

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121818\
 Data File : BG038699.D
 Acq On : 18 Dec 2018 18:00
 Operator : JU/SJ
 Sample : PB115750BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115750BL

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:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.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	5.138	345	349	359	rBV	17430	28687	1.65%	0.350%
2	5.608	422	429	439	rBV	349278	579424	33.41%	7.063%
3	7.212	694	702	714	rBV	360419	643476	37.10%	7.844%
4	7.582	758	765	773	rBV	335269	606187	34.95%	7.389%
5	8.058	839	846	854	rVB	84982	146508	8.45%	1.786%
6	8.382	892	901	909	rBV	256454	475747	27.43%	5.799%
7	9.233	1038	1046	1055	rBV	211774	393058	22.66%	4.791%
8	10.873	1317	1325	1336	rBV	108102	202446	11.67%	2.468%
9	13.311	1733	1740	1755	rBV	598609	965660	55.67%	11.771%
10	14.698	1969	1976	1990	rVB2	170053	301449	17.38%	3.675%
11	16.184	2223	2229	2244	rBV2	468076	762383	43.95%	9.293%
12	17.441	2436	2443	2452	rBV2	257506	414294	23.89%	5.050%
13	20.044	2879	2886	2897	rBV	1304360	1734490	100.00%	21.143%
14	21.719	3163	3171	3181	rBV	296589	490194	28.26%	5.975%
15	25.015	3722	3732	3751	rVB2	148041	459709	26.50%	5.604%

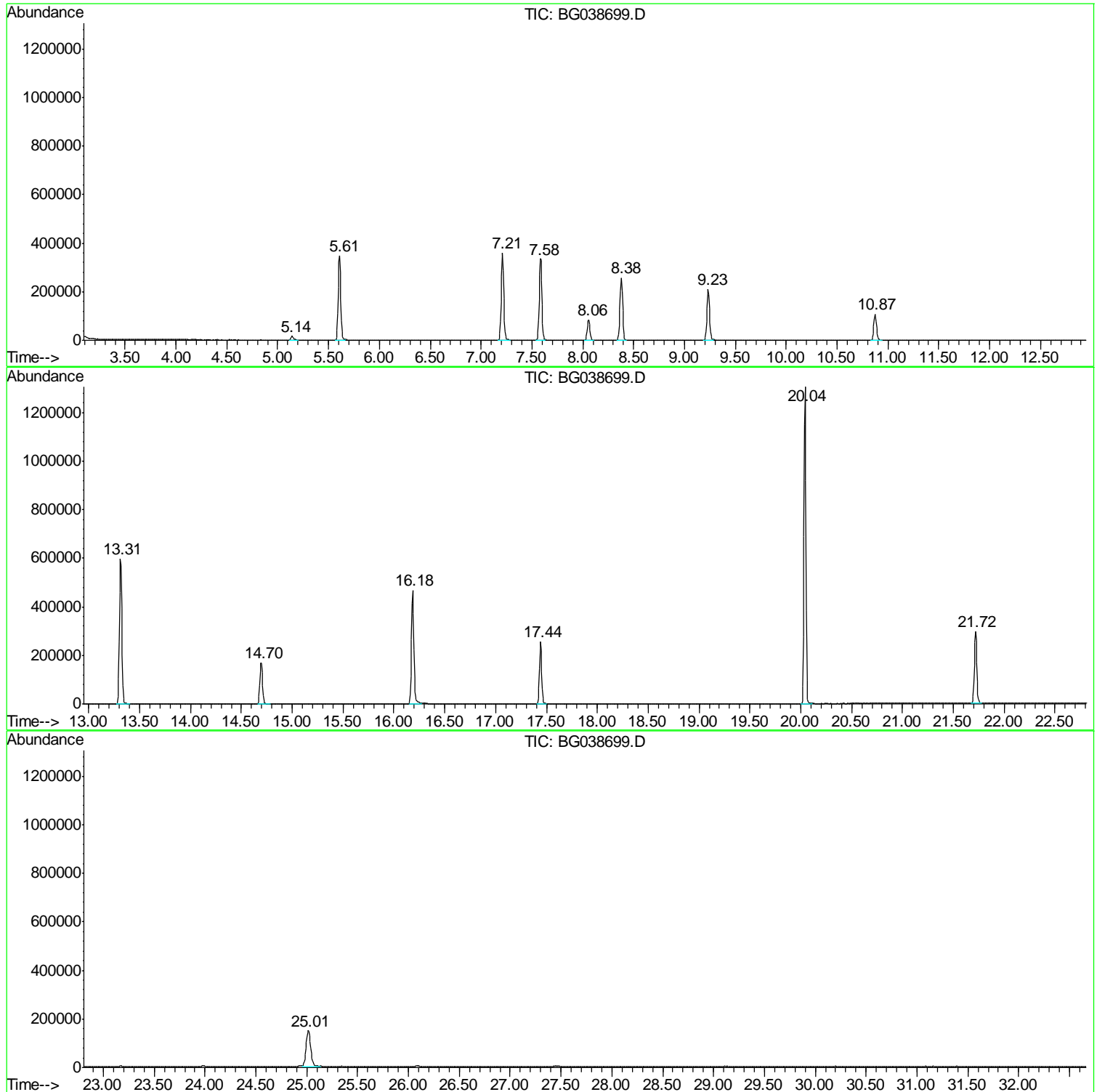
Sum of corrected areas: 8203712

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.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 :
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 ClientSampled :
 PB115750BL

