

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121518\
 Data File : BG038613.D
 Acq On : 15 Dec 2018 20:46
 Operator : JU/SJ
 Sample : PB115720BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115720BL

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	5.139	345	349	358	rVB2	17461	28354	1.86%	0.395%
2	5.609	422	429	439	rVB	322136	537445	35.21%	7.479%
3	7.213	694	702	711	rBV	319749	572646	37.52%	7.968%
4	7.583	758	765	777	rVB	301691	545260	35.72%	7.587%
5	8.059	839	846	856	rVB	80255	136155	8.92%	1.895%
6	8.382	890	901	909	rBV	234095	426287	27.93%	5.932%
7	9.234	1038	1046	1056	rBV	190619	359464	23.55%	5.002%
8	10.879	1318	1326	1337	rBV	95334	182783	11.97%	2.543%
9	13.311	1733	1740	1750	rBV	492448	817312	53.54%	11.373%
10	14.692	1969	1975	1984	rBV2	149000	256846	16.83%	3.574%
11	16.185	2222	2229	2243	rBV	376313	599992	39.31%	8.349%
12	17.442	2437	2443	2454	rBV	226378	351295	23.01%	4.888%
13	20.045	2879	2886	2895	rBV	1143573	1526422	100.00%	21.240%
14	21.719	3164	3171	3183	rBV	256023	440250	28.84%	6.126%
15	25.015	3721	3732	3744	rVB3	140416	405895	26.59%	5.648%

