

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122618\
 Data File : BG038836.D
 Acq On : 26 Dec 2018 18:50
 Operator : JU/SJ
 Sample : PB115917BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115917BL

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	342	349	357	rBV	23371	38354	1.53%	0.384%
2	5.608	422	429	441	rBV	402048	661014	26.39%	6.615%
3	7.218	694	703	713	rBV	429316	752418	30.04%	7.530%
4	7.588	758	766	778	rBV	397678	705123	28.15%	7.057%
5	8.053	836	845	853	rVB	69516	120248	4.80%	1.203%
6	8.376	892	900	910	rVB	290426	526731	21.03%	5.271%
7	9.234	1038	1046	1057	rBV	234838	443340	17.70%	4.437%
8	10.873	1316	1325	1333	rBV	90668	167903	6.70%	1.680%
9	13.311	1733	1740	1754	rBV	731881	1209468	48.29%	12.104%
10	14.692	1969	1975	1985	rBV2	162619	274361	10.95%	2.746%
11	16.184	2222	2229	2247	rBV3	720806	1178605	47.06%	11.795%
12	17.442	2436	2443	2451	rBV	263138	412271	16.46%	4.126%
13	20.044	2879	2886	2893	rBV	1892677	2504575	100.00%	25.065%
14	21.719	3164	3171	3179	rBV2	298596	514135	20.53%	5.145%
15	25.015	3722	3732	3743	rVB	161913	483646	19.31%	4.840%

