

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011119\
Data File : BG039092.D
Acq On : 11 Jan 2019 19:27
Operator : JU/SJ
Sample : PB116292BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB116292BL

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-BG010919.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	5.110	338	344	350	rBV	18179	31756	1.55%	0.361%
2	5.580	417	424	440	rVB	346521	594461	28.95%	6.751%
3	7.184	687	697	706	rBV	368701	649294	31.62%	7.374%
4	7.554	752	760	771	rBV	357341	617763	30.08%	7.016%
5	8.024	833	840	849	rVB	78189	133379	6.49%	1.515%
6	8.347	887	895	904	rVB	284455	497225	24.21%	5.647%
7	9.205	1033	1041	1055	rBV	196948	382789	18.64%	4.347%
8	10.844	1311	1320	1332	rBV	97847	191691	9.33%	2.177%
9	13.282	1728	1735	1746	rBV	645811	1030998	50.20%	11.709%
10	14.669	1963	1971	1981	rBV2	190521	335571	16.34%	3.811%
11	16.155	2218	2224	2246	rBV	459069	787054	38.32%	8.938%
12	17.419	2432	2439	2449	rBV	281710	451333	21.98%	5.126%
13	20.016	2875	2881	2891	rBV	1537134	2053685	100.00%	23.323%
14	21.690	3159	3166	3175	rBV2	324007	537726	26.18%	6.107%
15	24.963	3712	3723	3738	rBV2	164130	510587	24.86%	5.799%

