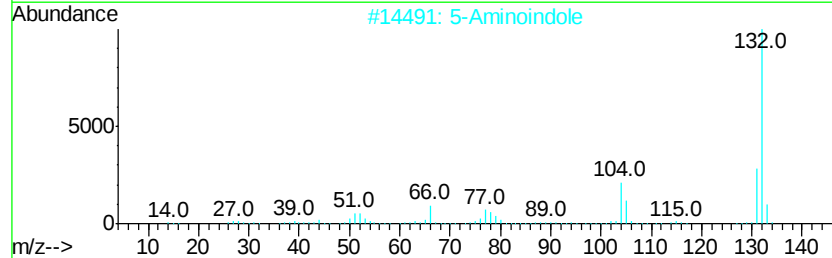
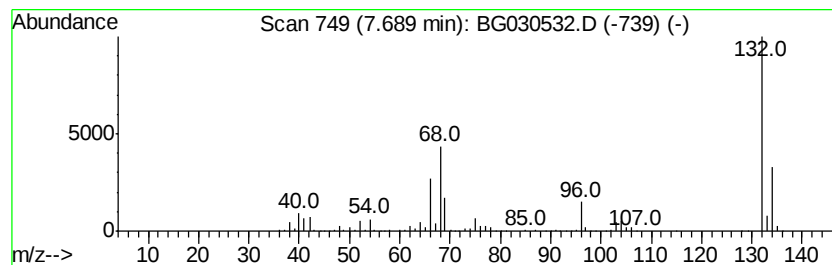
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Peak Number 2 unknown7.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	94.61 ng	1936400	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	15.1	ng	308933	1	8.16	409346	20.0
unknown7.69	7.69	94.6	ng	1936400	1	8.16	409346	20.0