

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101518\
 Data File : BG037346.D
 Acq On : 15 Oct 2018 23:14
 Operator : JU/SJ
 Sample : PB113837BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

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-BG101118.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

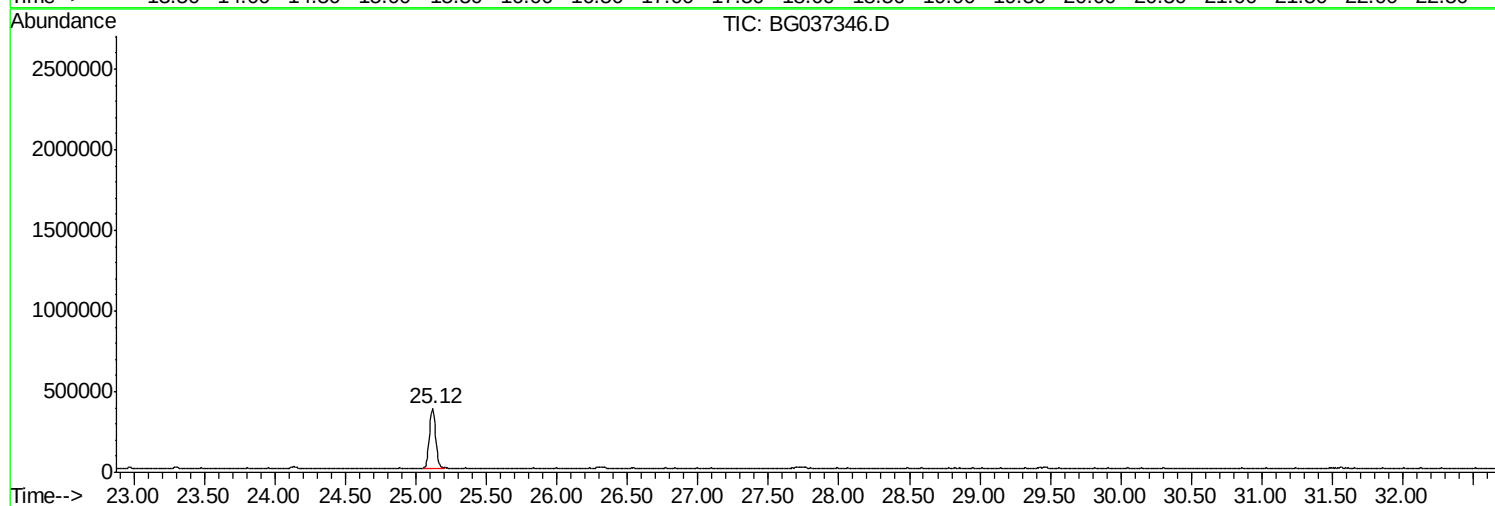
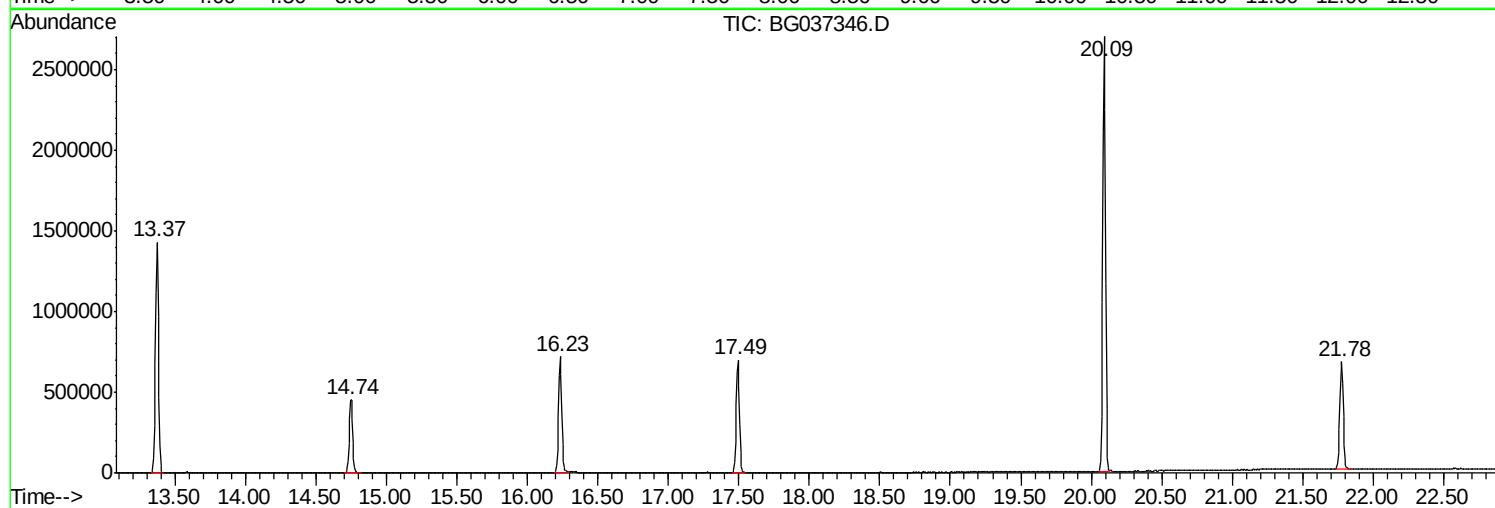
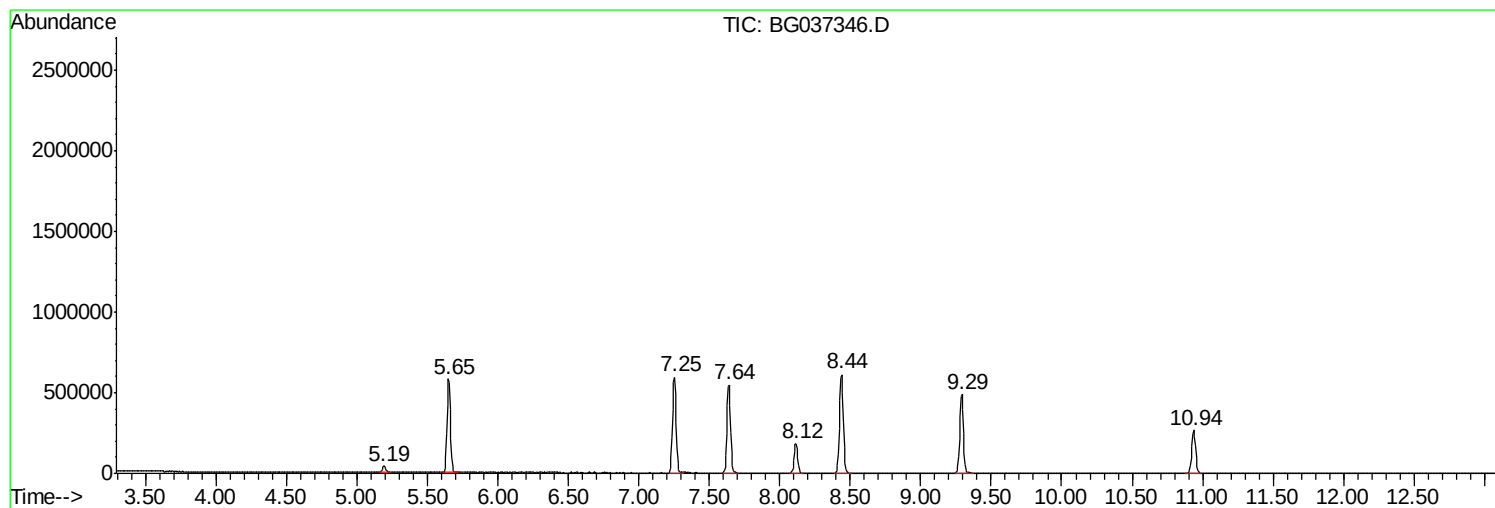
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.191	319	324	330	rVB	39303	62542	1.74%	0.369%
2	5.649	395	402	414	rBV	581798	945100	26.32%	5.569%
3	7.253	668	675	685	rBV	593245	1028978	28.66%	6.064%
4	7.641	733	741	751	rBV	545597	975417	27.17%	5.748%
