

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101918\
Data File : BG037470.D
Acq On : 20 Oct 2018 3:40
Operator : JU/SJ
Sample : PB113548BL
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB113548BL

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-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	5.186	317	323	333	rVB	155298	245770	5.39%	1.230%
2	5.644	394	401	413	rBV	724252	1192356	26.15%	5.969%
3	7.248	666	674	685	rBV	766318	1338509	29.35%	6.701%
4	7.630	731	739	749	rBV	688688	1238007	27.15%	6.198%
5	8.106	813	820	828	rBV	190197	332144	7.28%	1.663%
6	8.429	866	875	886	rBV	744038	1339365	29.37%	6.705%
7	9.287	1008	1021	1033	rBV	625776	1168003	25.61%	5.847%
8	10.932	1291	1301	1309	rBV	271623	496664	10.89%	2.486%
9	13.364	1708	1715	1723	rBV	1808999	2877960	63.11%	14.408%
10	14.739	1942	1949	1958	rBV2	462173	765845	16.79%	3.834%
11	16.226	2195	2202	2211	rBV	899555	1408404	30.89%	7.051%
12	17.489	2409	2417	2429	rBV	639772	1023940	22.45%	5.126%
13	20.086	2853	2859	2871	rBV	3246388	4559997	100.00%	22.829%
14	21.772	3139	3146	3155	rBV2	599462	1054557	23.13%	5.279%
15	25.104	3703	3713	3730	rBV	281488	933044	20.46%	4.671%

