

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110818\
 Data File : BG037899.D
 Acq On : 8 Nov 2018 23:47
 Operator : JU/SJ
 Sample : PB114622BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BG101818.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

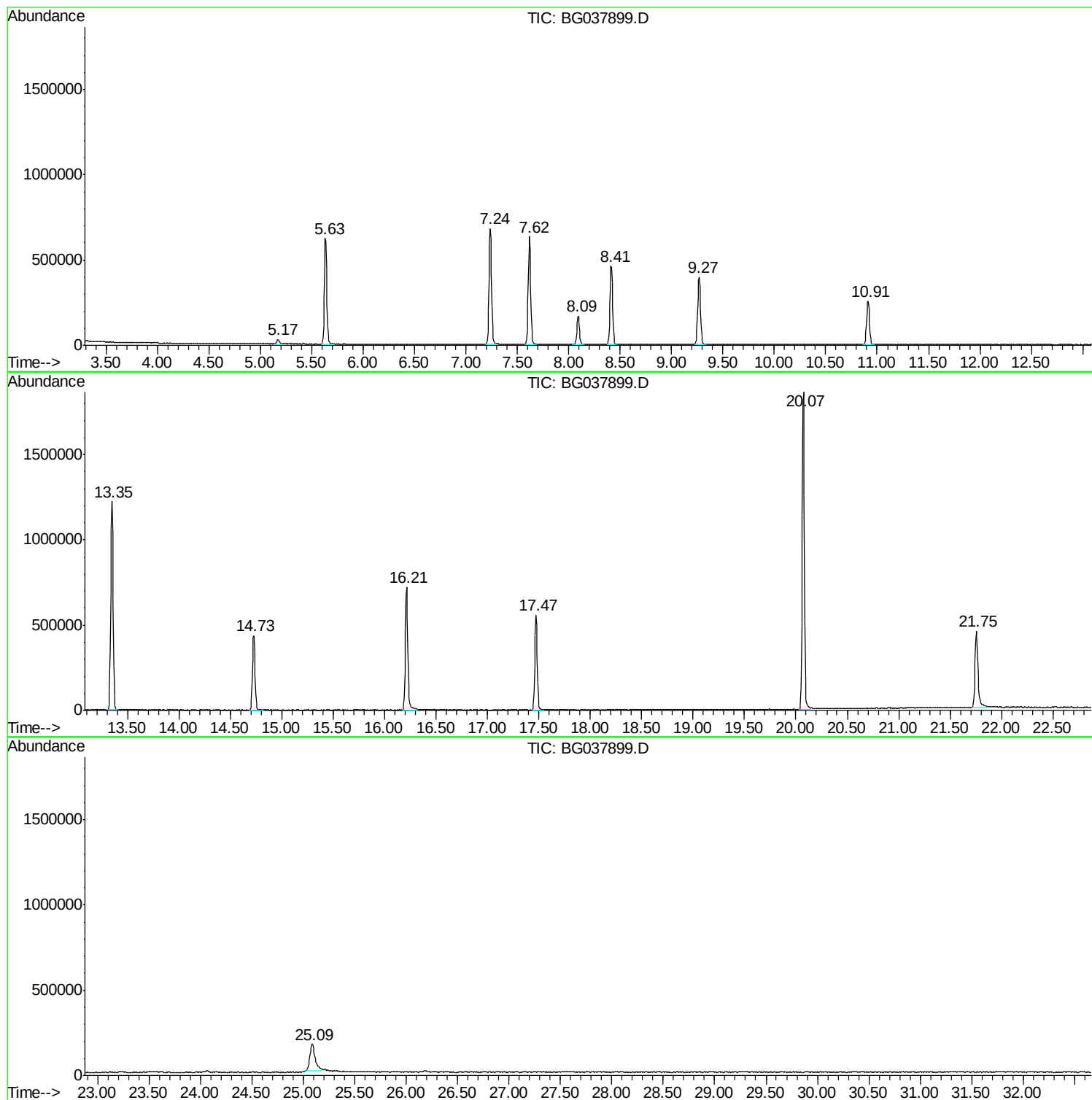
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.174	315	321	327	rBV	26827	46442	1.74%	0.316%
2	5.632	392	399	413	rBV	621737	1051504	39.51%	7.152%
3	7.236	664	672	681	rBV	680131	1201622	45.15%	8.173%
4	7.618	729	737	747	rBV	634246	1128000	42.38%	7.672%
5	8.088	808	817	827	rVB	165265	294407	11.06%	2.002%
6	8.411	865	872	880	rVB	459886	851826	32.01%	5.794%
7	9.269	1010	1018	1029	rBV	395379	741330	27.85%	5.042%
8	10.914	1288	1298	1307	rBV	255484	485107	18.23%	3.300%
9	13.347	1705	1712	1721	rBV	1220919	1936925	72.78%	13.174%
10	14.728	1939	1947	1962	rBV2	436162	739682	27.79%	5.031%
11	16.214	2193	2200	2217	rBV5	721611	1212621	45.56%	8.248%
12	17.471	2407	2414	2426	rBV	555867	875012	32.88%	5.952%
13	20.074	2850	2857	2869	rBV	1860498	2661439	100.00%	18.102%
14	21.755	3136	3143	3161	rBV2	447857	886546	33.31%	6.030%
15	25.086	3698	3710	3724	rBV2	161281	589727	22.16%	4.011%