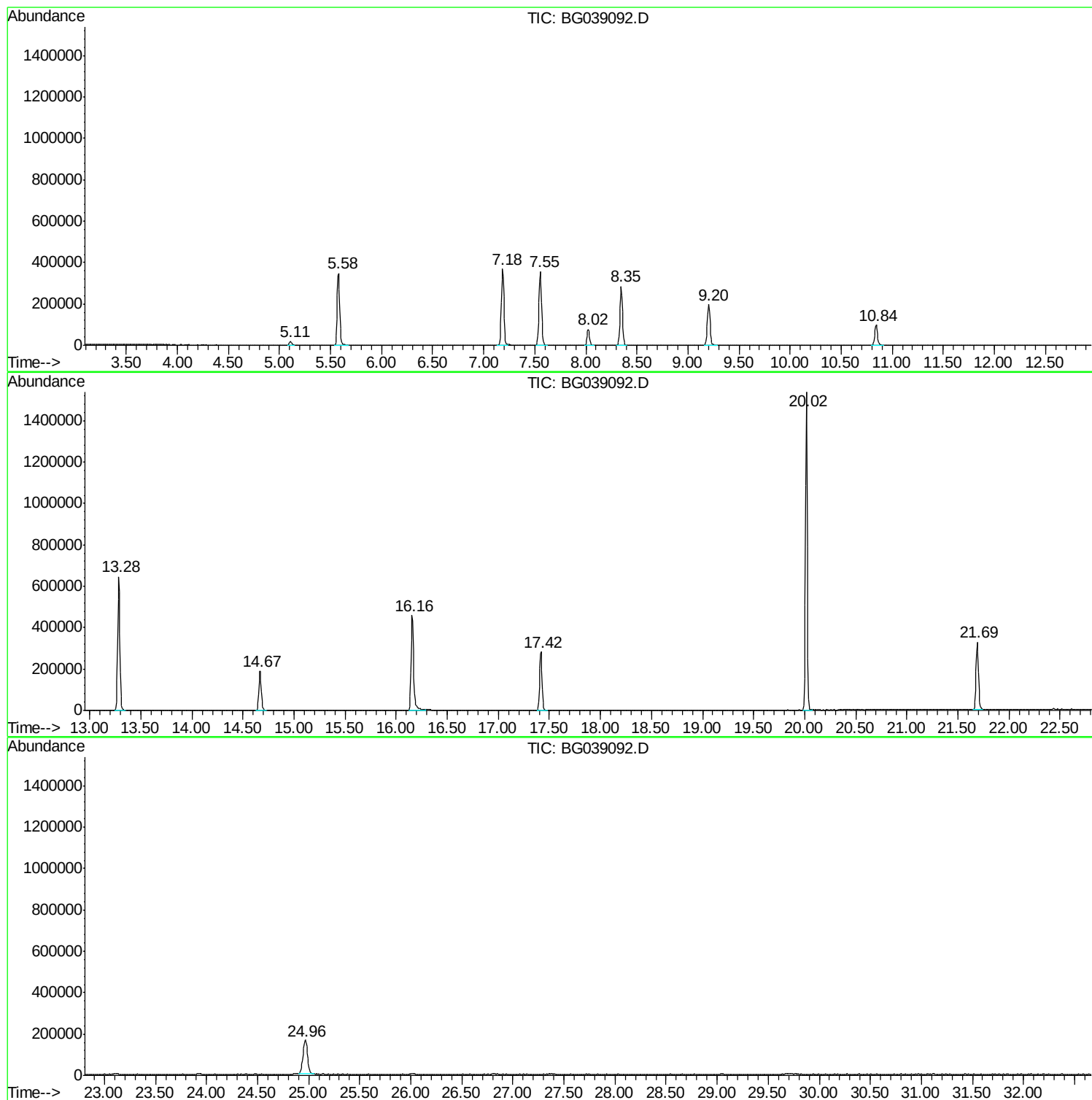
Sum of corrected areas: 8805312

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.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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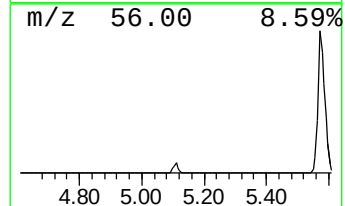
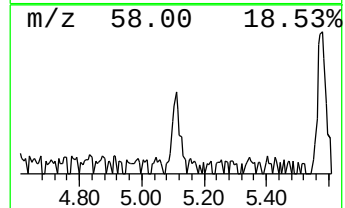
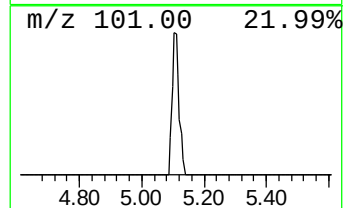
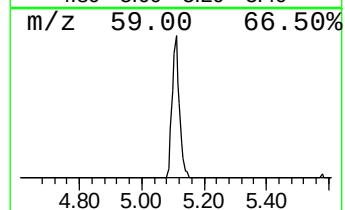
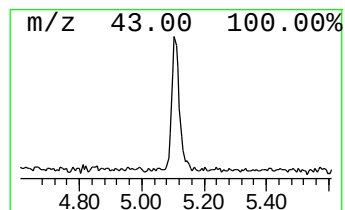
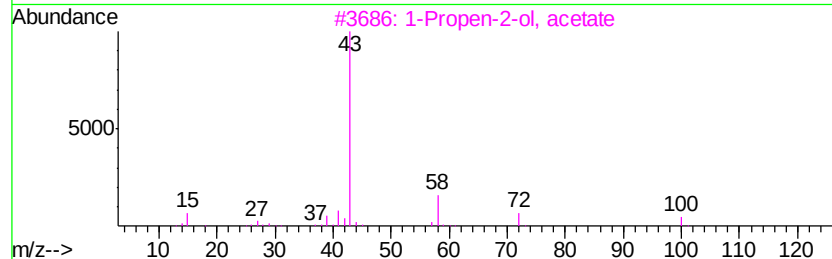
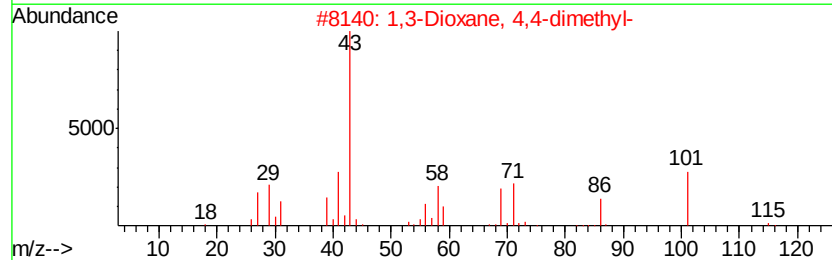
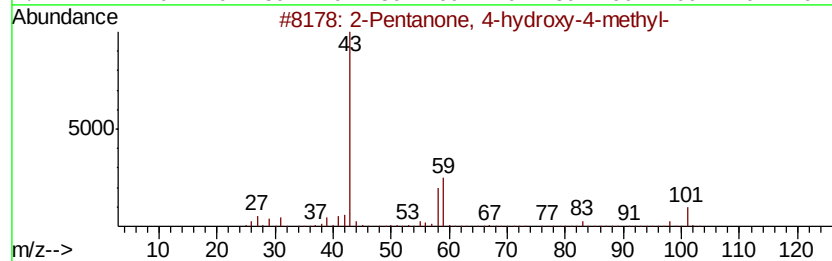
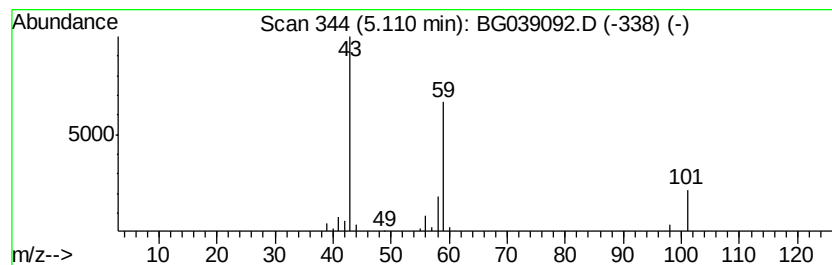
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	4.76 ng	31756	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		1,3-Dioxane, 4,4-dimethyl-	116	C6H12O2	000766-15-4	9
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	9
5		1,2-Butanediol	90	C4H10O2	000584-03-2	9



