

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112818\
 Data File : BG038205.D
 Acq On : 28 Nov 2018 18:34
 Operator : JU/SJ
 Sample : PB115117BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115117BL

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	5.167	315	320	328	rVB	21964	36402	1.90%	0.346%
2	5.631	392	399	410	rVB	470845	788692	41.22%	7.489%
3	7.230	663	671	682	rBV	496982	905442	47.32%	8.598%
4	7.611	728	736	747	rBV	474647	843708	44.10%	8.012%
5	8.082	810	816	826	rBV	122318	213612	11.16%	2.028%
6	8.405	864	871	883	rVB	340962	622289	32.52%	5.909%
7	9.262	1009	1017	1026	rBV	305714	563736	29.46%	5.353%
8	10.908	1287	1297	1304	rBV	171974	332587	17.38%	3.158%
9	13.340	1704	1711	1718	rBV	823896	1318154	68.89%	12.517%
10	14.721	1937	1946	1956	rBV2	286092	489337	25.57%	4.647%
11	16.207	2192	2199	2220	rBV2	442627	773559	40.43%	7.345%
12	17.465	2407	2413	2424	rVB2	366520	578664	30.24%	5.495%
13	20.067	2850	2856	2870	rBV	1429202	1913382	100.00%	18.169%
14	21.754	3136	3143	3156	rBV	354691	641888	33.55%	6.095%
15	25.079	3694	3709	3729	rBV2	128460	509638	26.64%	4.839%