Sum of corrected areas: 14702190

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



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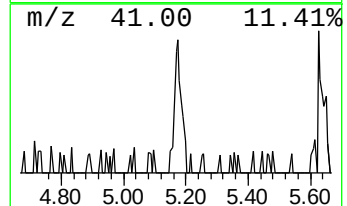
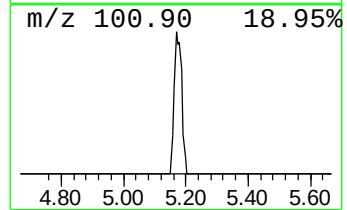
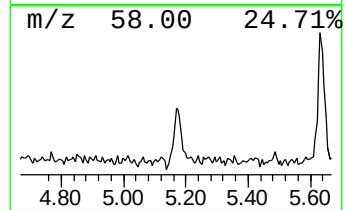
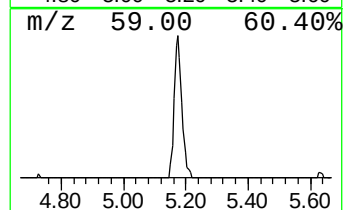
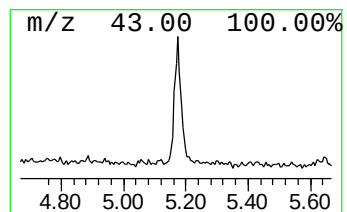
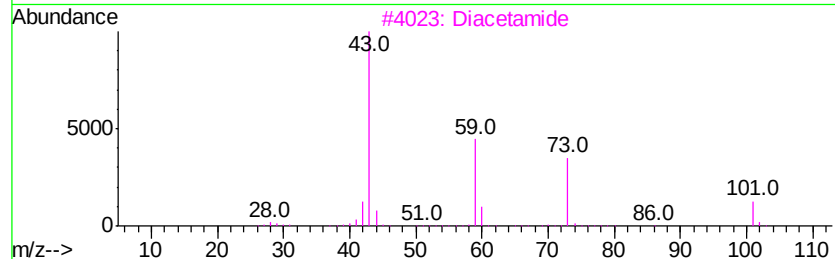
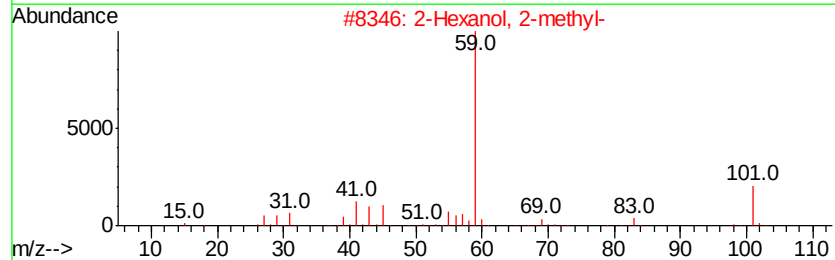
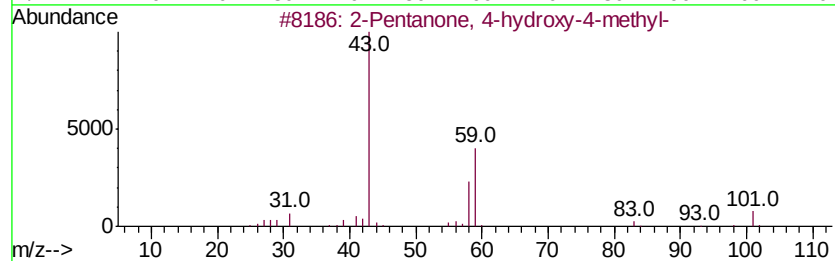
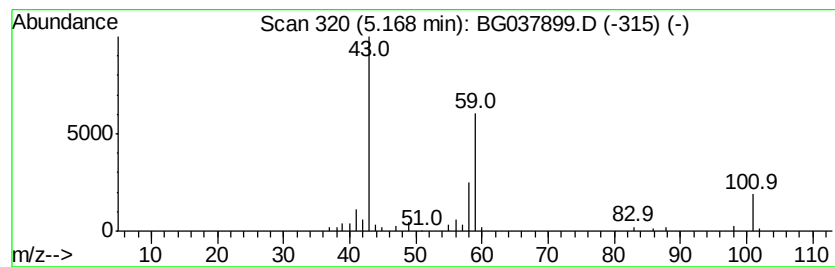
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	3.15 ng	46442	1,4-Dichlorobenzene-d4	8.09