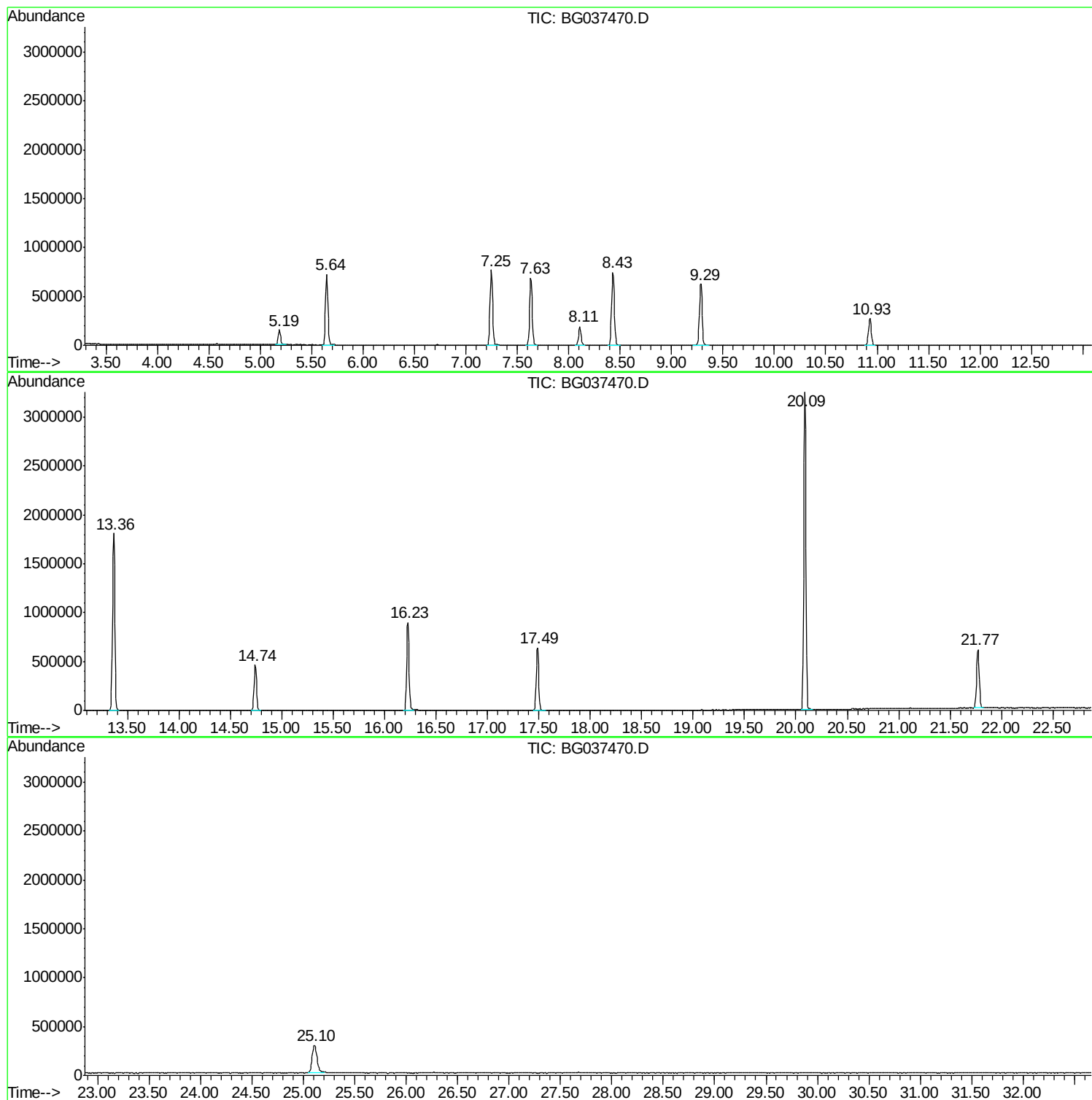
Sum of corrected areas: 19974565

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.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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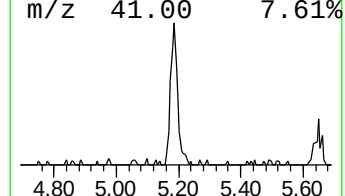
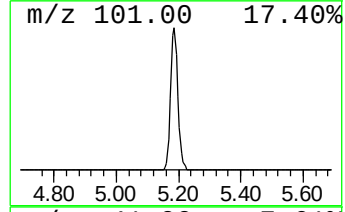
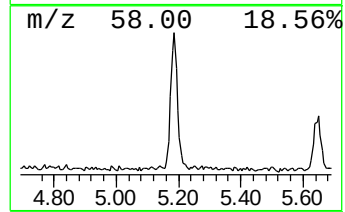
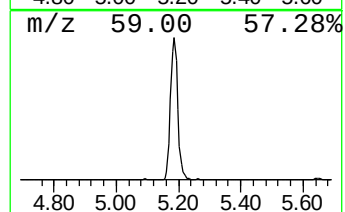
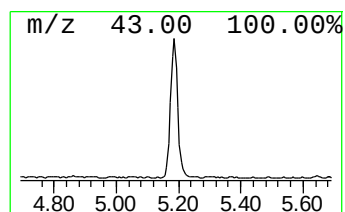
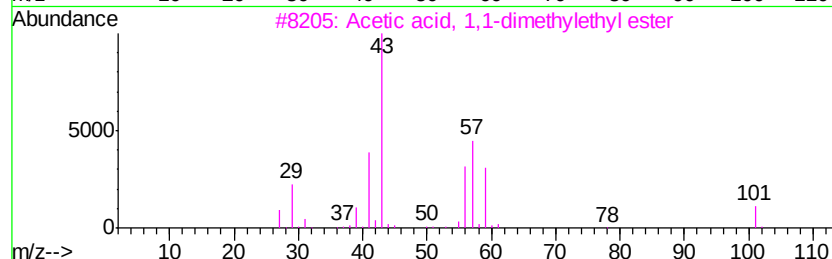
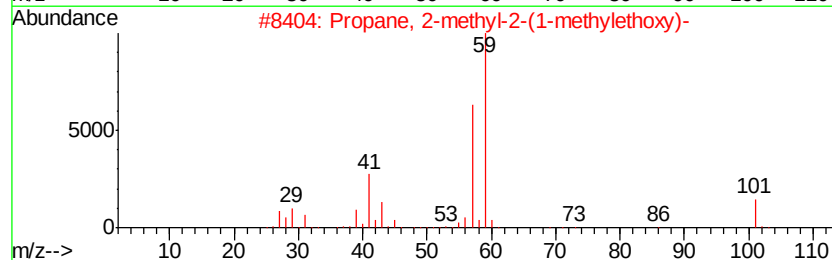
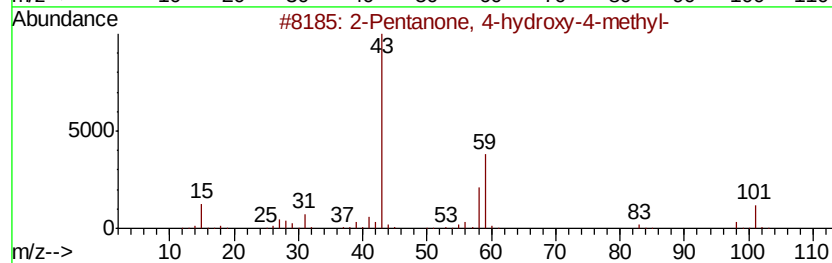
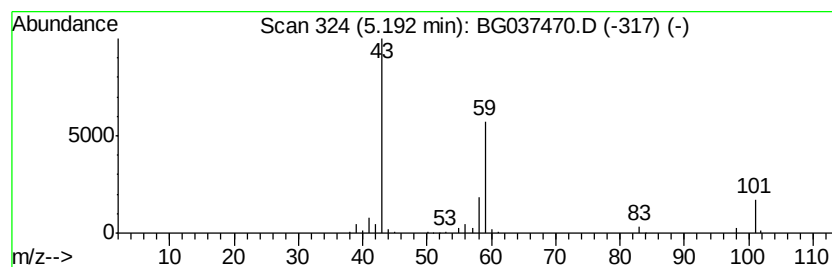
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	14.80 ng	245770	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	78
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	40
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	25
4			Acetone	58	C3H6O	000067-64-1	22