5	8.117	814	822	829	rBV	182864	323560	9.01%	1.907%
6	8.440	868	877	888	rVB	609881	1114778	31.05%	6.569%
7	9.292	1014	1022	1035	rBV	489244	887450	24.72%	5.230%
8	10.937	1292	1302	1310	rBV	265168	493145	13.74%	2.906%
9	13.369	1709	1716	1723	rBV	1427189	2353454	65.55%	13.869%
10	14.744	1943	1950	1963	rVB2	453211	763884	21.28%	4.502%
11	16.231	2196	2203	2213	rBV	716788	1109842	30.91%	6.540%
12	17.494	2411	2418	2426	rBV	698100	1066788	29.71%	6.287%
13	20.091	2854	2860	2867	rBV	2697648	3590159	100.00%	21.157%
14	21.777	3140	3147	3157	rVB	661919	1136269	31.65%	6.696%
15	25.120	3704	3716	3729	rBV2	370384	1118077	31.14%	6.589%

Sum of corrected areas: 16969443

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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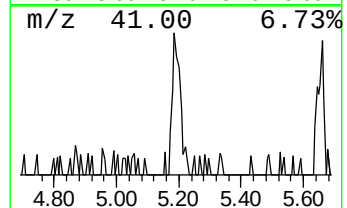
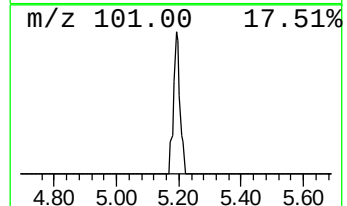
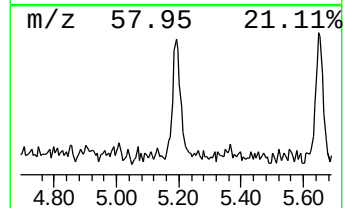
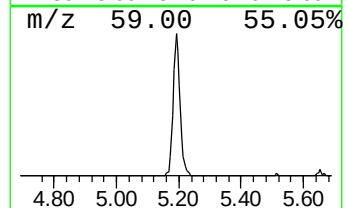
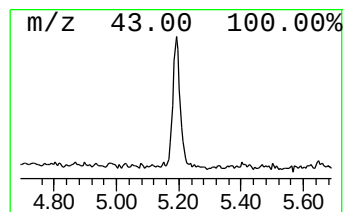
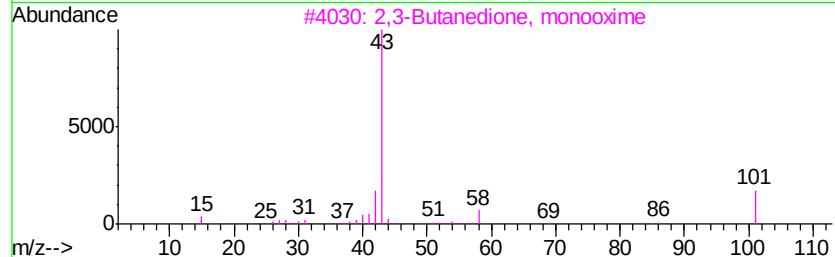
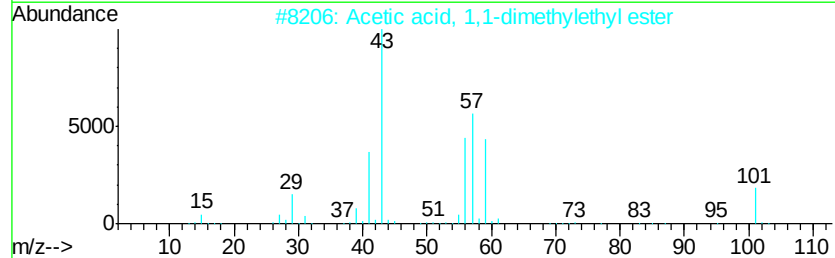
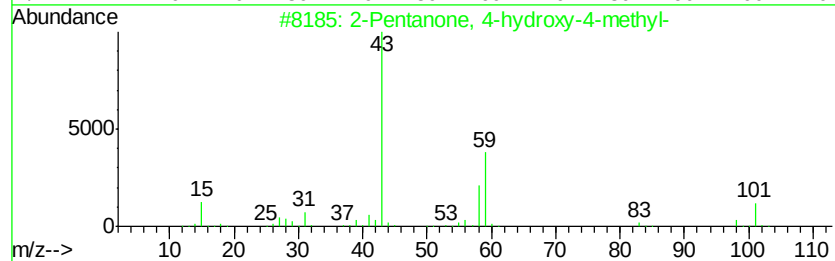
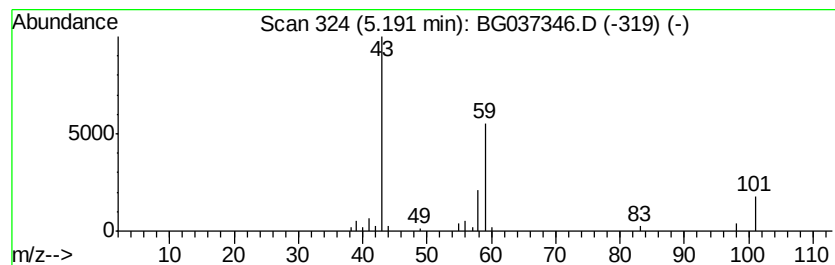
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.87 ng	62542	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Diacetamide	101	C4H7NO2	000625-77-4	9
5			Acetone	58	C3H6O	000067-64-1	9



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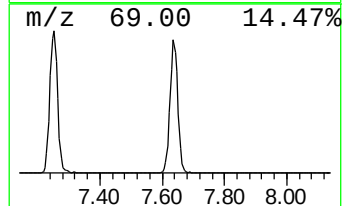
