

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
 Data File : BG038045.D
 Acq On : 16 Nov 2018 22:34
 Operator : JU/SJ
 Sample : PB114870BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114870BL

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	5.169	314	320	328	rBV	37119	64871	2.06%	0.399%
2	5.627	391	398	409	rBV	772000	1288738	40.85%	7.935%
3	7.231	663	671	682	rBV	794431	1419570	45.00%	8.741%
4	7.613	728	736	744	rBV	753795	1346957	42.70%	8.294%
5	8.083	809	816	824	rVB	149151	267486	8.48%	1.647%
6	8.406	864	871	879	rVB	572405	1017240	32.25%	6.263%
7	9.264	1008	1017	1033	rVB	472348	887071	28.12%	5.462%
8	10.903	1289	1296	1305	rVB	216390	394739	12.51%	2.431%
9	13.342	1703	1711	1723	rBV	1338736	2087186	66.16%	12.851%
10	14.722	1939	1946	1954	rBV2	330827	561507	17.80%	3.457%
11	16.209	2192	2199	2215	rBV3	937134	1468389	46.55%	9.041%
12	17.466	2404	2413	2421	rBV2	458819	709964	22.51%	4.371%
13	20.069	2850	2856	2863	rBV	2429972	3154663	100.00%	19.424%
14	21.750	3135	3142	3152	rBV	470105	798676	25.32%	4.918%
15	25.075	3692	3708	3722	rBV2	244711	773794	24.53%	4.764%