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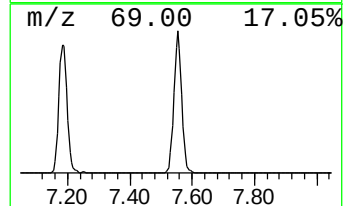
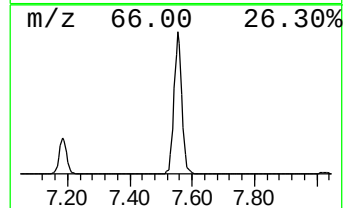
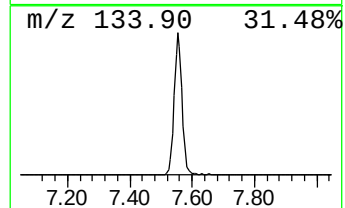
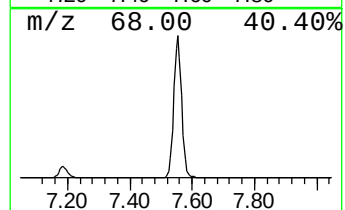
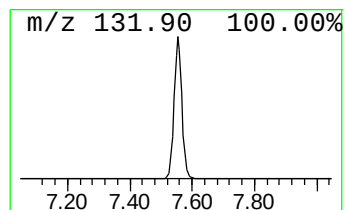
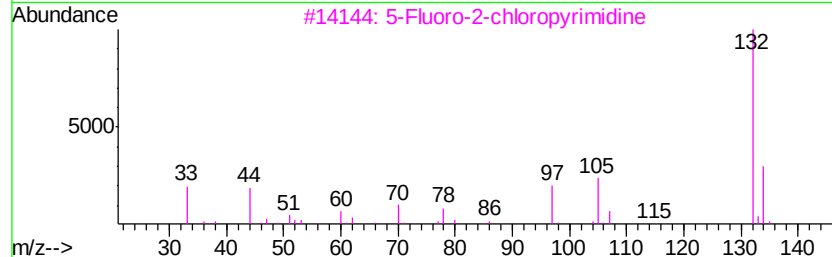
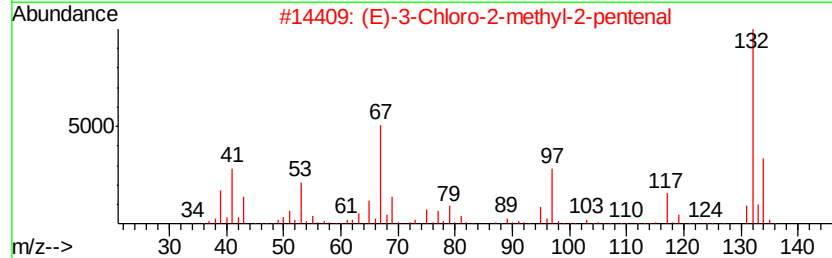
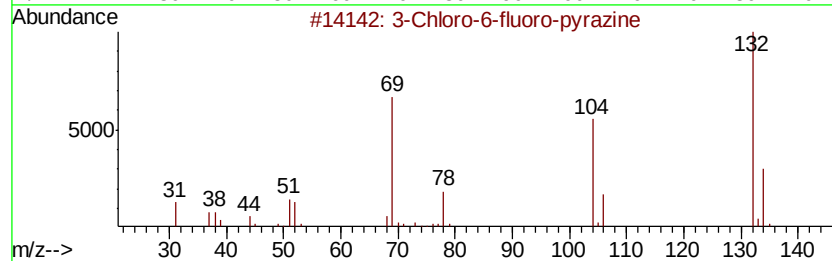
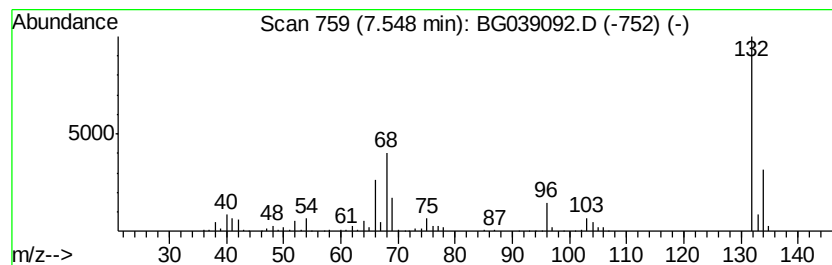
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Peak Number 2 unknown7.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.55	92.63 ng	617763	1,4-Dichlorobenzene-d4	8.02

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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.11	4.8	ng	31756	1	8.02	133379	20.0
unknown7.55	7.55	92.6	ng	617763	1	8.02	133379	20.0