

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112718\
 Data File : BG038183.D
 Acq On : 28 Nov 2018 1:33
 Operator : JU/SJ
 Sample : PB115087BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115087BL

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	5.170	315	320	331	rVB	39863	71782	2.17%	0.429%
2	5.634	391	399	409	rBV	835284	1398921	42.20%	8.354%
3	7.238	664	672	685	rBV	876112	1567374	47.28%	9.360%
4	7.614	727	736	747	rBV	819434	1460005	44.04%	8.718%
5	8.084	809	816	824	rBV	152960	266840	8.05%	1.593%
6	8.407	864	871	881	rVB	595533	1075298	32.43%	6.421%
7	9.265	1009	1017	1025	rBV	520977	970593	29.28%	5.796%
8	10.904	1288	1296	1304	rVB	210801	392928	11.85%	2.346%
9	13.343	1703	1711	1720	rBV	1398336	2257624	68.10%	13.481%
10	14.724	1937	1946	1956	rBV2	339824	563823	17.01%	3.367%
11	16.210	2192	2199	2216	rBV3	788163	1355102	40.87%	8.092%
12	17.473	2407	2414	2423	rVB	428201	686821	20.72%	4.101%
13	20.070	2850	2856	2867	rBV	2466028	3315278	100.00%	19.797%
14	21.751	3135	3142	3154	rBV2	385948	755390	22.79%	4.511%
15	25.082	3693	3709	3727	rBV2	154647	608324	18.35%	3.633%

