

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080618\
 Data File : BG036157.D
 Acq On : 7 Aug 2018 11:09
 Operator : JU/SJ
 Sample : J4335-03
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CINNAMINSON-8-3-18

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-BG071218.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	5.126	308	313	327	rVB	121102	193801	5.36%	1.362%
2	5.596	386	393	408	rBV	565900	973292	26.91%	6.841%
3	7.182	655	663	677	rBV	488766	938038	25.94%	6.593%
4	7.547	716	725	736	rBV	651051	1108689	30.66%	7.792%
5	8.011	796	804	811	rBV	108589	191553	5.30%	1.346%
6	8.334	851	859	870	rVB	476321	865797	23.94%	6.085%
7	9.174	994	1002	1014	rBV	362516	720367	19.92%	5.063%
8	10.813	1273	1281	1293	rBV2	100540	225813	6.24%	1.587%
9	13.256	1690	1697	1711	rBV	1064595	1750120	48.39%	12.301%
10	14.637	1925	1932	1948	rBV2	189470	358705	9.92%	2.521%
11	16.123	2178	2185	2198	rBV	911057	1557761	43.07%	10.949%
12	17.380	2393	2399	2410	rBV2	282733	520393	14.39%	3.658%
13	19.989	2836	2843	2855	rBV	2571516	3616411	100.00%	25.418%
14	21.639	3117	3124	3138	rBV	326545	645077	17.84%	4.534%
15	24.829	3656	3667	3677	rBV2	173895	562066	15.54%	3.950%

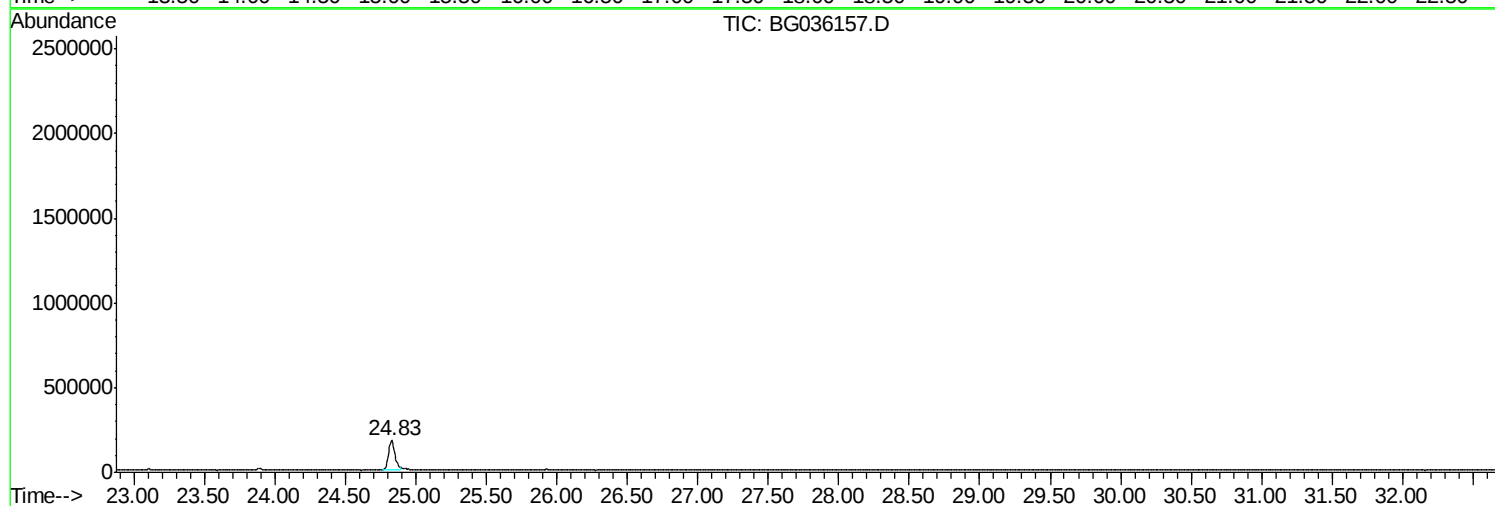
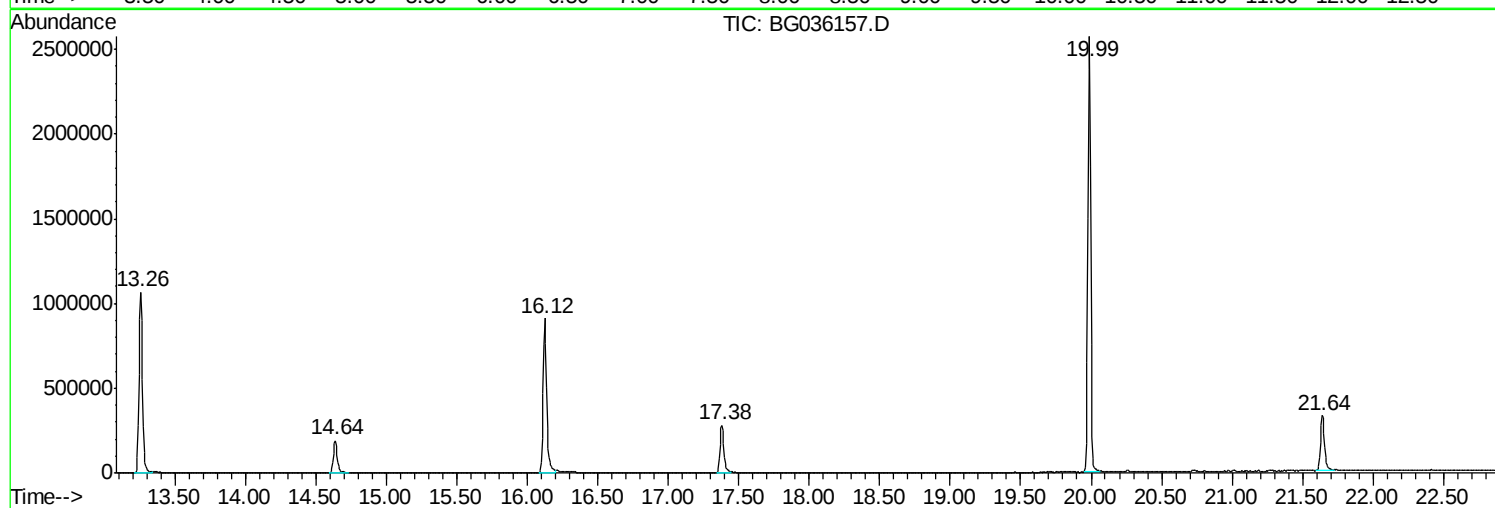
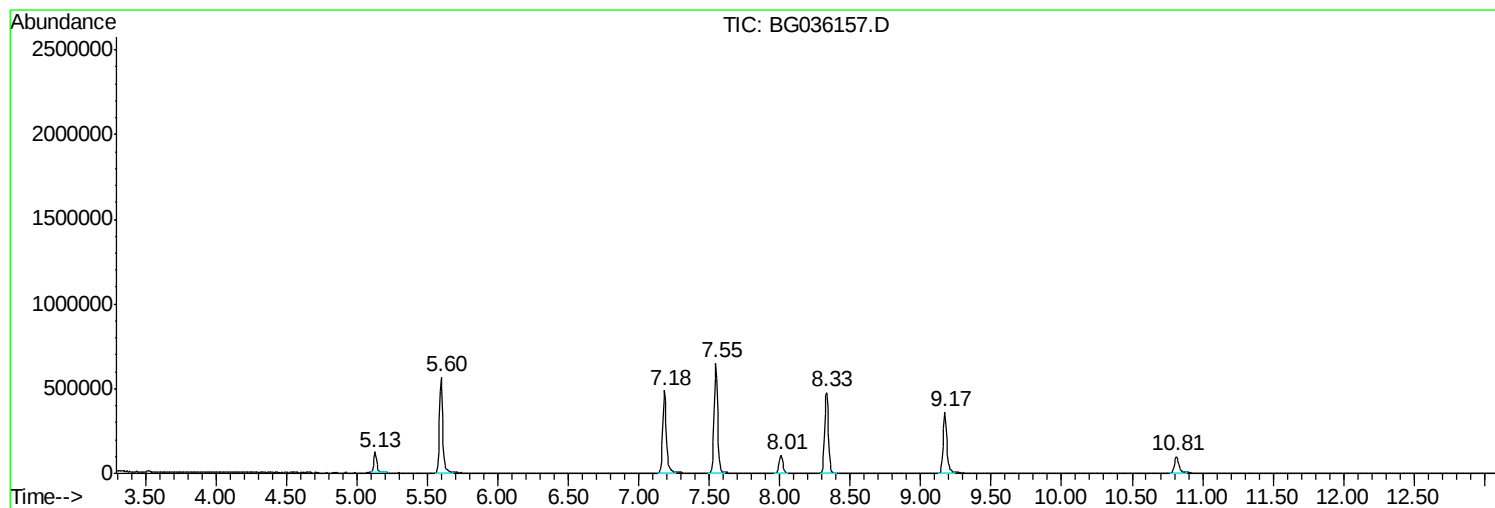
Sum of corrected areas: 14227883