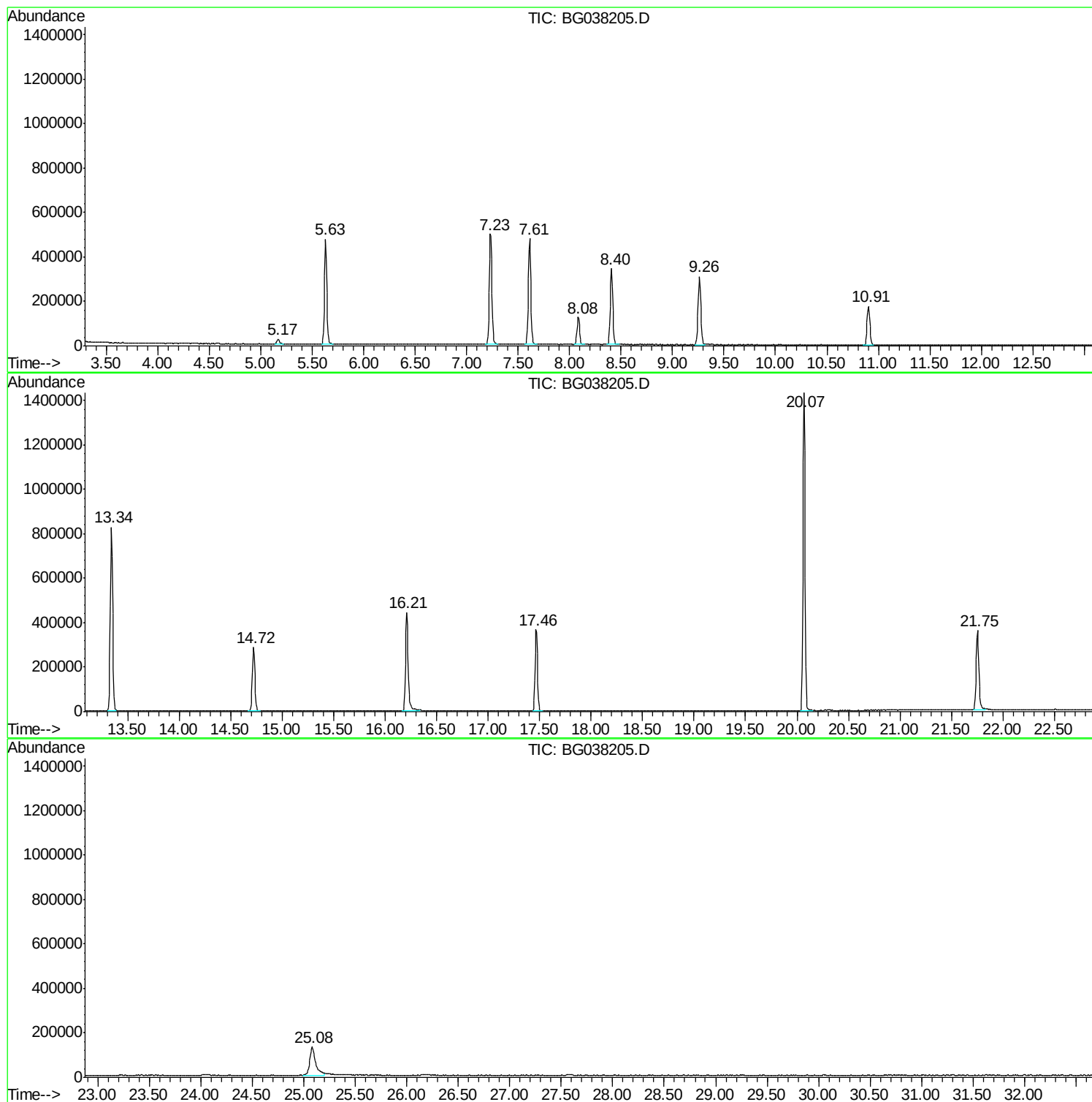
Sum of corrected areas: 10531090

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



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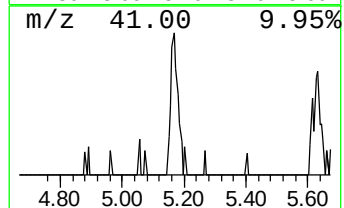
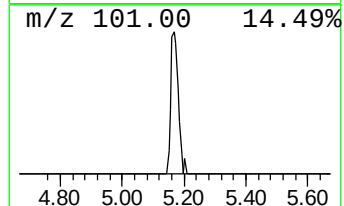
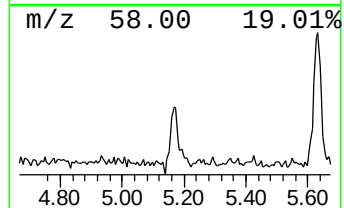
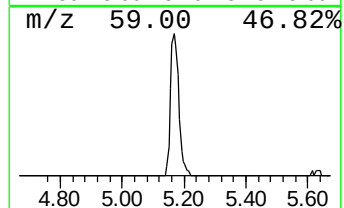
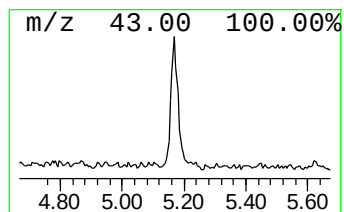
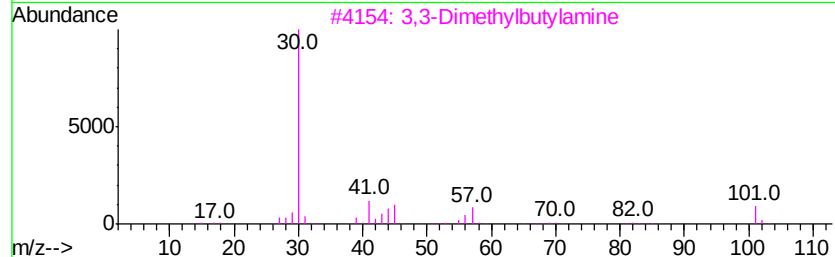
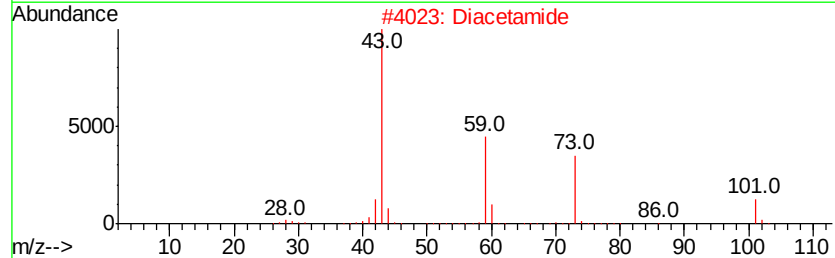
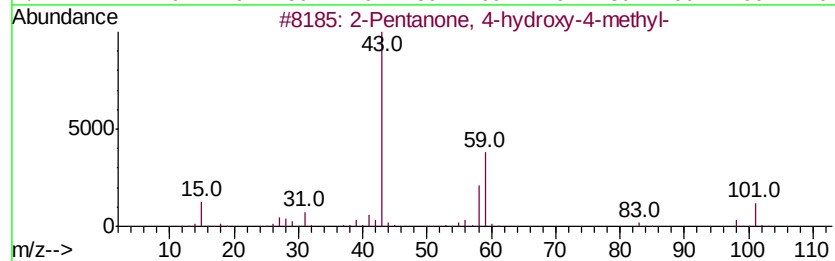
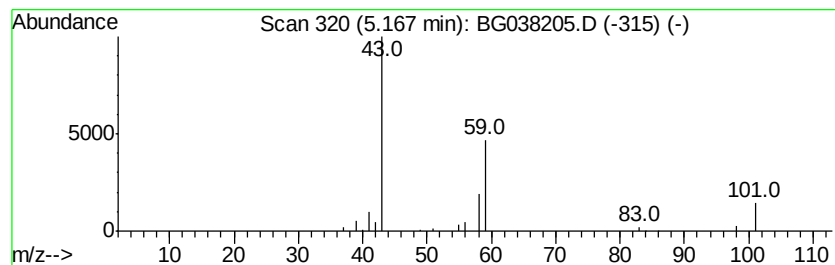
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	3.41 ng	36402	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	78
2			Diacetamide	101	C4H7NO2	000625-77-4	7
3			3,3-Dimethylbutylamine	101	C6H15N	015673-00-4	7
4			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	5
5			Acetone	58	C3H6O	000067-64-1	4