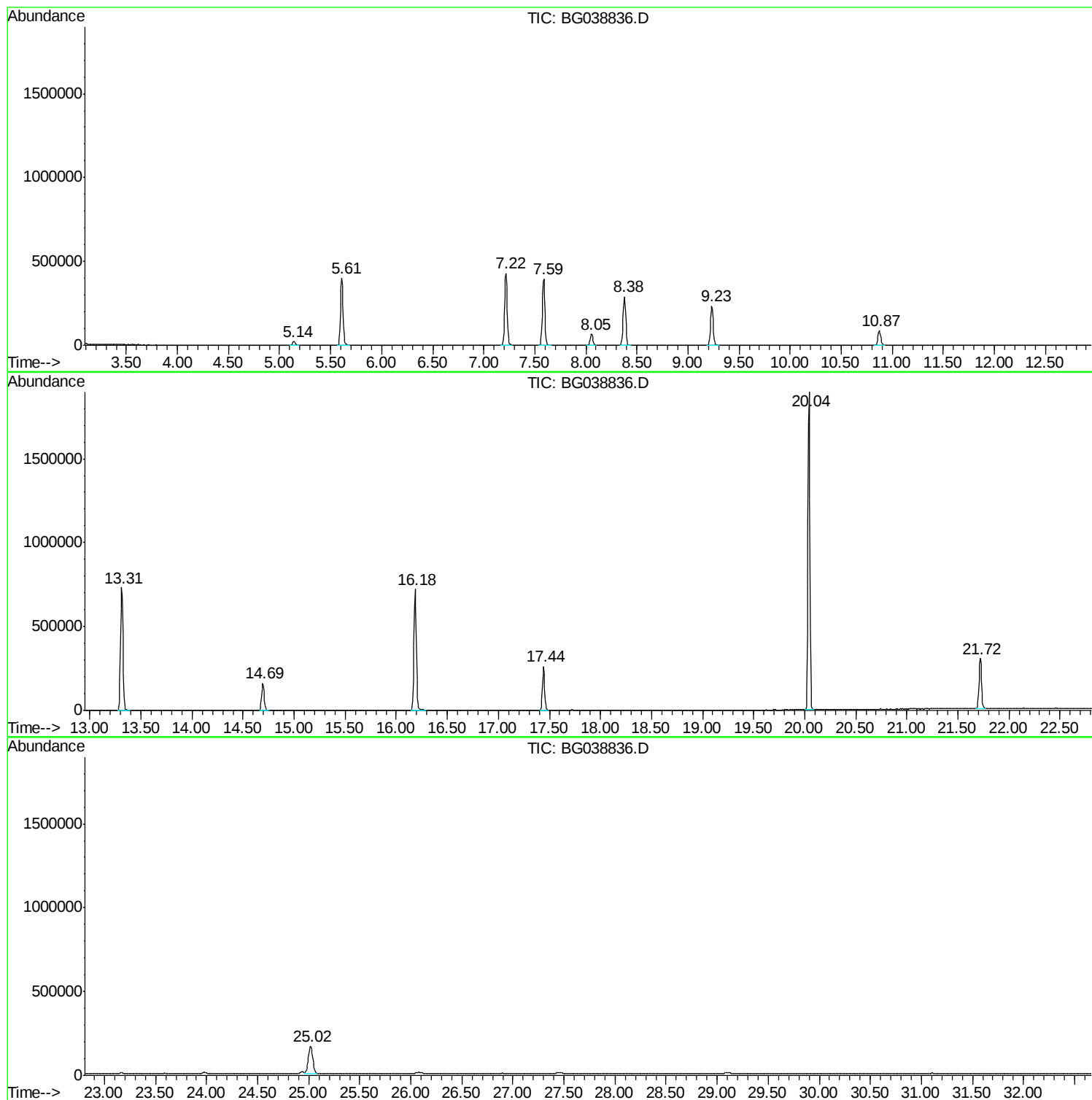
Sum of corrected areas: 9992192

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



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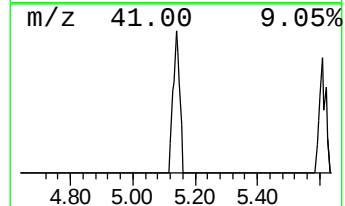
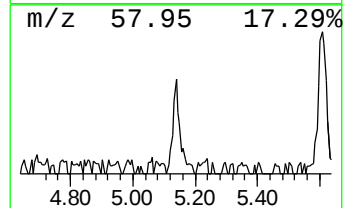
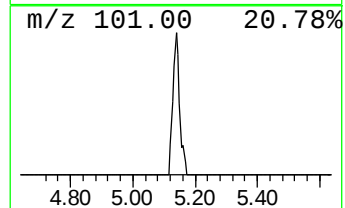
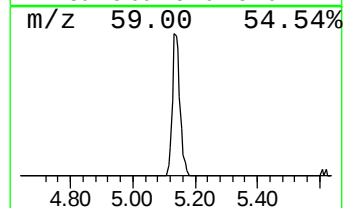
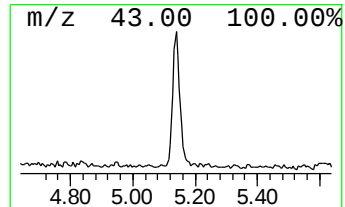
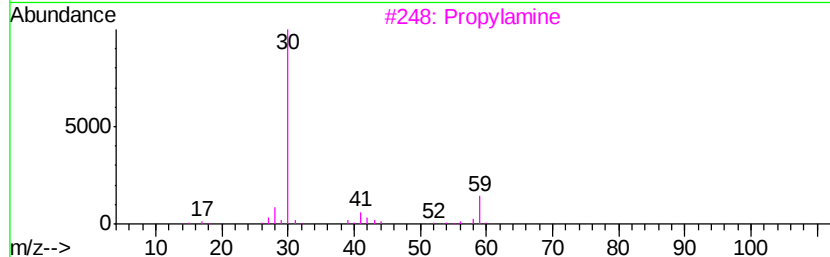
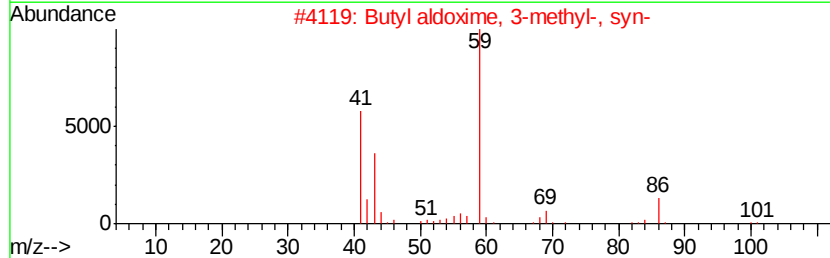
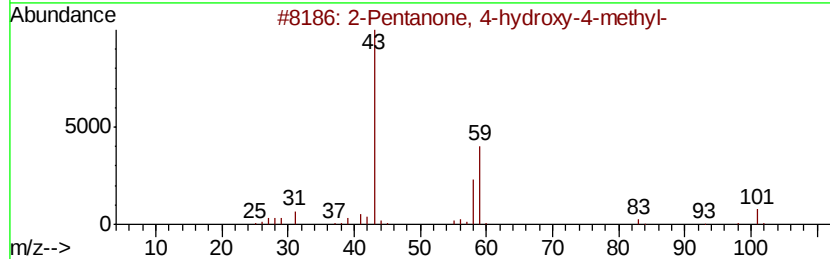
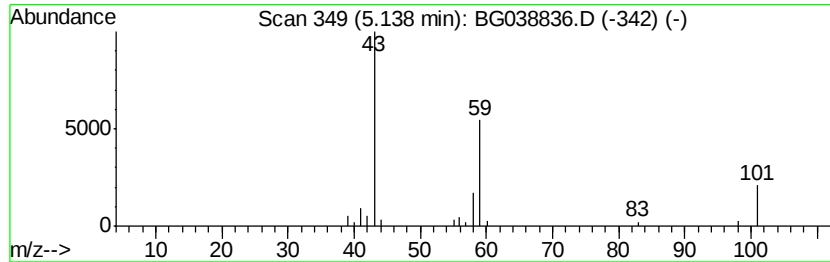
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	6.38 ng	38354	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
3			Propylamine	59	C3H9N	000107-10-8	9
4			Tetramethylammonium acetate	133	C6H15N02	010581-12-1	9