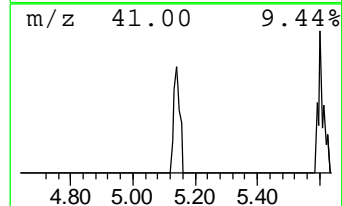
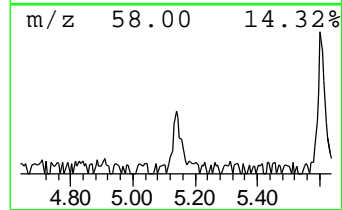
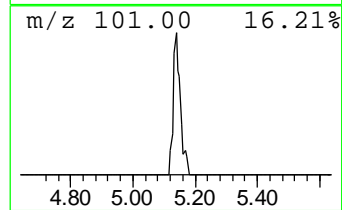
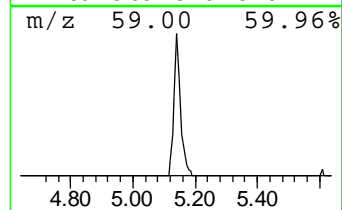
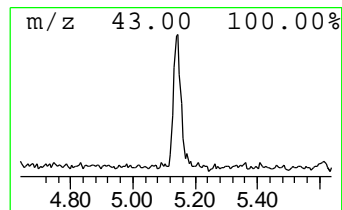
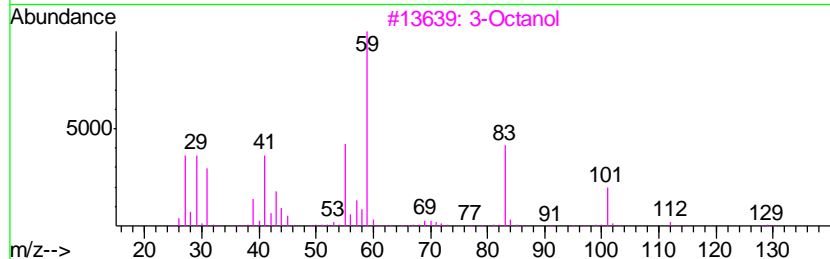
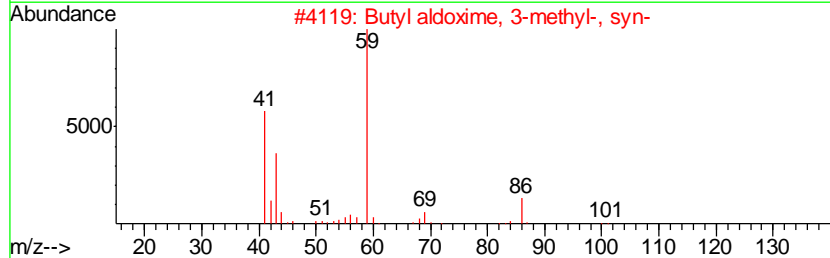
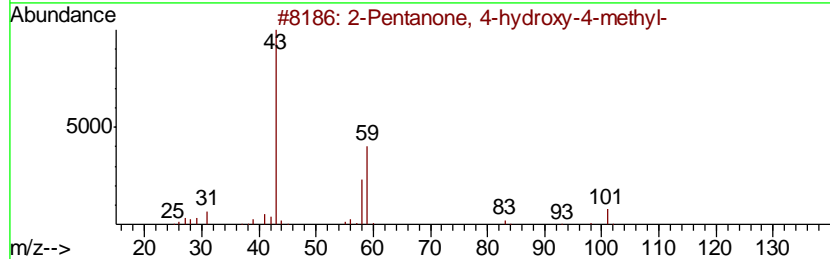
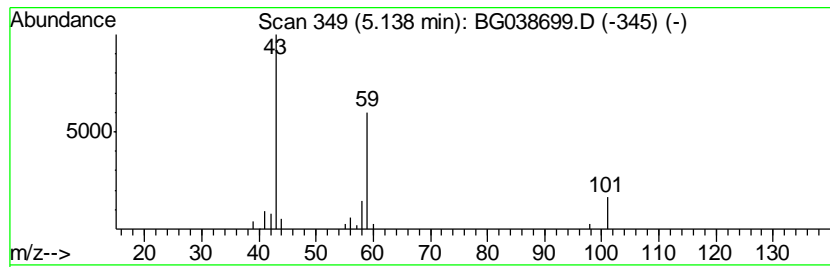
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 Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	3.92 ng	28687	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	37
3		3-Octanol	130	C8H18O	000589-98-0	33
4		Butane, 1-(1-methylethoxy)-	116	C7H16O	001860-27-1	32
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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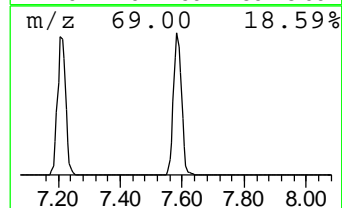