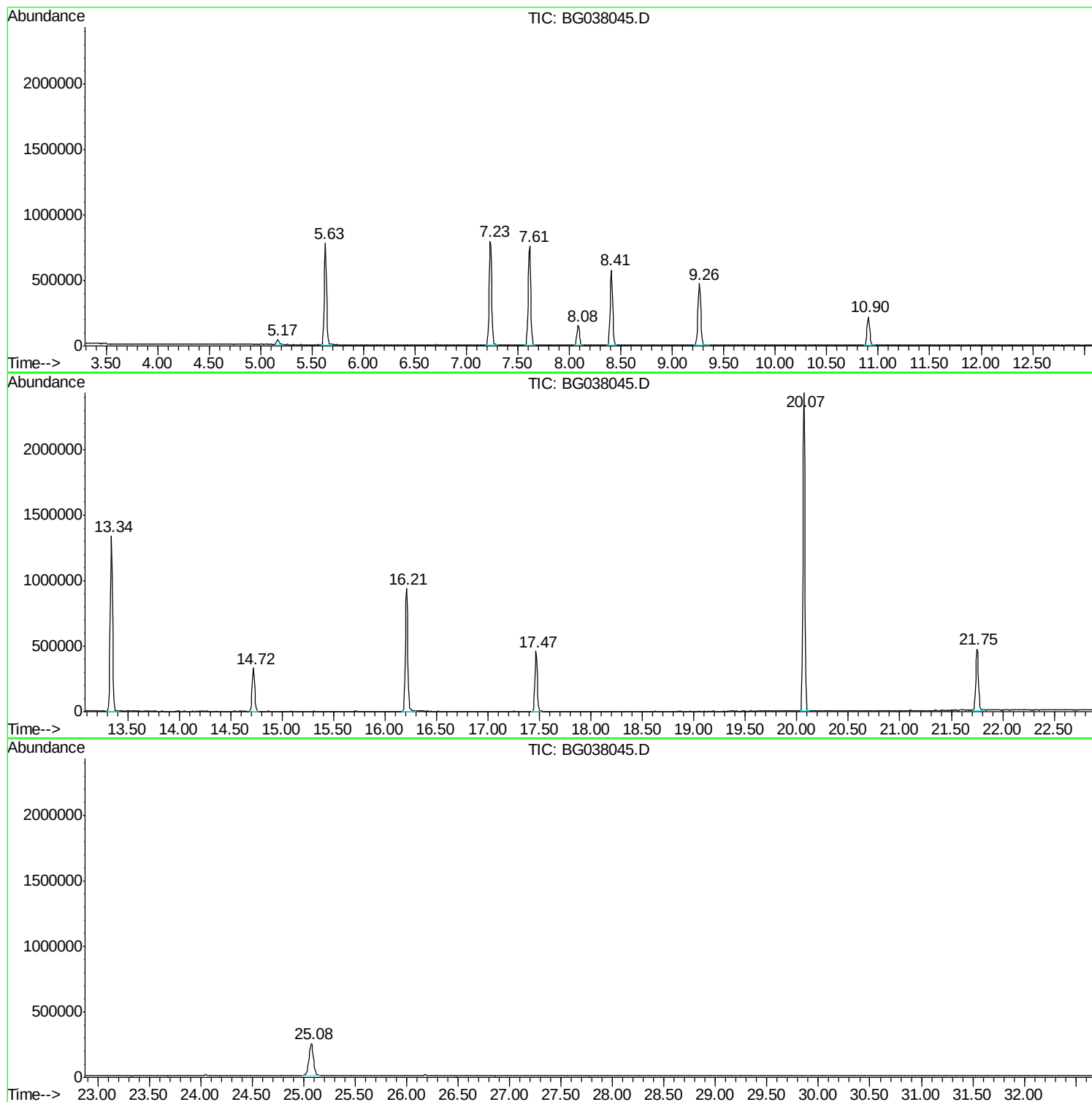
Sum of corrected areas: 16240851

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



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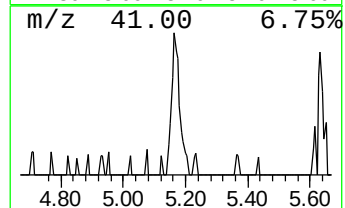
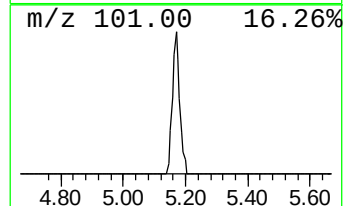
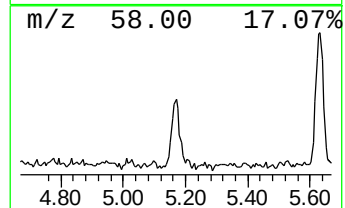
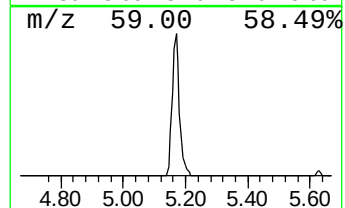
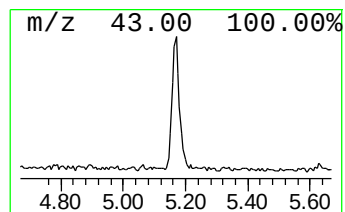
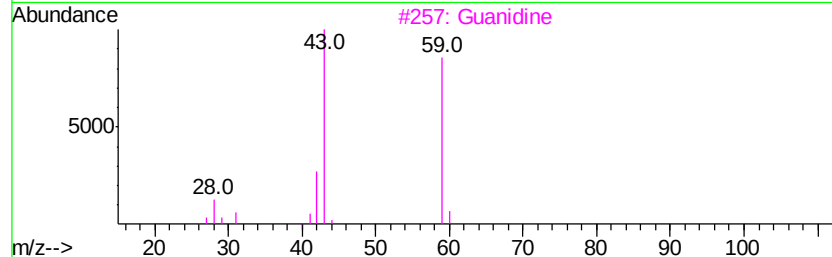
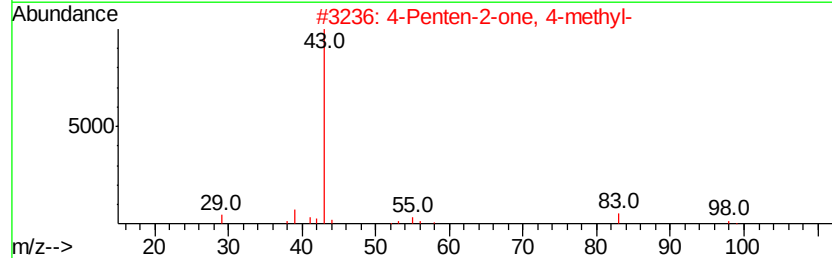
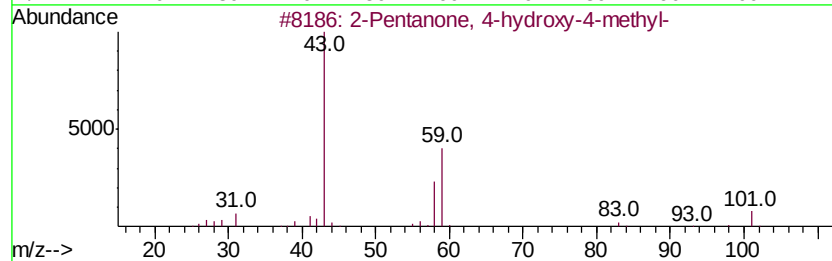
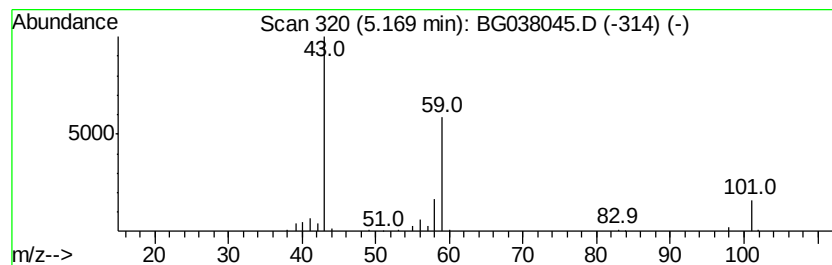
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	4.85 ng	64871	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	64
2			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
3			Guanidine	59	CH5N3	000113-00-8	9
4			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9