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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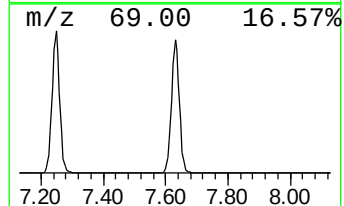
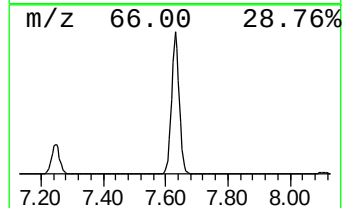
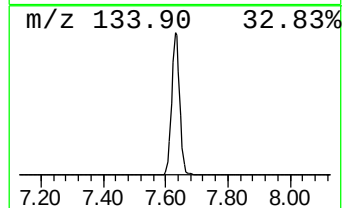
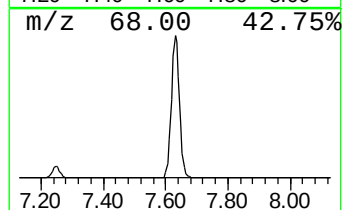
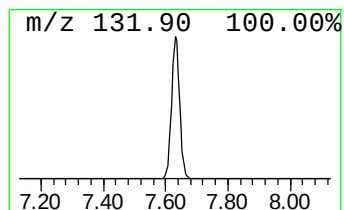
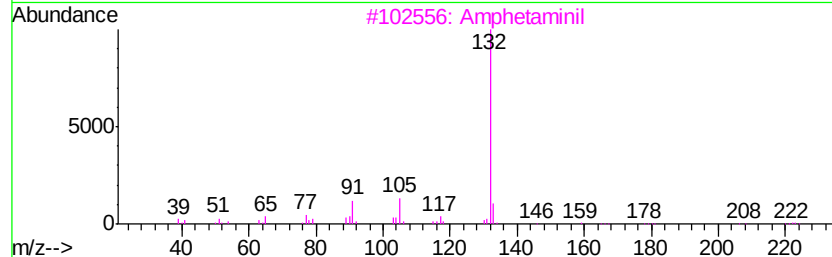
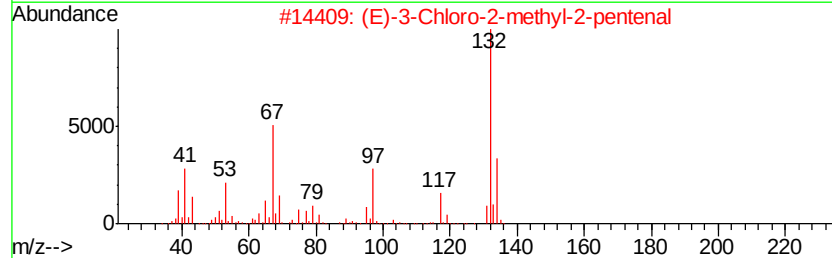
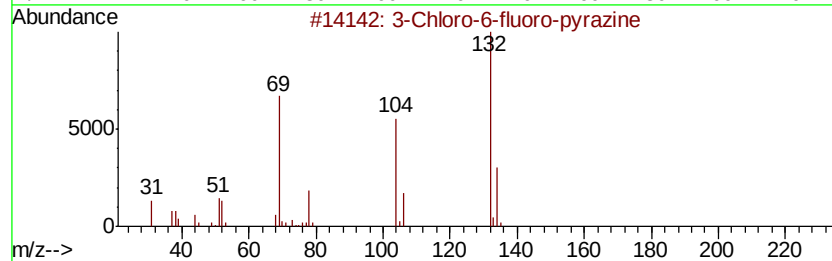
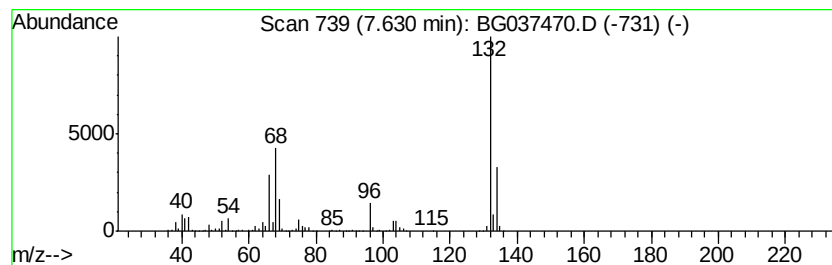
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Peak Number 2 unknown7.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	74.55 ng	1238010	1,4-Dichlorobenzene-d4	8.11

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	27
3		Amphetaminil	250	C17H18N2	017590-01-1	10
4		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	14.8	ng	245770	1	8.11	332144	20.0
unknown7.63	7.63	74.5	ng	1238010	1	8.11	332144	20.0