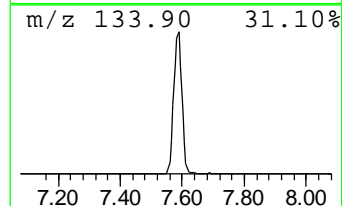
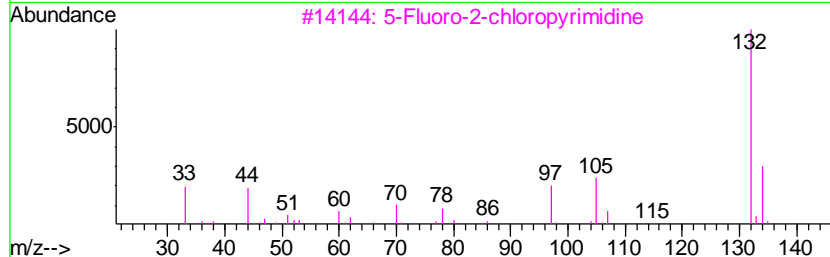
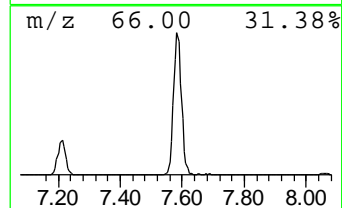
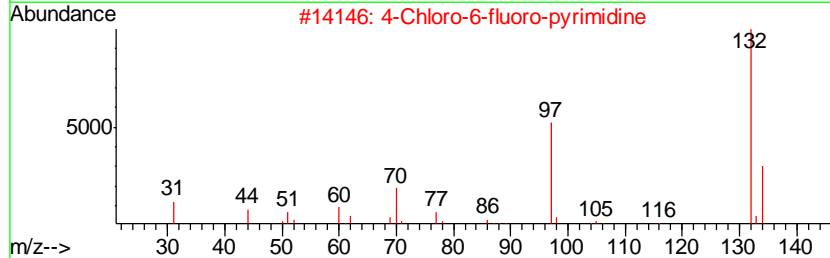
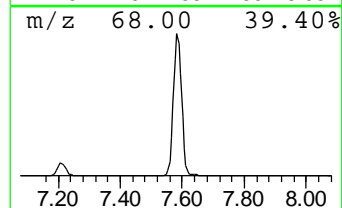
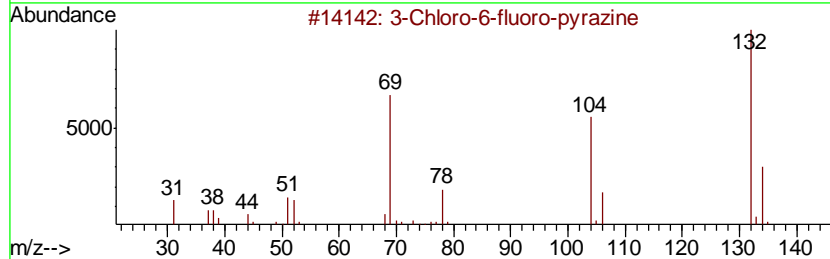
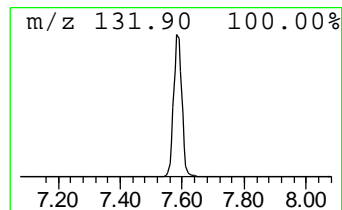
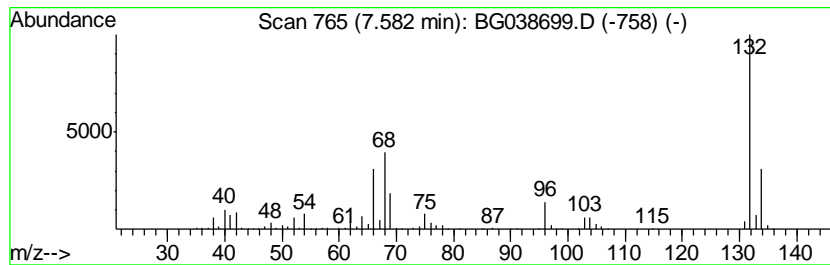
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 Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	82.75 ng	606187	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	12
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	3.9	ng	28687	1	8.06	146508	20.0
unknown7.58	7.58	82.8	ng	606187	1	8.06	146508	20.0