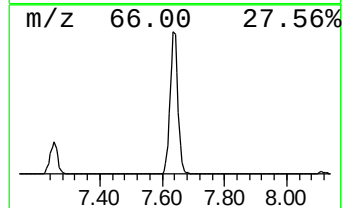
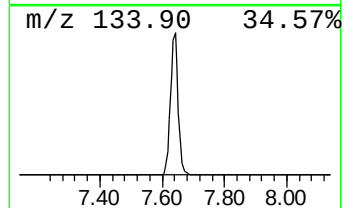
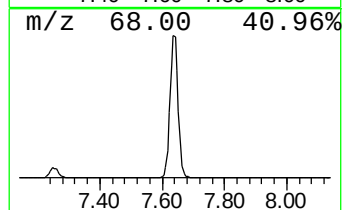
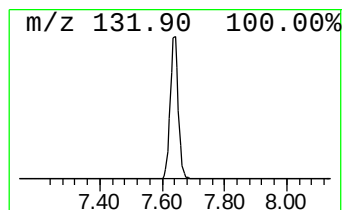
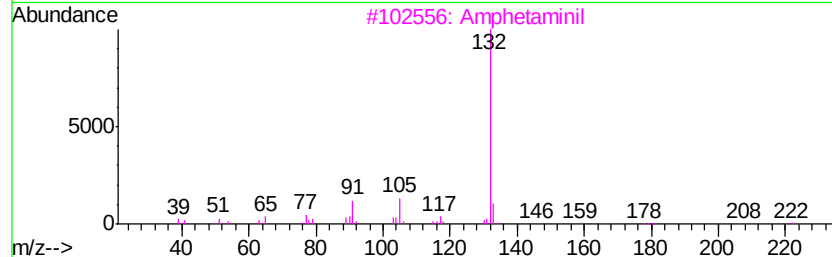
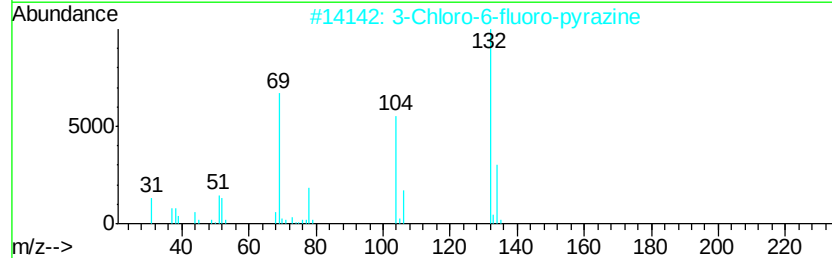
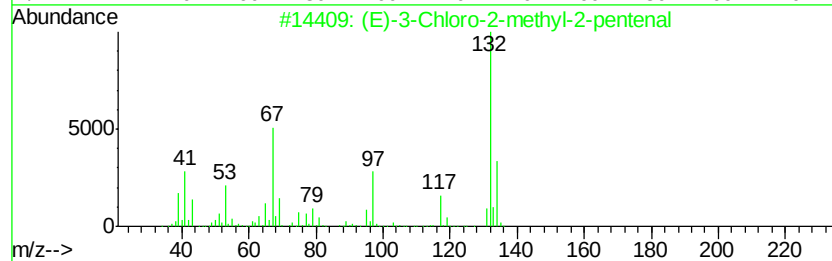
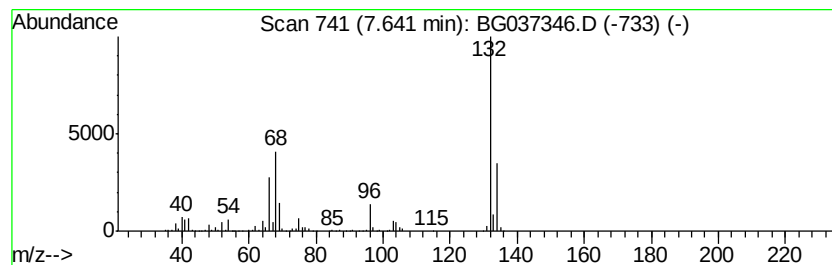
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	60.29 ng	975417	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Amphetaminil	250	C17H18N2	017590-01-1	10
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
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...	5.19	3.9	ng	62542	1	8.11	323560	20.0
unknown7.64	7.64	60.3	ng	975417	1	8.11	323560	20.0