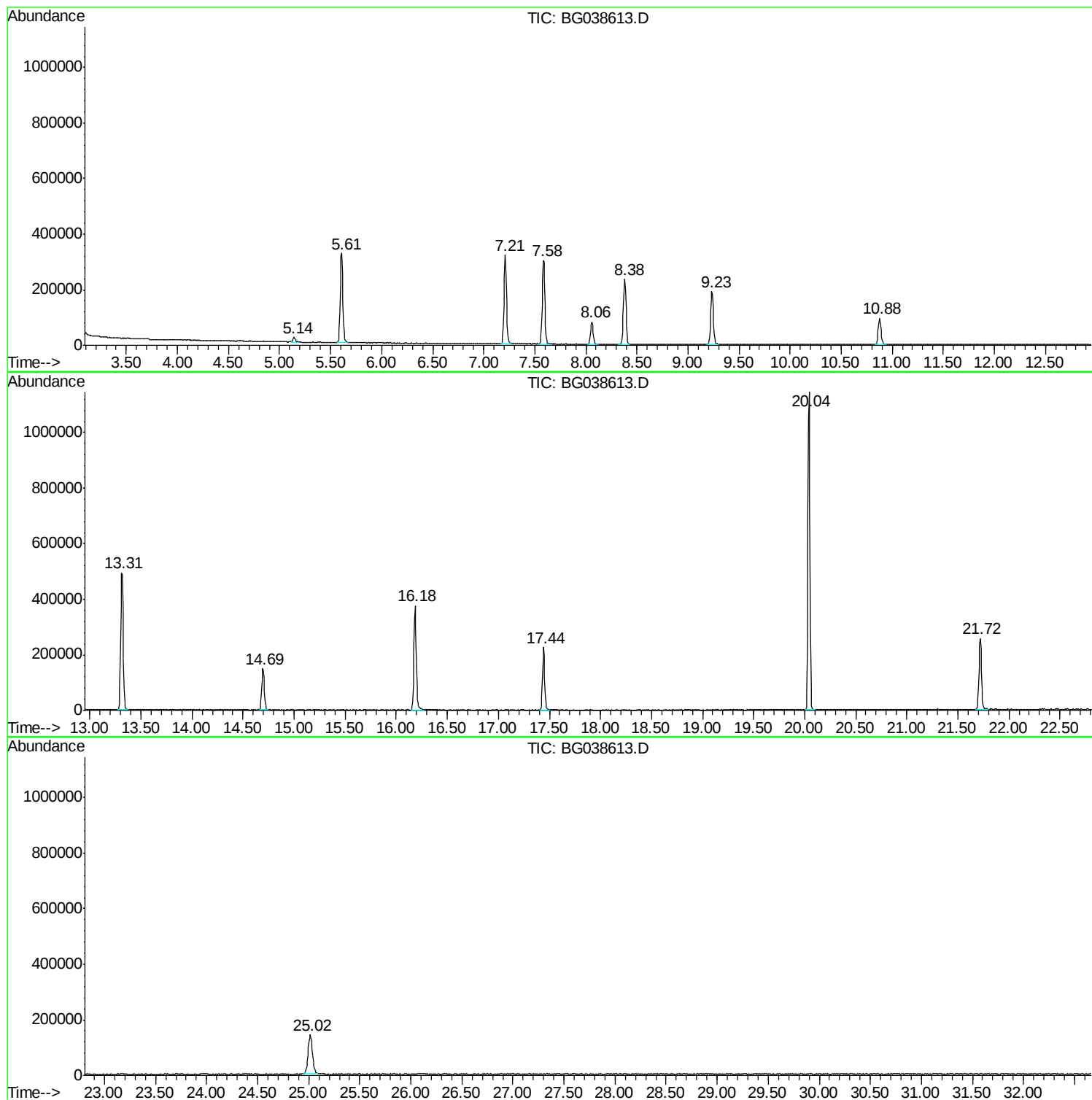
Sum of corrected areas: 7186406

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



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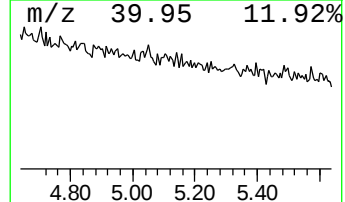
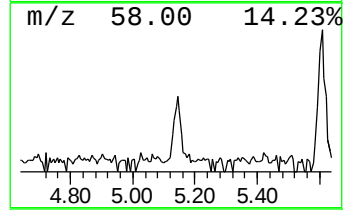
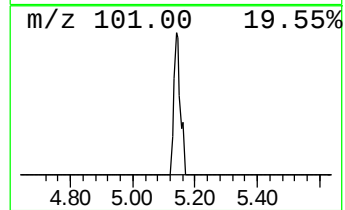
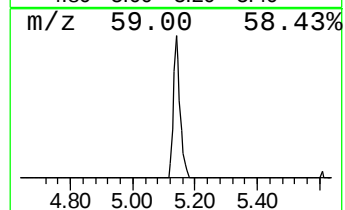
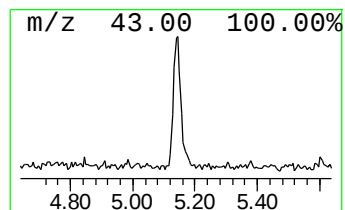
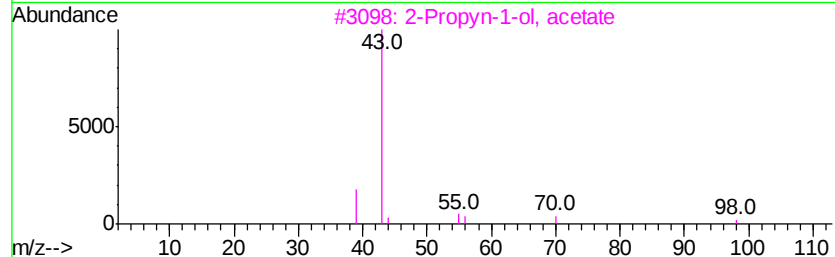
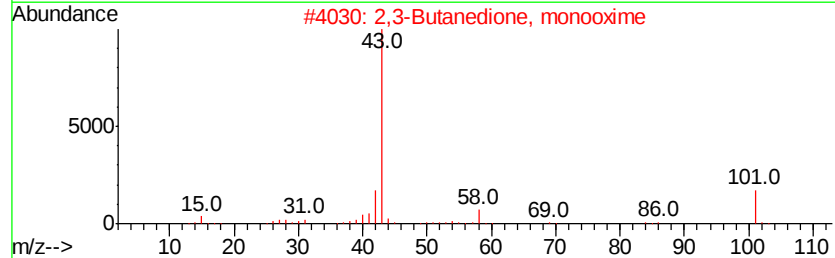
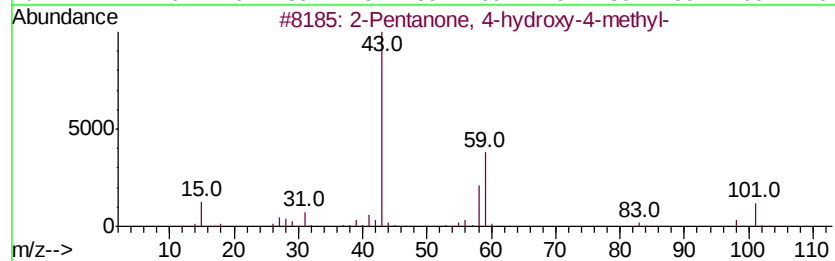
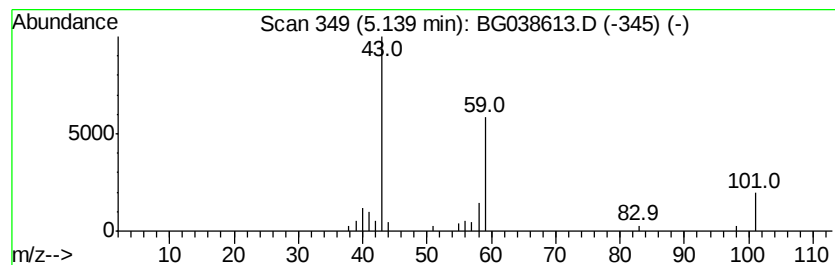
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	4.16 ng	28354	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3			2-Propyn-1-ol, acetate	98	C5H6O2	000627-09-8	9
4			2-Heptanone, 7,7-dichloro-	182	C7H12Cl2O	066241-43-8	9
5			3-Hexanol	102	C6H14O	000623-37-0	9



