

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112018\
 Data File : BG038092.D
 Acq On : 20 Nov 2018 19:27
 Operator : JU/SJ
 Sample : PB114970BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114970BL

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-BG111318.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	5.164	315	319	326	rVB	25956	43779	1.99%	0.368%
2	5.629	391	398	413	rVB	553515	904356	41.07%	7.597%
3	7.233	663	671	682	rBV	571505	1004255	45.60%	8.436%
4	7.609	728	735	747	rBV	524996	951647	43.21%	7.994%
5	8.084	809	816	824	rVB2	132053	231870	10.53%	1.948%
6	8.408	863	871	879	rBV	395487	719797	32.69%	6.047%
7	9.260	1009	1016	1026	rVB	329399	616772	28.01%	5.181%
8	10.905	1289	1296	1305	rVB	187578	339329	15.41%	2.851%
9	13.343	1702	1711	1718	rBV	925555	1464845	66.52%	12.306%
10	14.718	1937	1945	1954	rVB3	287597	487780	22.15%	4.098%
11	16.210	2191	2199	2217	rBV2	599668	997147	45.28%	8.377%
12	17.468	2406	2413	2421	rBV2	402835	610662	27.73%	5.130%
13	20.065	2849	2855	2866	rBV	1606217	2202220	100.00%	18.500%
14	21.751	3135	3142	3154	rVB	401905	681877	30.96%	5.728%
15	25.070	3694	3707	3721	rBV	205343	647470	29.40%	5.439%

