

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101817\
 Data File : BG029353.D
 Acq On : 19 Oct 2017 5:08
 Operator : SJ/JU
 Sample : PB103395BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103395BL

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	5.238	327	332	343	rVB	145048	222401	5.15%	1.108%
2	5.720	407	414	423	rBV	832250	1340547	31.05%	6.681%
3	7.318	678	686	695	rBV	841934	1401740	32.46%	6.986%
4	7.694	743	750	758	rBV	807008	1391666	32.23%	6.936%
5	8.164	823	830	838	rVB	190281	324179	7.51%	1.616%
6	8.493	878	886	893	rBV	625629	1071640	24.82%	5.341%
7	9.328	1020	1028	1039	rBV	467622	827742	19.17%	4.125%
8	10.984	1301	1310	1317	rBV	270192	478159	11.07%	2.383%
9	13.411	1715	1723	1731	rBV	1463061	2204037	51.05%	10.985%
10	14.786	1948	1957	1967	rBV2	438374	718285	16.64%	3.580%
11	16.266	2201	2209	2222	rBV	1290076	2019208	46.77%	10.064%
12	17.518	2415	2422	2434	rVB2	649710	1029185	23.84%	5.129%
13	20.115	2857	2864	2869	rBV	3223546	4317702	100.00%	21.519%
14	21.789	3142	3149	3162	rBV	842202	1365297	31.62%	6.805%
15	25.097	3701	3712	3725	rBV2	462454	1352737	31.33%	6.742%