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.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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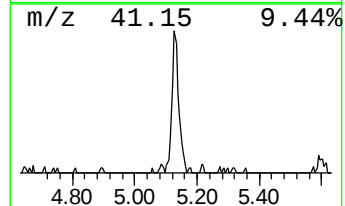
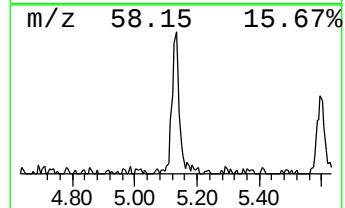
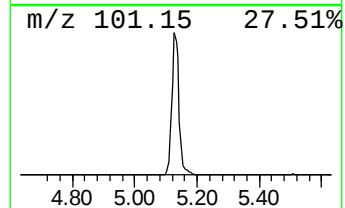
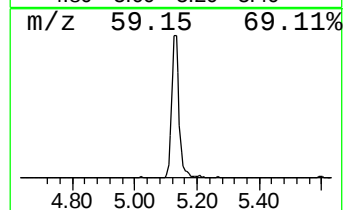
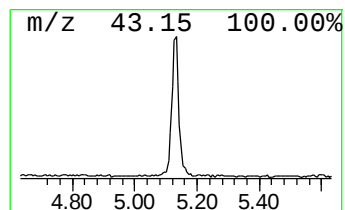
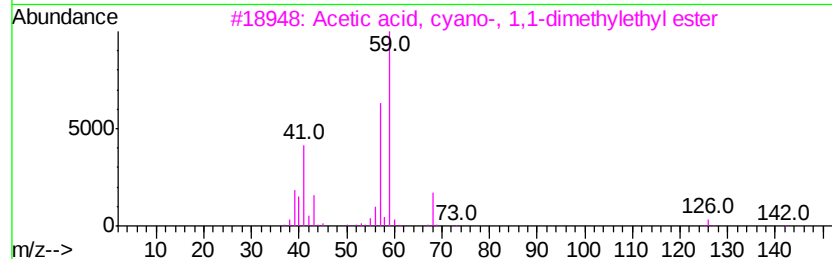
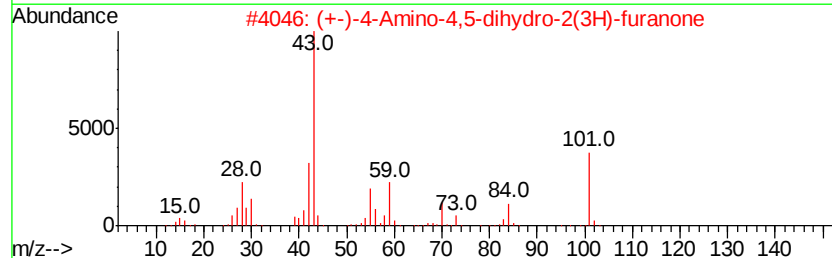
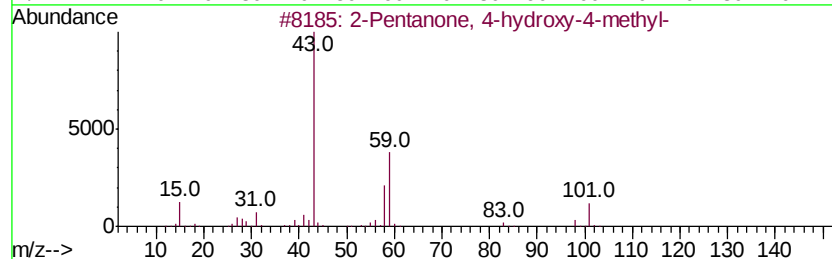
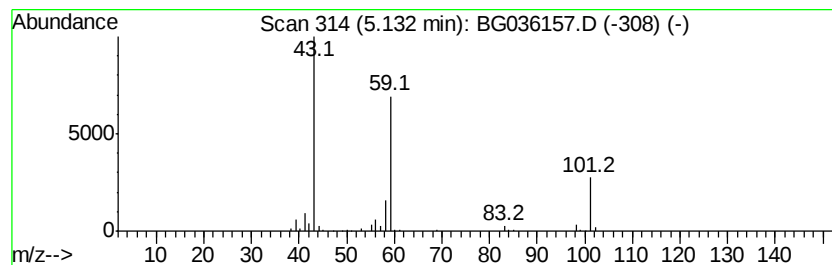
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	20.23 ng	193801	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	38
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	12