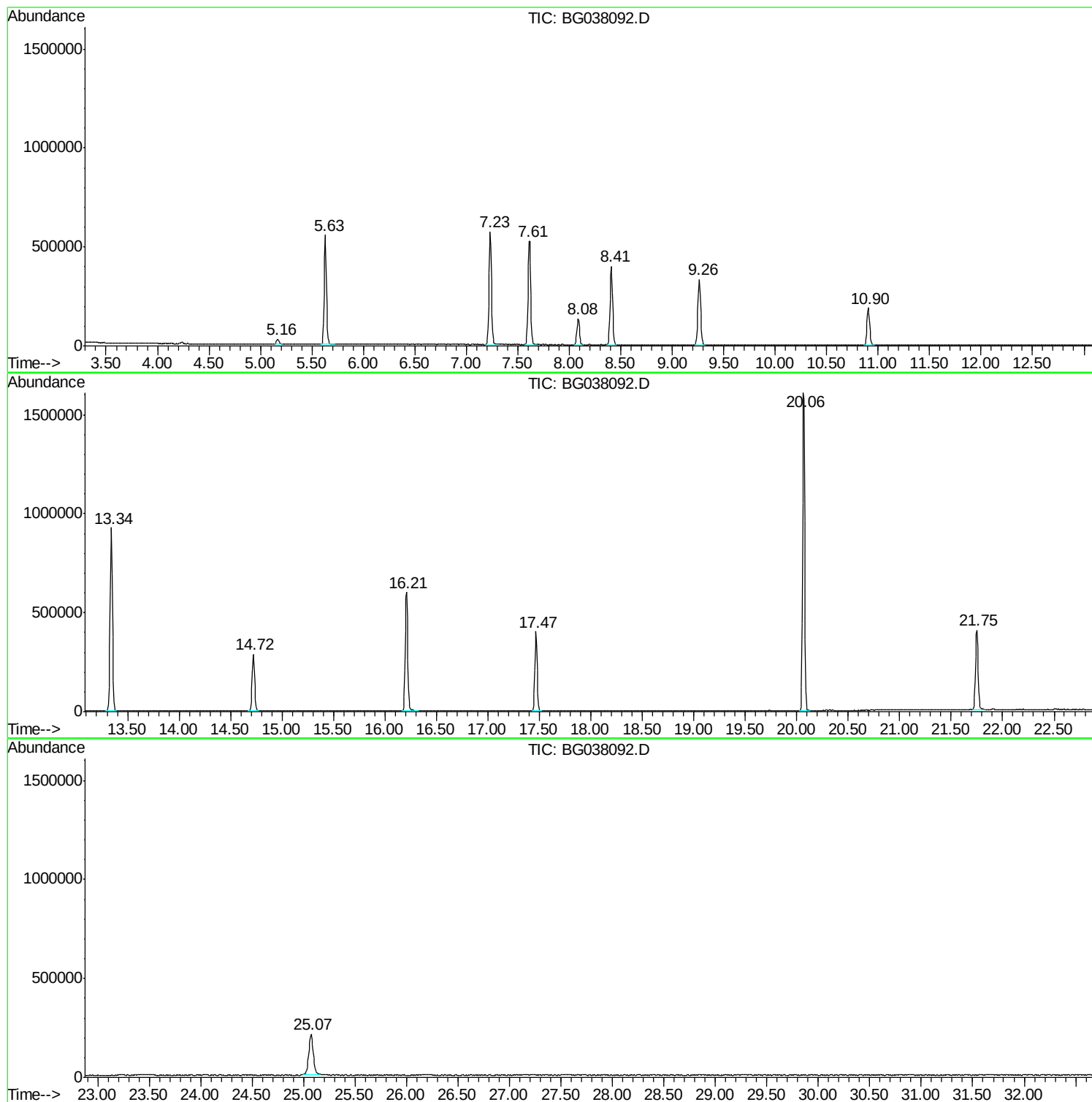
Sum of corrected areas: 11903806

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



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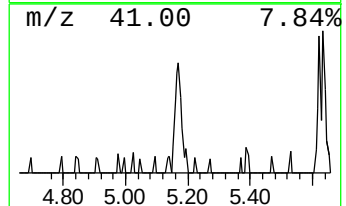
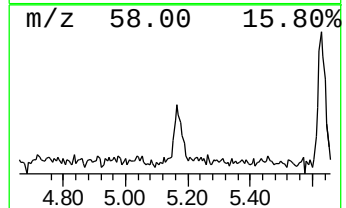
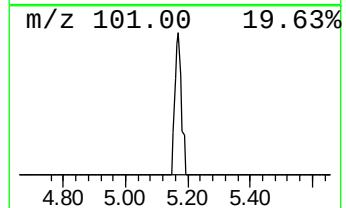
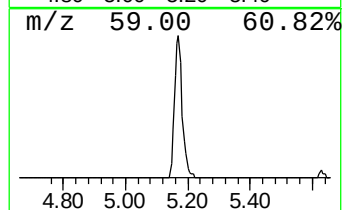
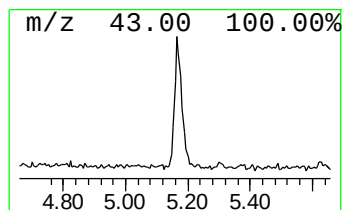
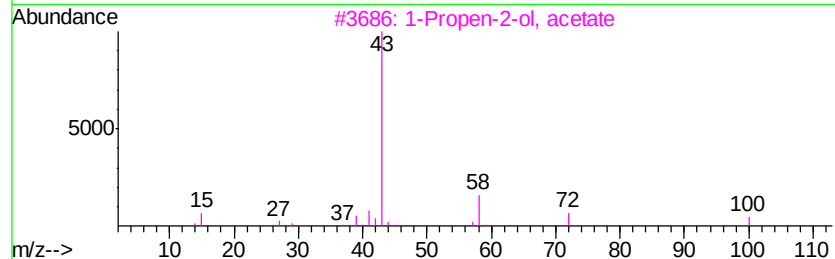
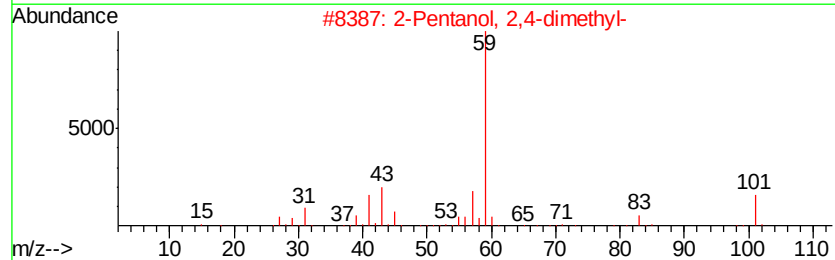
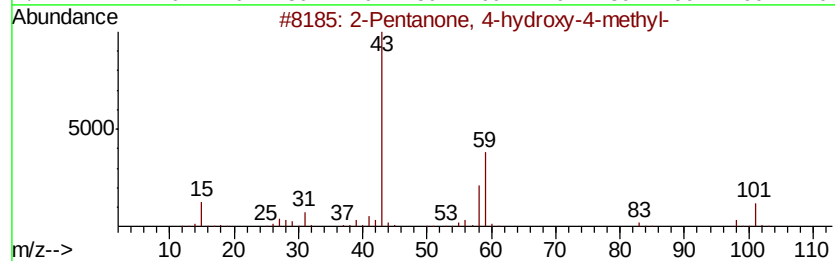
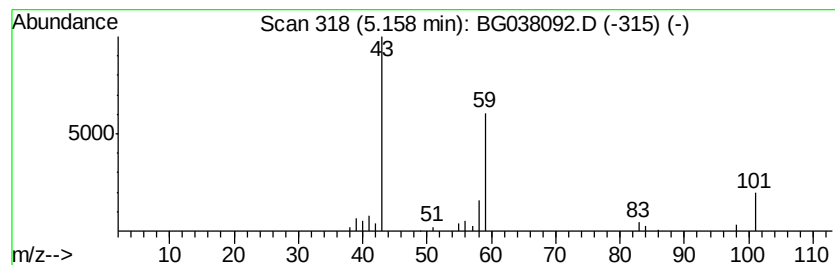
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	3.78 ng	43779	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	9
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4			Diacetamide	101	C4H7NO2	000625-77-4	7
5			5-Hexen-2-one	98	C6H10O	000109-49-9	7



