

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120118\
 Data File : BG038291.D
 Acq On : 1 Dec 2018 17:26
 Operator : JU/SJ
 Sample : PB115166BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115166BL

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	5.158	313	318	328	rVB2	23769	39625	1.76%	0.385%
2	5.622	390	397	408	rVB	473663	782156	34.81%	7.600%
3	7.226	662	670	682	rBV	482983	852771	37.96%	8.286%
4	7.602	726	734	745	rBV	448634	800526	35.63%	7.779%
5	8.072	808	814	822	rBV	93000	162898	7.25%	1.583%
6	8.395	860	869	878	rBV	333344	618226	27.52%	6.007%
7	9.253	1007	1015	1028	rBV	285108	535011	23.81%	5.199%
8	10.898	1287	1295	1303	rBV	115220	225235	10.03%	2.189%
9	13.331	1702	1709	1723	rBV	747122	1222937	54.43%	11.883%
10	14.711	1936	1944	1953	rBV3	189731	326773	14.55%	3.175%
11	16.198	2191	2197	2211	rBV	598772	953156	42.43%	9.262%
12	17.461	2405	2412	2426	rBV	291191	462423	20.58%	4.493%
13	20.064	2848	2855	2864	rBV	1629146	2246631	100.00%	21.831%
14	21.744	3133	3141	3151	rBV2	296654	534628	23.80%	5.195%
15	25.052	3685	3704	3721	rBV2	160192	528175	23.51%	5.132%