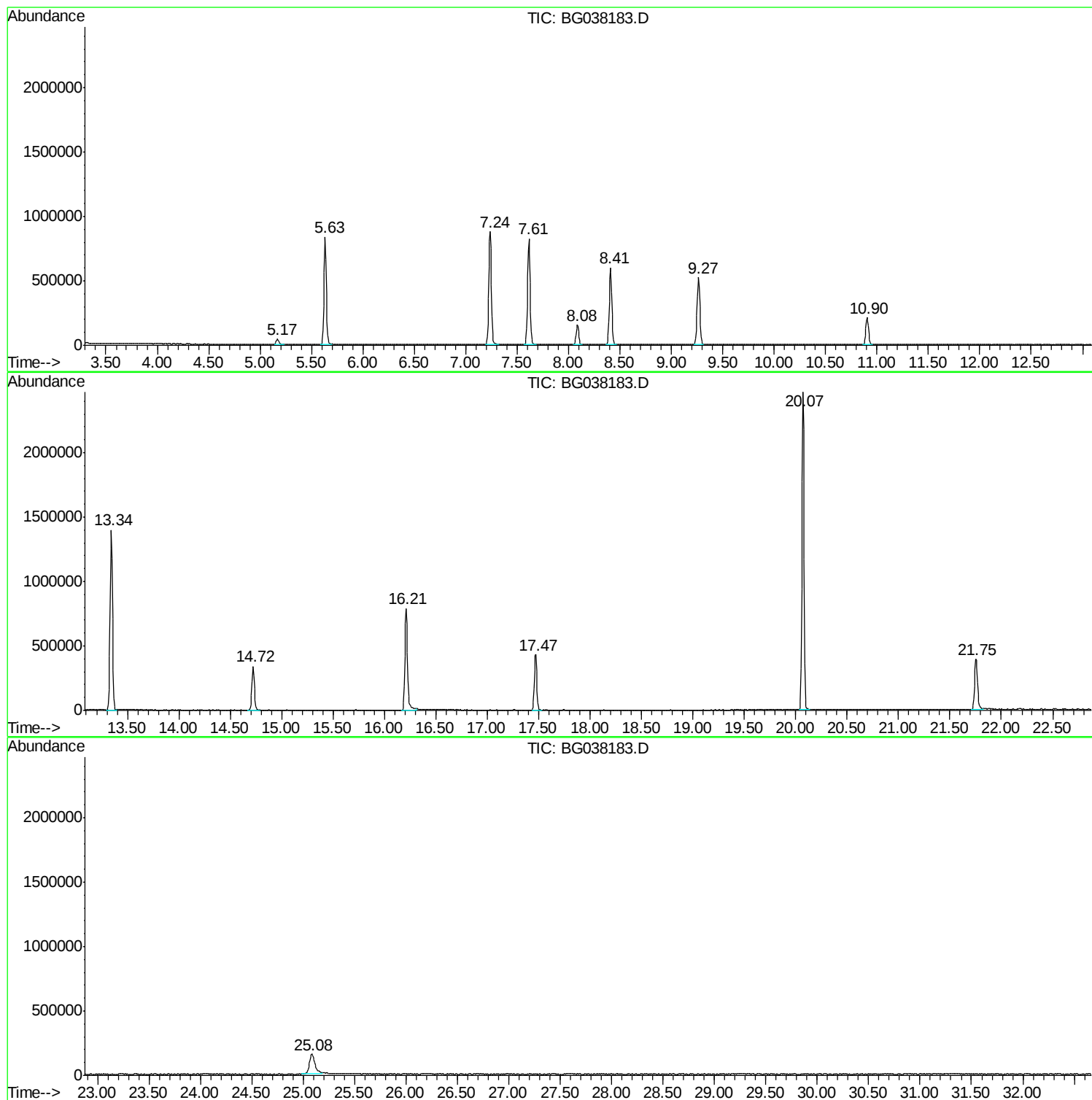
Sum of corrected areas: 16746103

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



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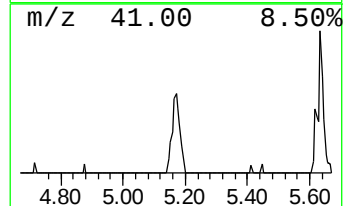
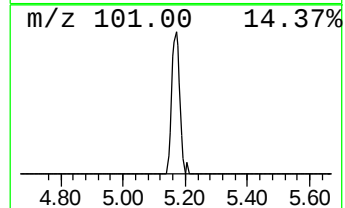
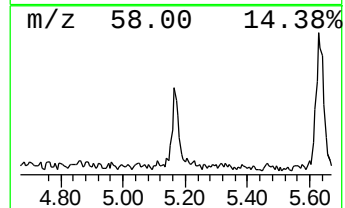
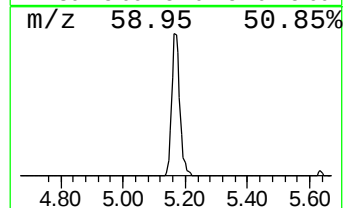
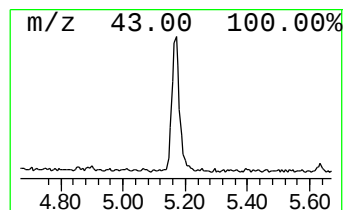
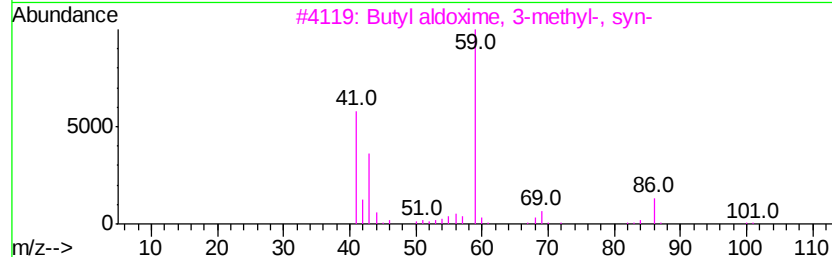
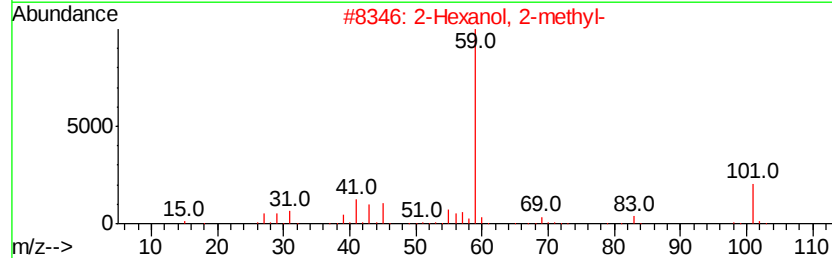
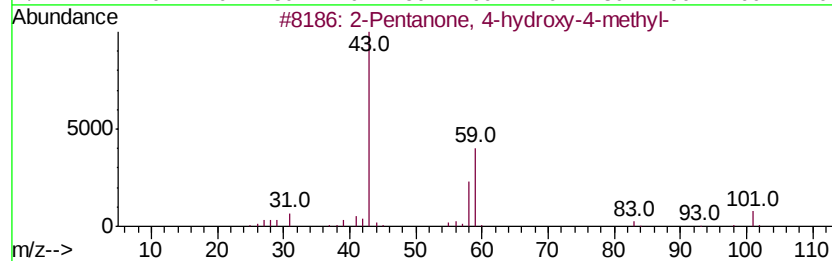
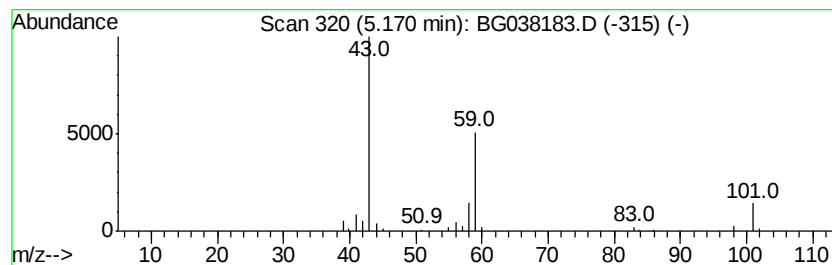
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	5.38 ng	71782	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	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
4			Acetamide, N-(aminocarbonyl)-	102	C3H6N2O2	000591-07-1	9
5			Butane	58	C4H10	000106-97-8	9