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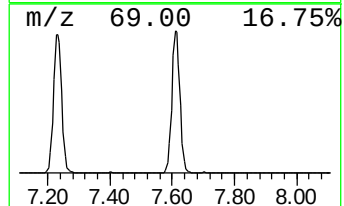
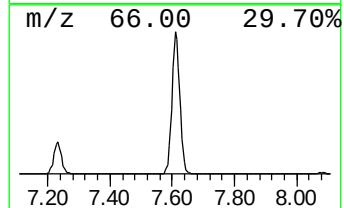
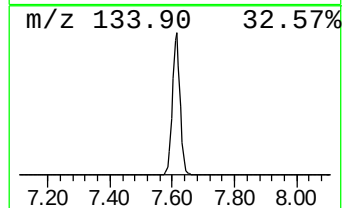
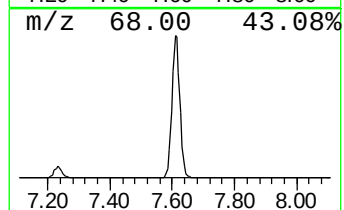
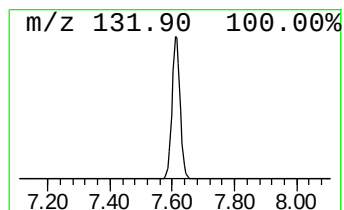
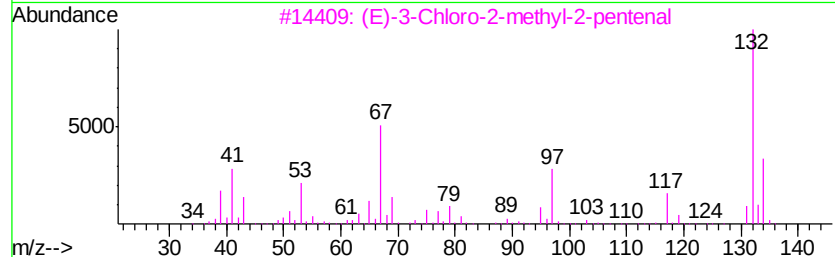
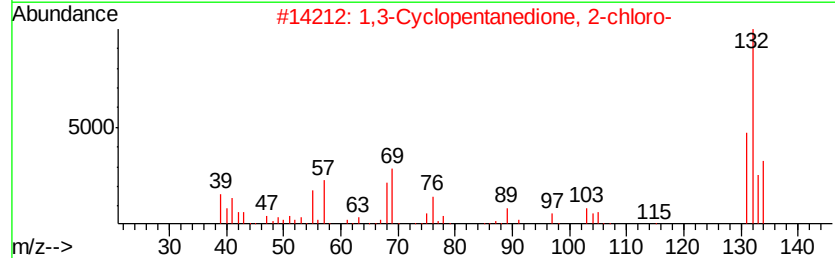
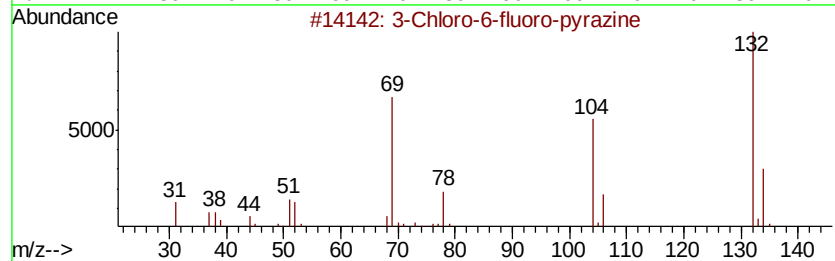
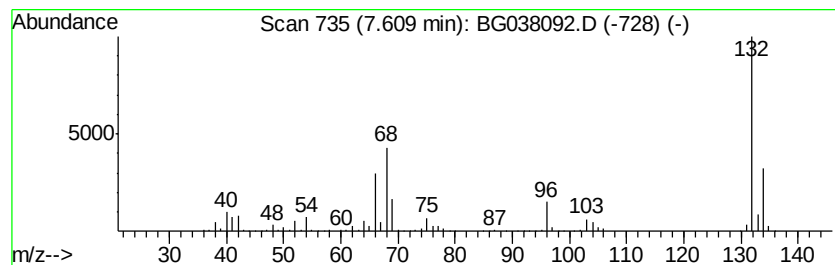
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	82.08 ng	951647	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		5-Aminoindole	132	C8H8N2	005192-03-0	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	3.8	ng	43779	1	8.08	231870	20.0
unknown7.61	7.61	82.1	ng	951647	1	8.08	231870	20.0