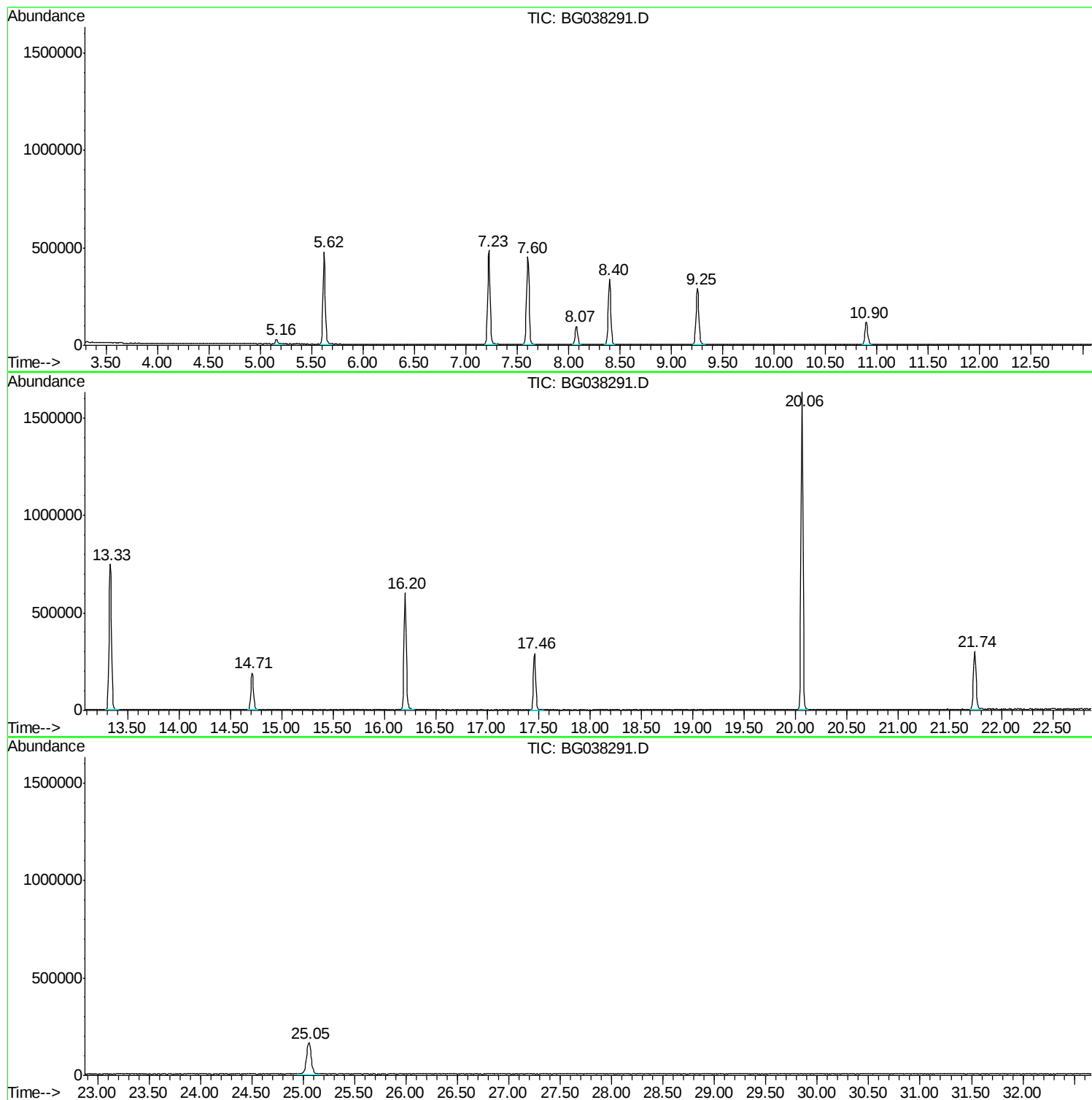
Sum of corrected areas: 10291171

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



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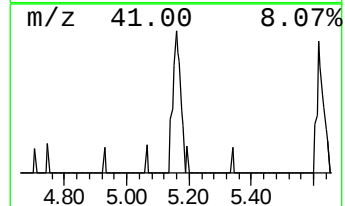
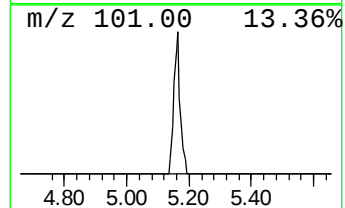
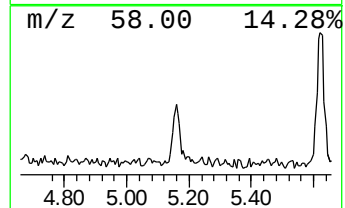
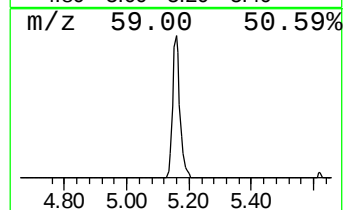
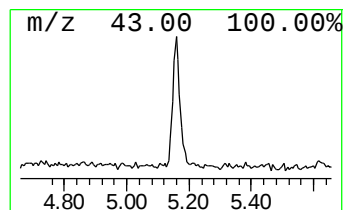
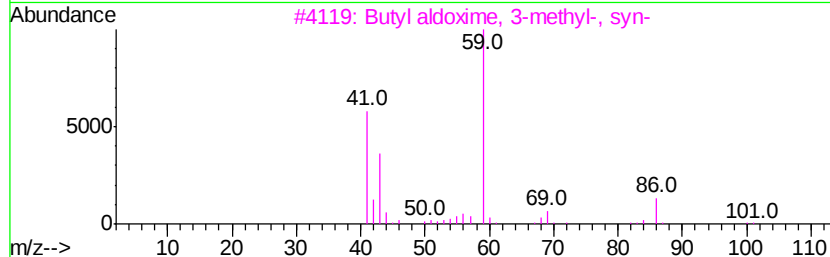
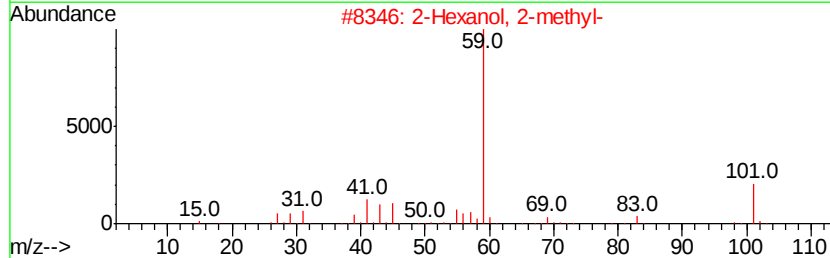
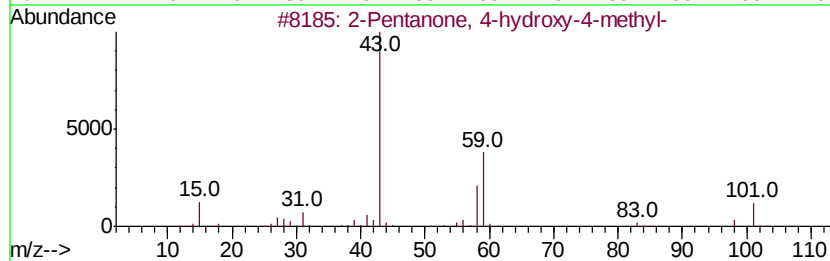
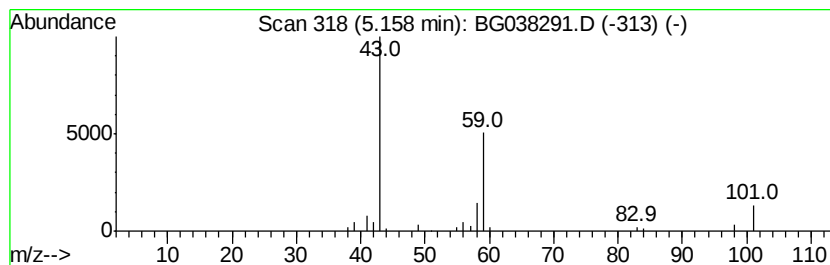
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	4.87 ng	39625	1,4-Dichlorobenzene-d4	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
4			Propane, 2-methyl-2-(1-methylethyl)-	116	C7H16O	017348-59-3	9