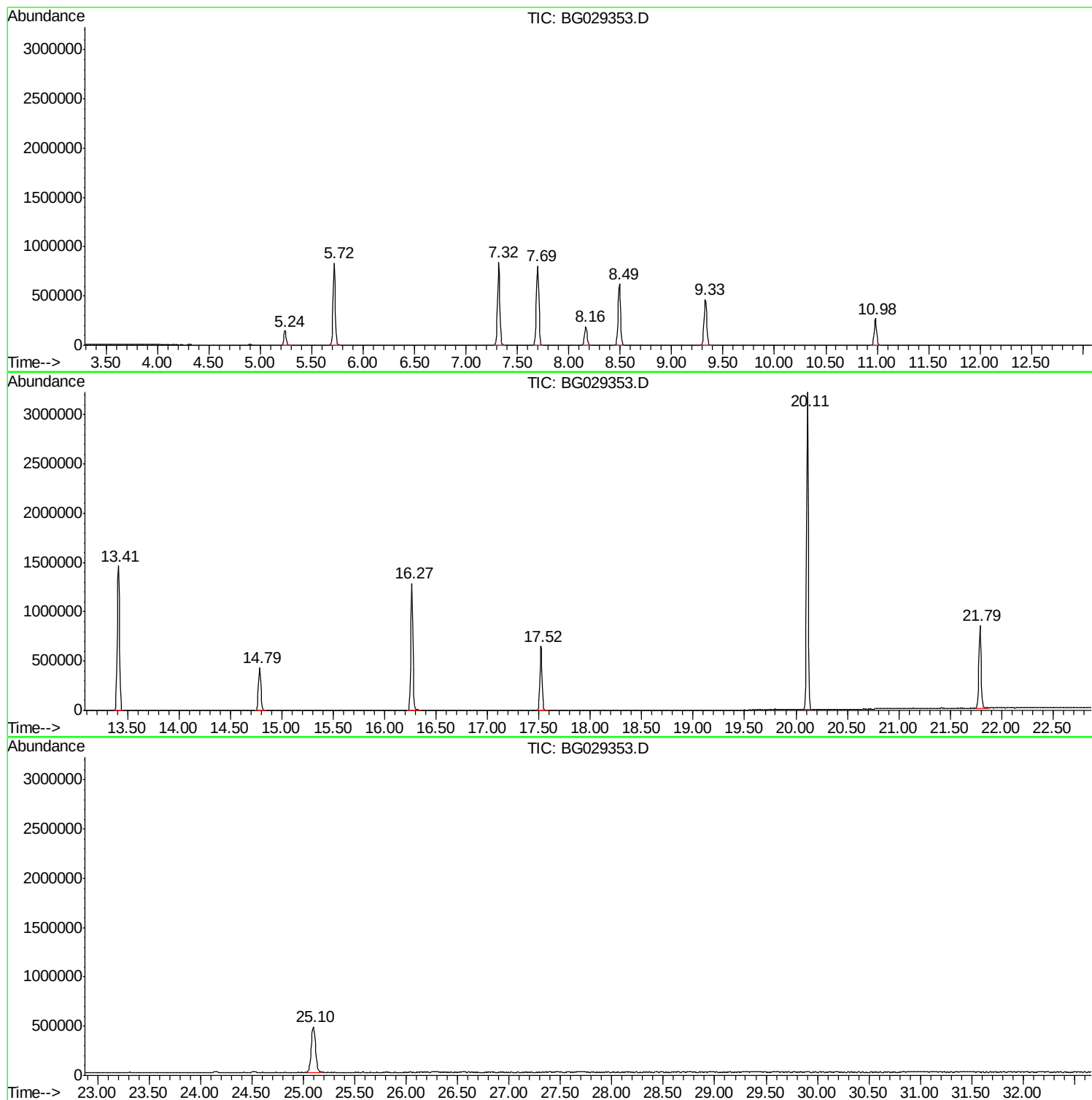
Sum of corrected areas: 20064525

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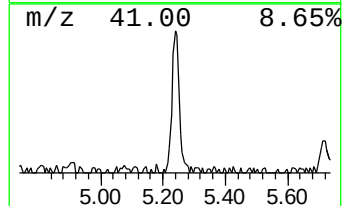
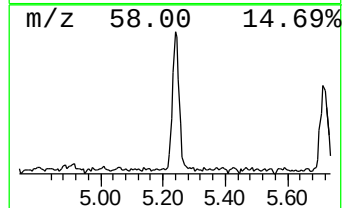
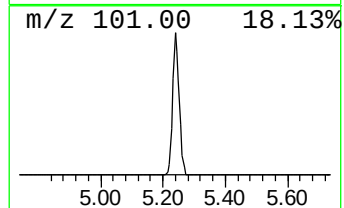
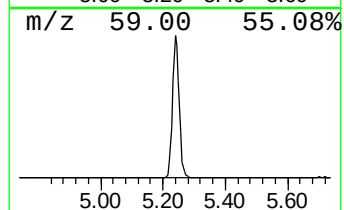
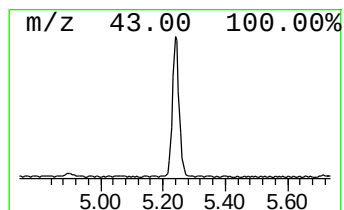
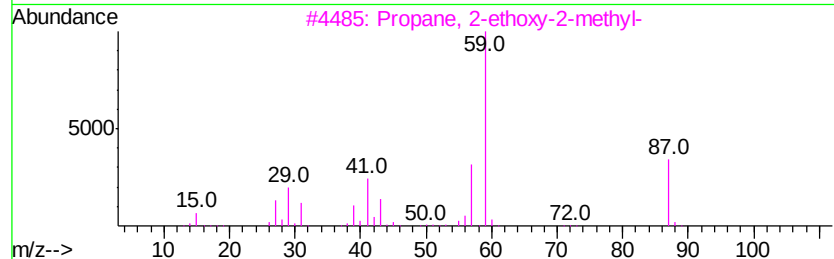
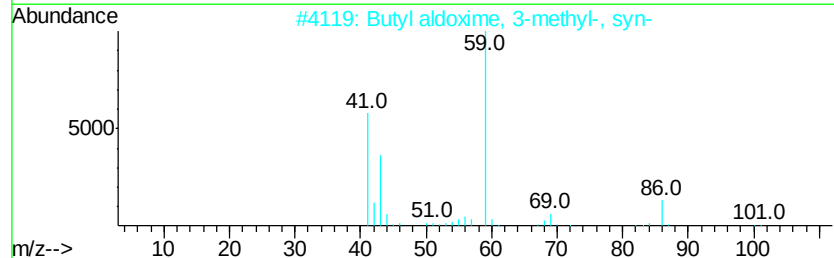
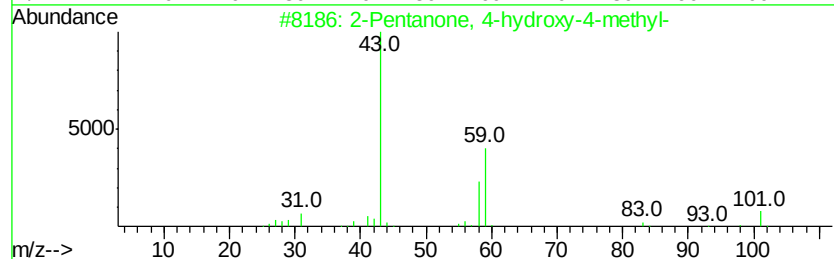
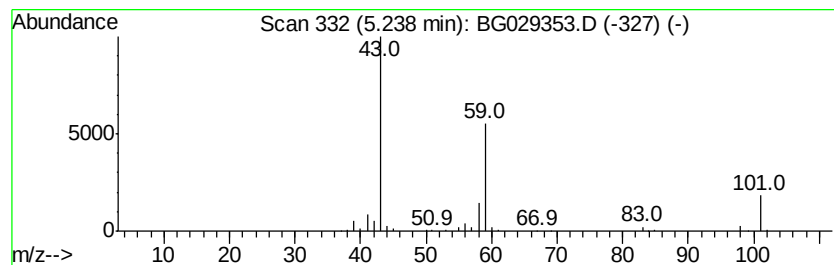
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	13.72 ng	222401	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	45
2		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
3		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	23
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9