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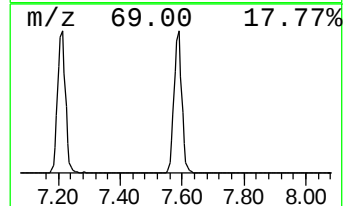
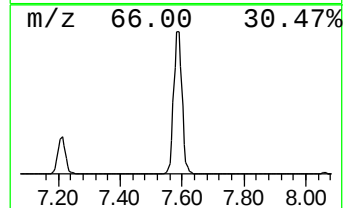
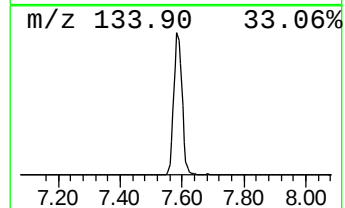
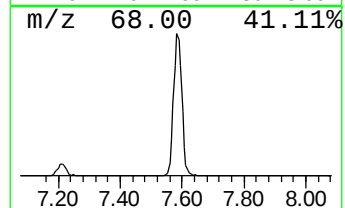
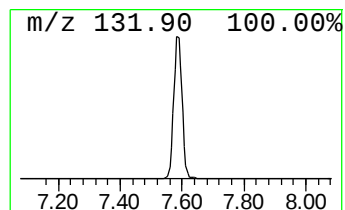
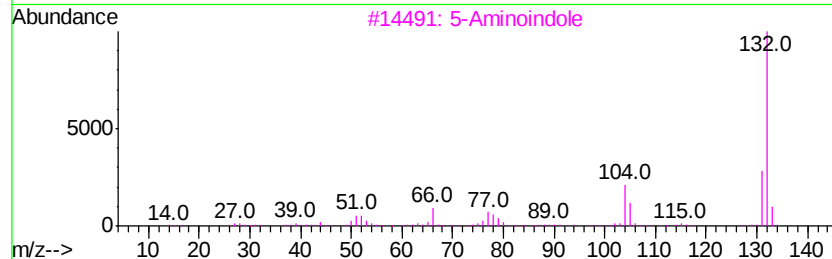
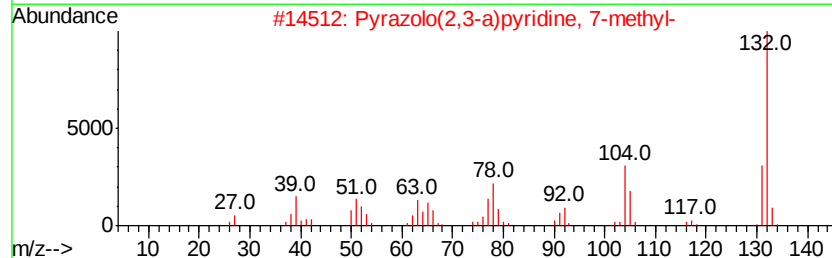
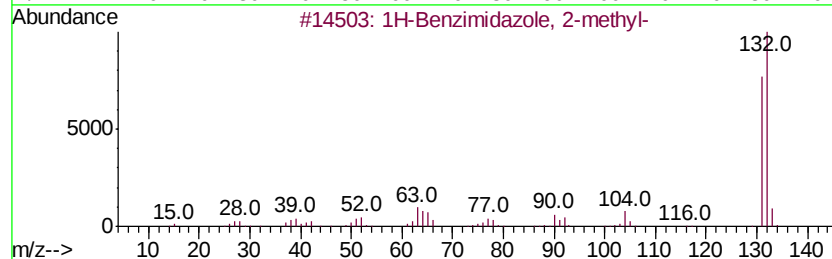
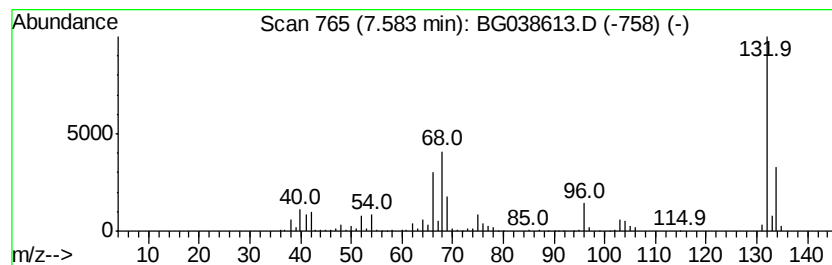
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	80.09 ng	545260	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	22
3		5-Aminoindole	132	C8H8N2	005192-03-0	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	4.2	ng	28354	1	8.06	136155	20.0
unknown7.58	7.58	80.1	ng	545260	1	8.06	136155	20.0