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		Diacetamide	101	C4H7NO2	000625-77-4	9
4		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	9
5		2-Pentanol, 2,3-dimethyl-	116	C7H16O	004911-70-0	9



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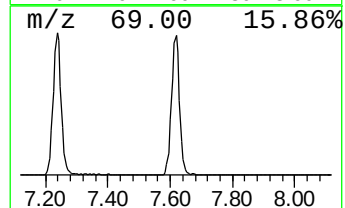
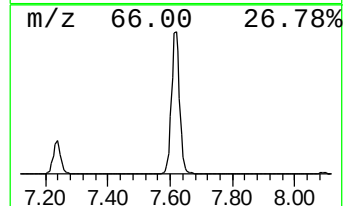
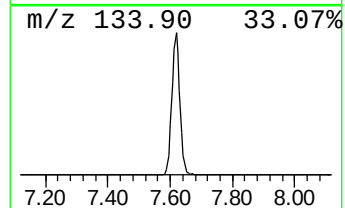
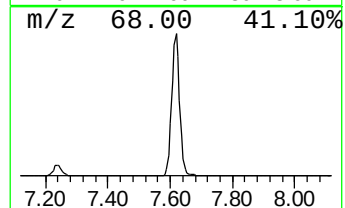
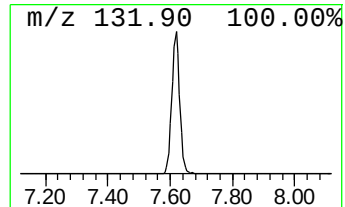
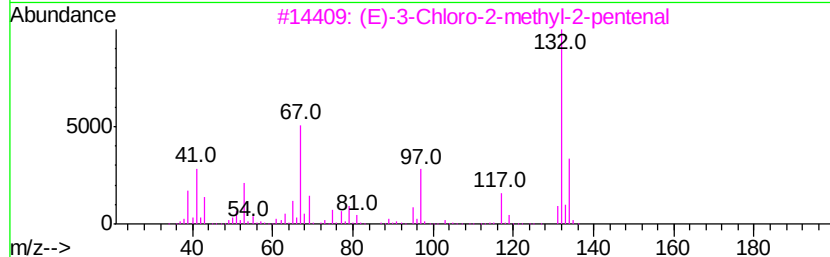
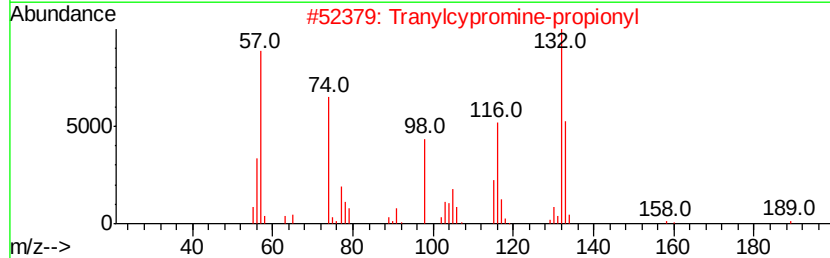
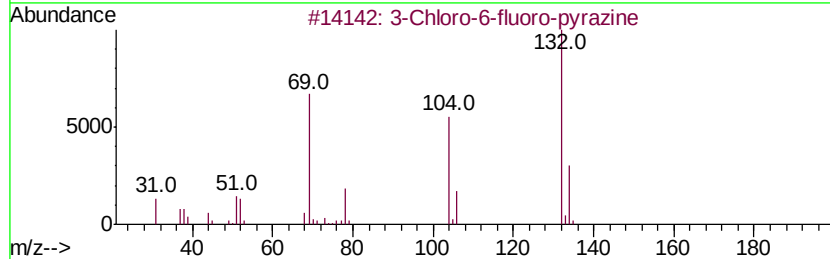
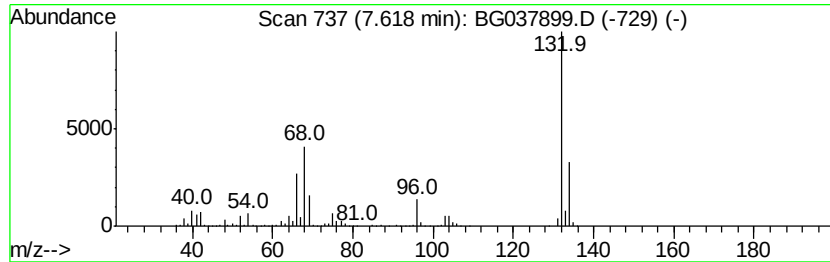
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	76.63 ng	1128000	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		Xenon	132	Xe	007440-63-3	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	3.1	ng	46442	1	8.09	294407	20.0
unknown7.62	7.62	76.6	ng	1128000	1	8.09	294407	20.0