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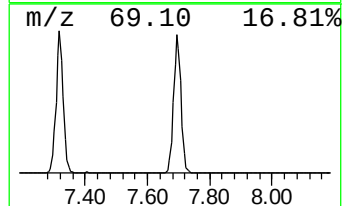
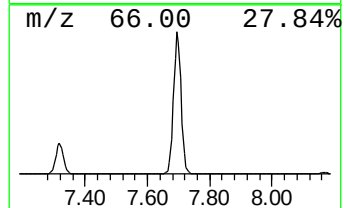
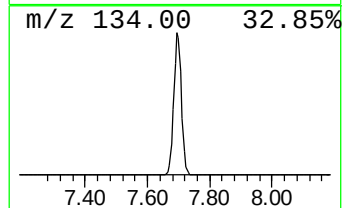
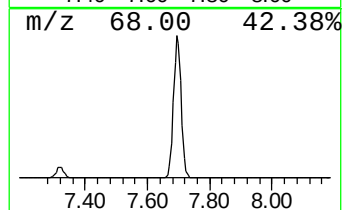
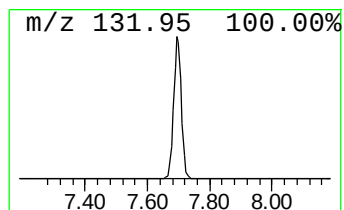
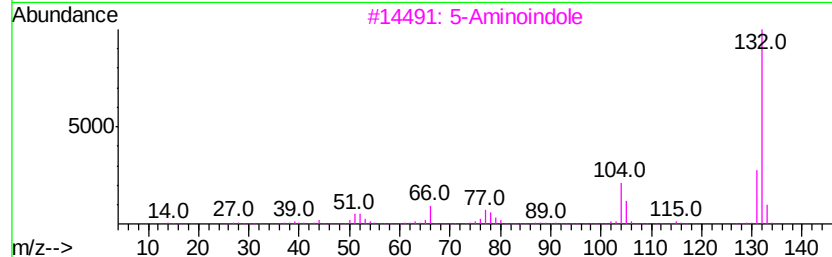
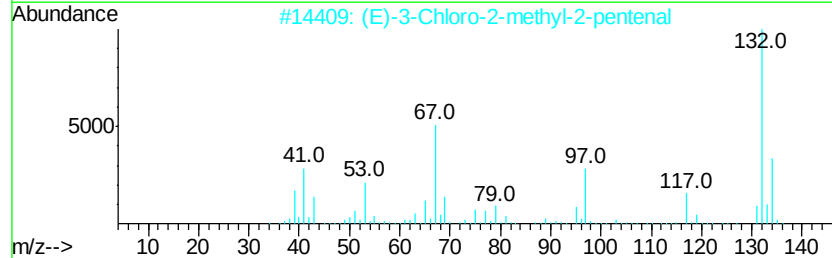
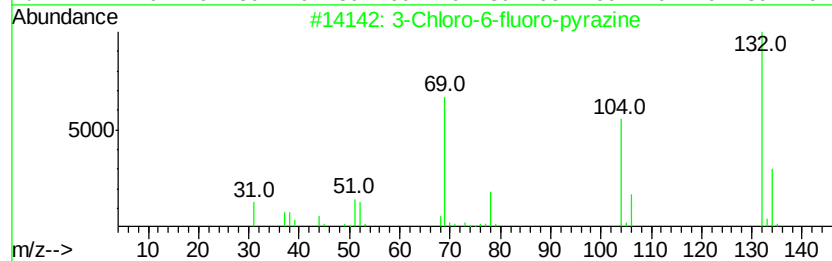
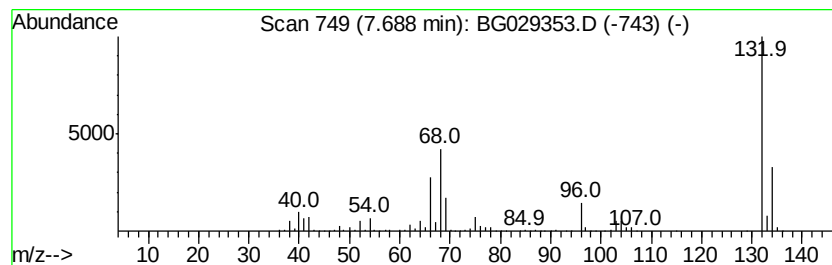
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Peak Number 2 unknown7.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	85.86 ng	1391670	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		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	13.7	ng	222401	1	8.16	324179	20.0
unknown7.69	7.69	85.9	ng	1391670	1	8.16	324179	20.0