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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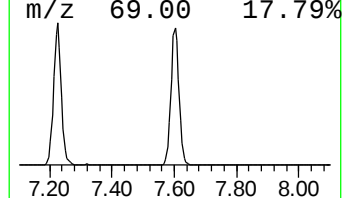
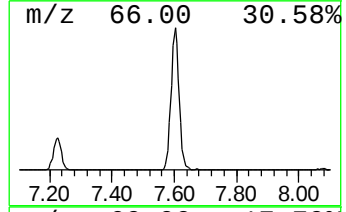
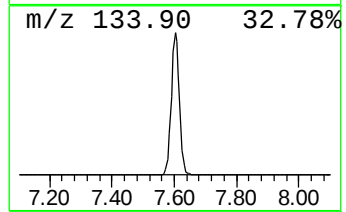
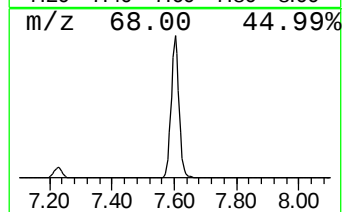
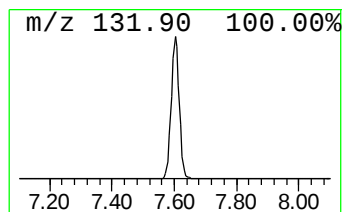
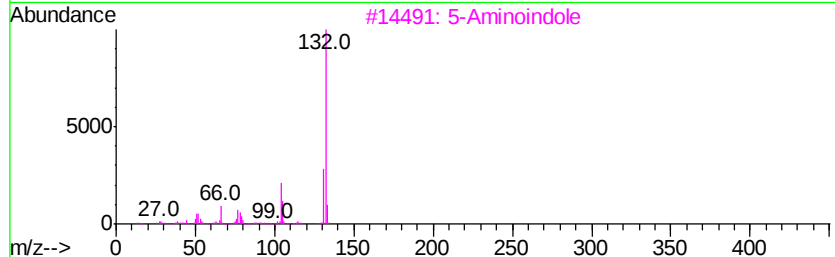
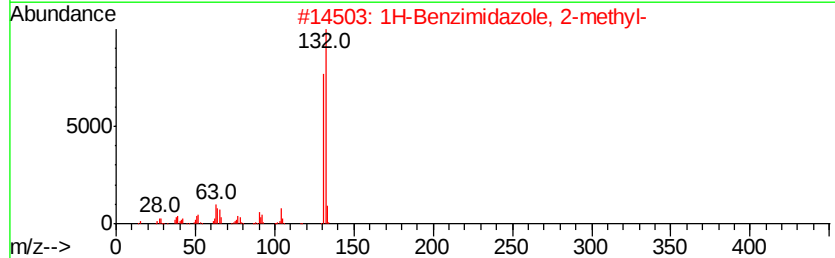
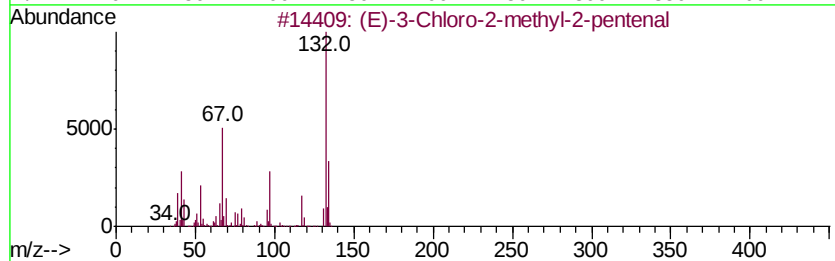
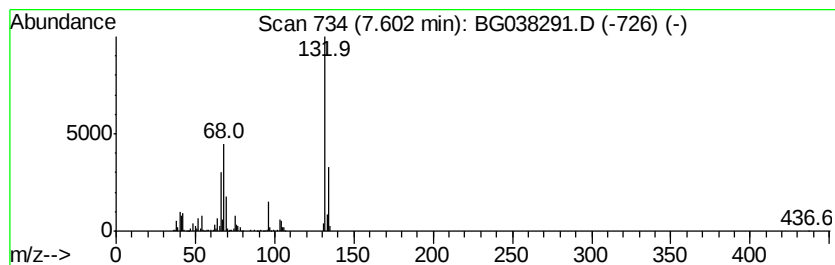
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Peak Number 2 unknown7.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	98.29 ng	800526	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	4.9	ng	39625	1	8.07	162898	20.0
unknown7.60	7.60	98.3	ng	800526	1	8.07	162898	20.0