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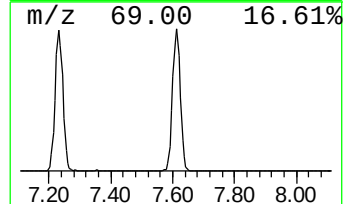
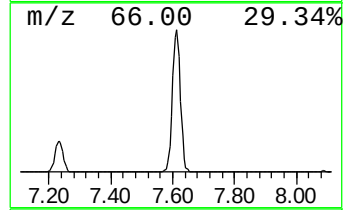
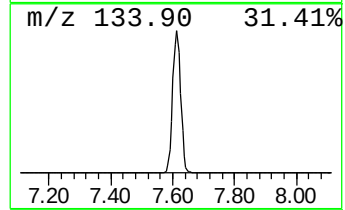
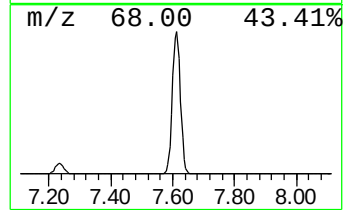
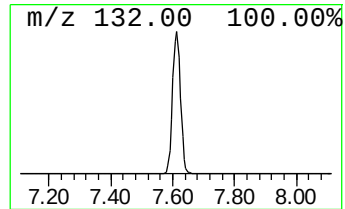
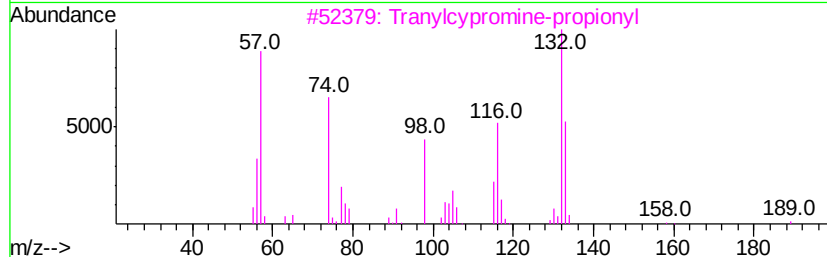
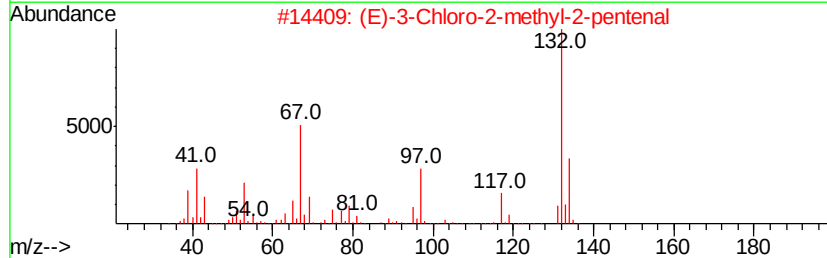
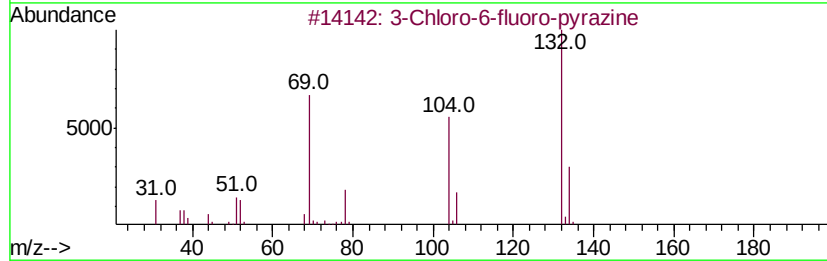
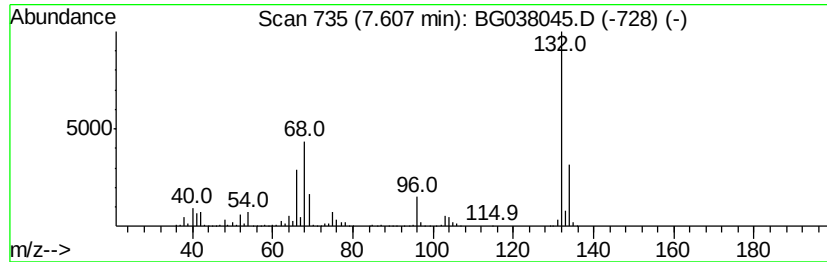
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	100.71 ng	1346960	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	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
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	4.8	ng	64871	1	8.08	267486	20.0
unknown7.61	7.61	100.7	ng	1346960	1	8.08	267486	20.0