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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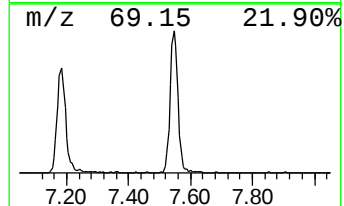
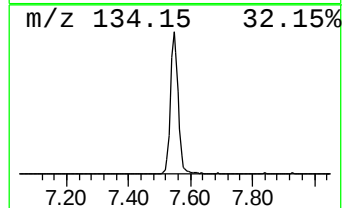
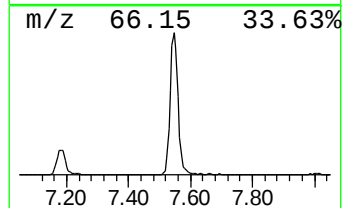
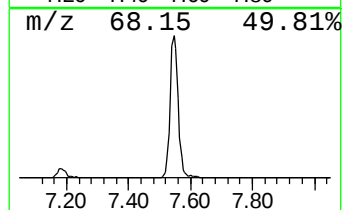
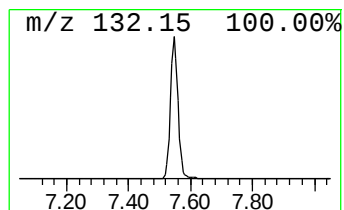
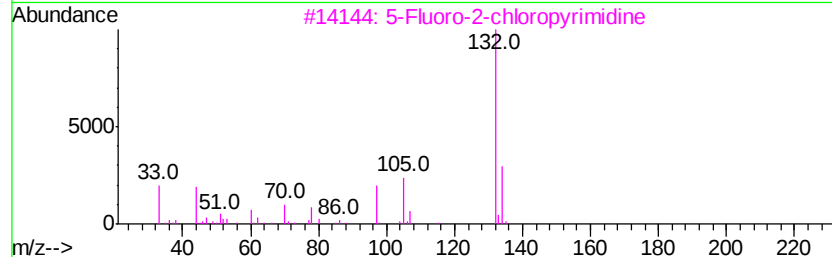
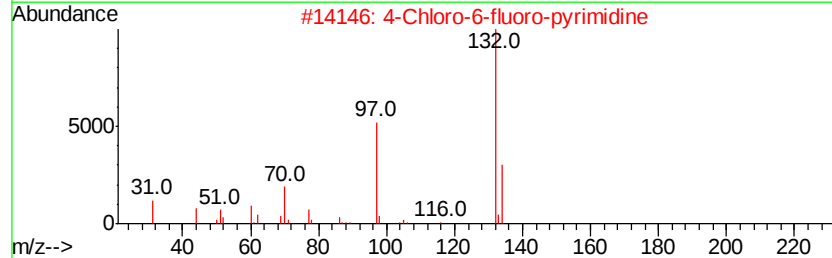
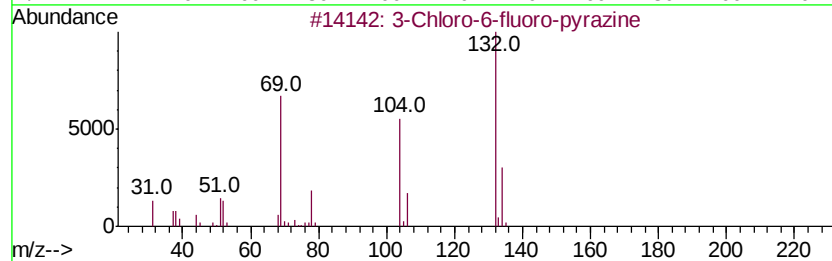
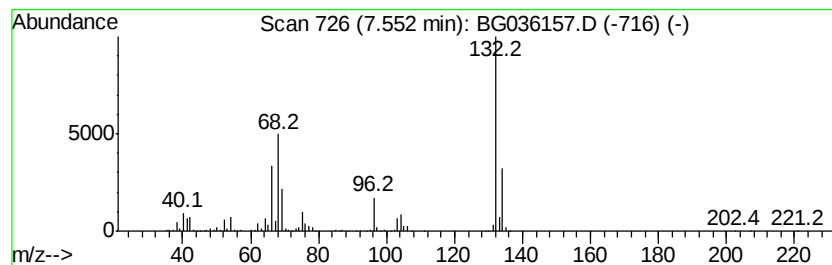
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Peak Number 2 unknown7.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.55	115.76 ng	1108690	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	20.2	ng	193801	1	8.01	191553	20.0
unknown7.55	7.55	115.8	ng	1108690	1	8.01	191553	20.0