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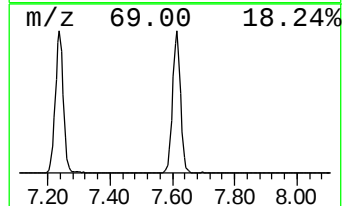
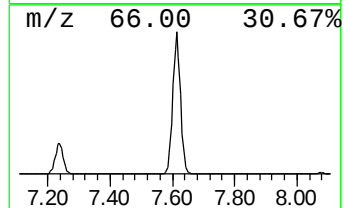
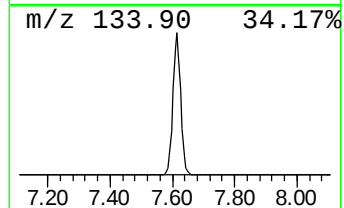
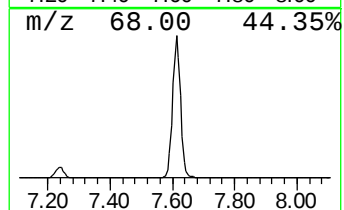
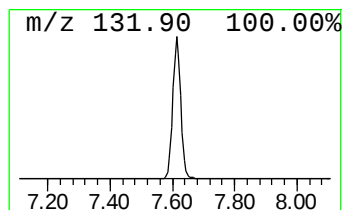
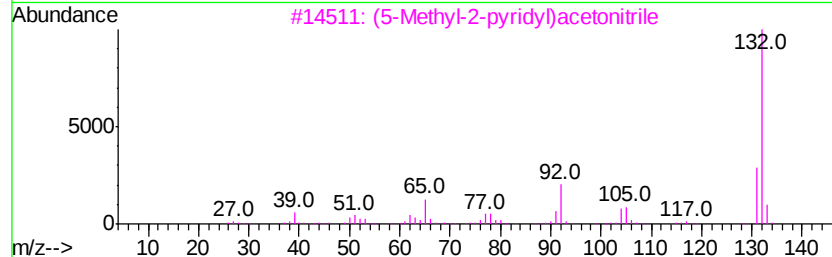
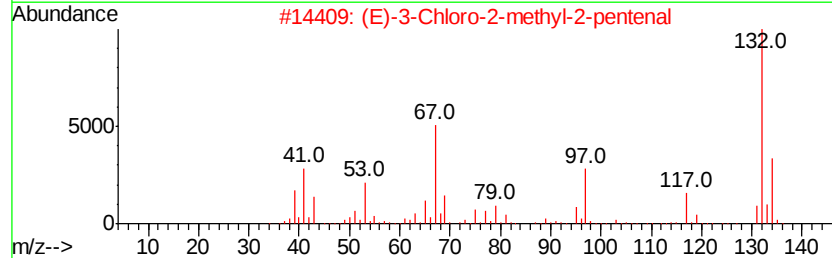
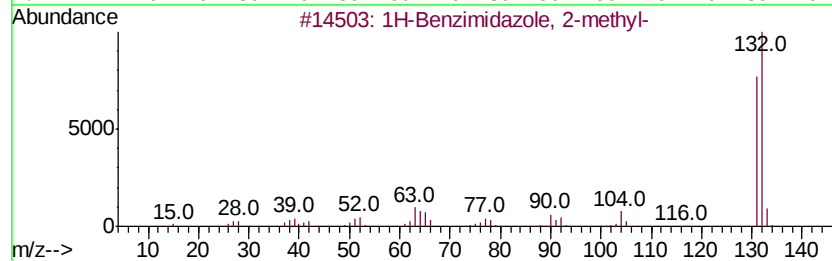
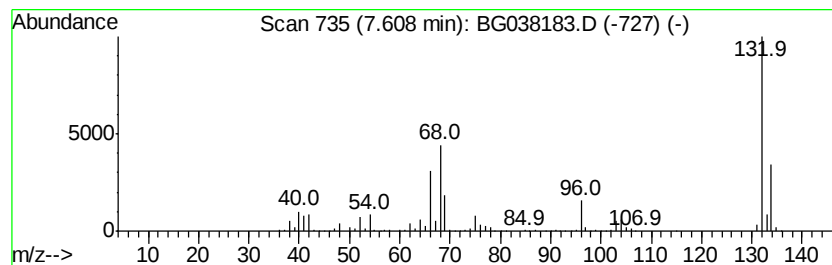
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	109.43 ng	1460010	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	5.4	ng	71782	1	8.08	266840	20.0
unknown7.61	7.61	109.4	ng	1460010	1	8.08	266840	20.0