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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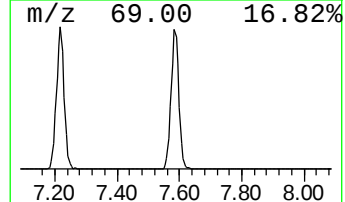
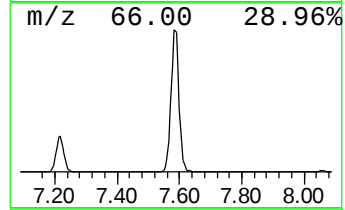
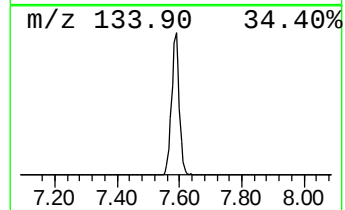
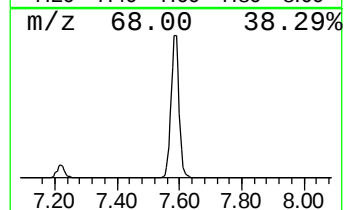
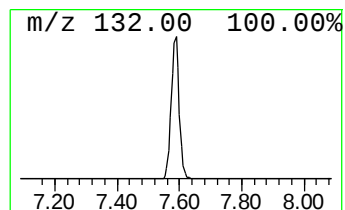
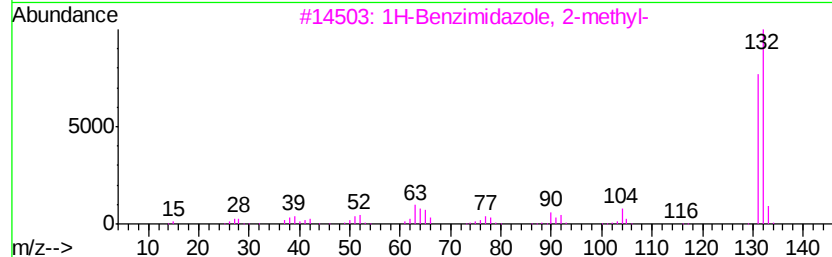
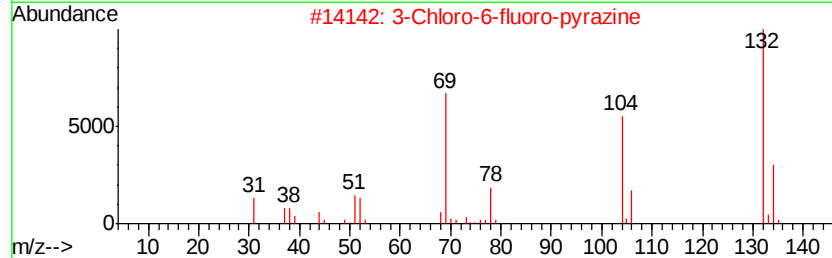
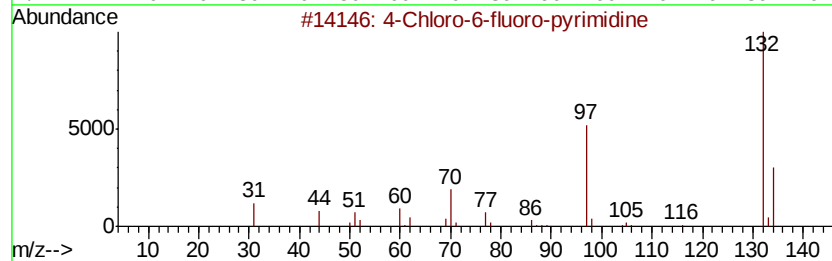
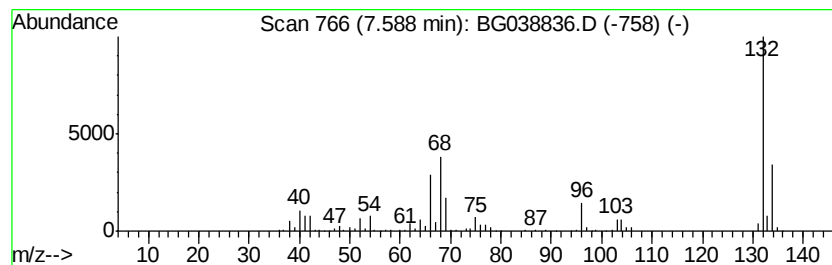
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	117.28 ng	705123	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	25
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	6.4	ng	38354	1	8.05	120248	20.0
unknown7.59	7.59	117.3	ng	705123	1	8.05	120248	20.0