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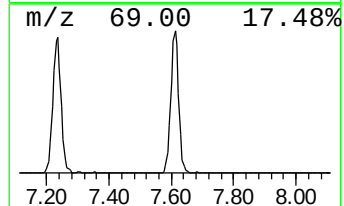
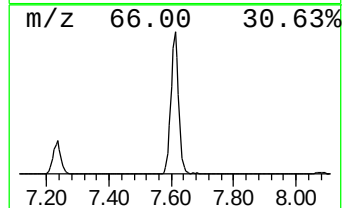
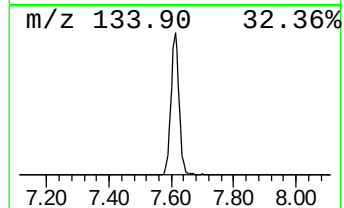
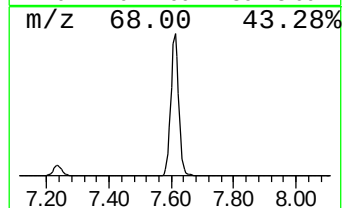
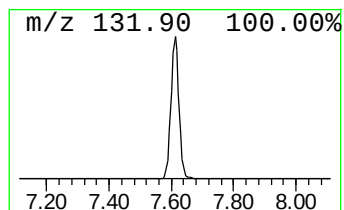
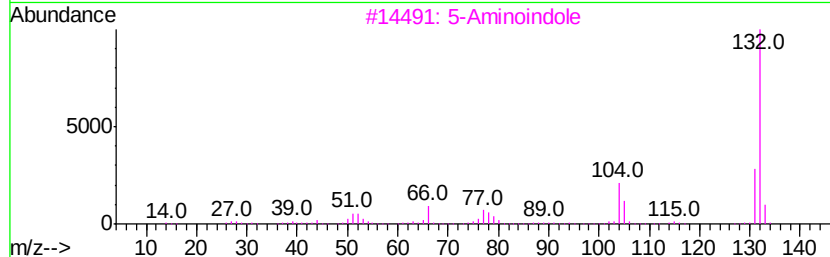
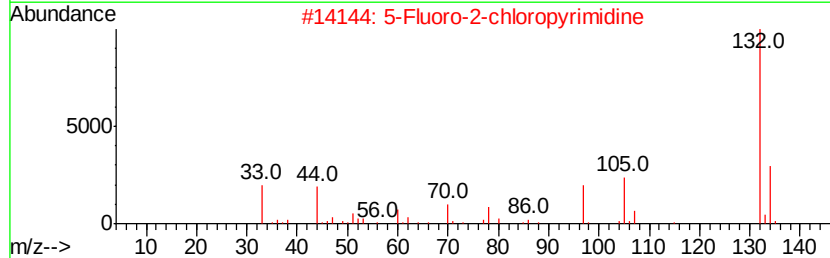
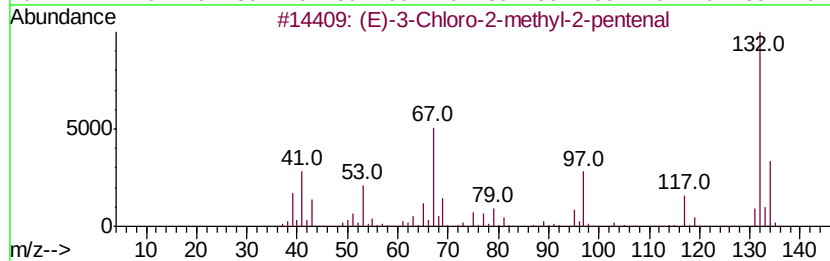
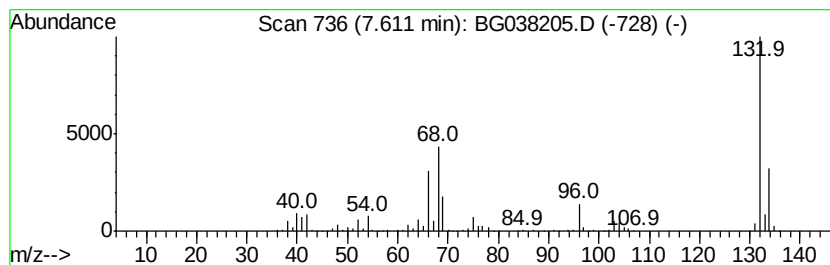
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	78.99 ng	843708	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	3.4	ng	36402	1	8.08	213612	20.0
unknown7.61	7.61	79.0	ng	843708	1	8.08	213612	20.0