

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112918\
Data File : BG038237.D
Acq On : 29 Nov 2018 17:43
Operator : JU/SJ
Sample : PB115157BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB115157BL

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-BG111318.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	5.165	313	319	326	rBV	26729	46020	2.04%	0.391%
2	5.629	391	398	412	rBV	570316	959637	42.64%	8.153%
3	7.233	663	671	686	rBV	613754	1096039	48.70%	9.312%
4	7.609	727	735	747	rBV	574742	1017189	45.20%	8.642%
5	8.085	808	816	825	rVB	115841	205881	9.15%	1.749%
6	8.408	863	871	882	rVB	409362	745872	33.14%	6.337%
7	9.260	1008	1016	1030	rBV	369271	680073	30.22%	5.778%
8	10.899	1288	1295	1305	rBV	163370	310486	13.80%	2.638%
9	13.338	1703	1710	1720	rBV	978928	1574593	69.97%	13.377%
10	14.718	1935	1945	1953	rBV2	273081	446002	19.82%	3.789%
11	16.205	2192	2198	2210	rBV3	537809	943267	41.91%	8.014%
12	17.468	2406	2413	2423	rBV2	327152	528127	23.47%	4.487%
13	20.065	2849	2855	2866	rBV	1665566	2250482	100.00%	19.119%
14	21.751	3135	3142	3154	rBV	301791	566054	25.15%	4.809%
15	25.077	3696	3708	3725	rBV2	106389	400988	17.82%	3.407%

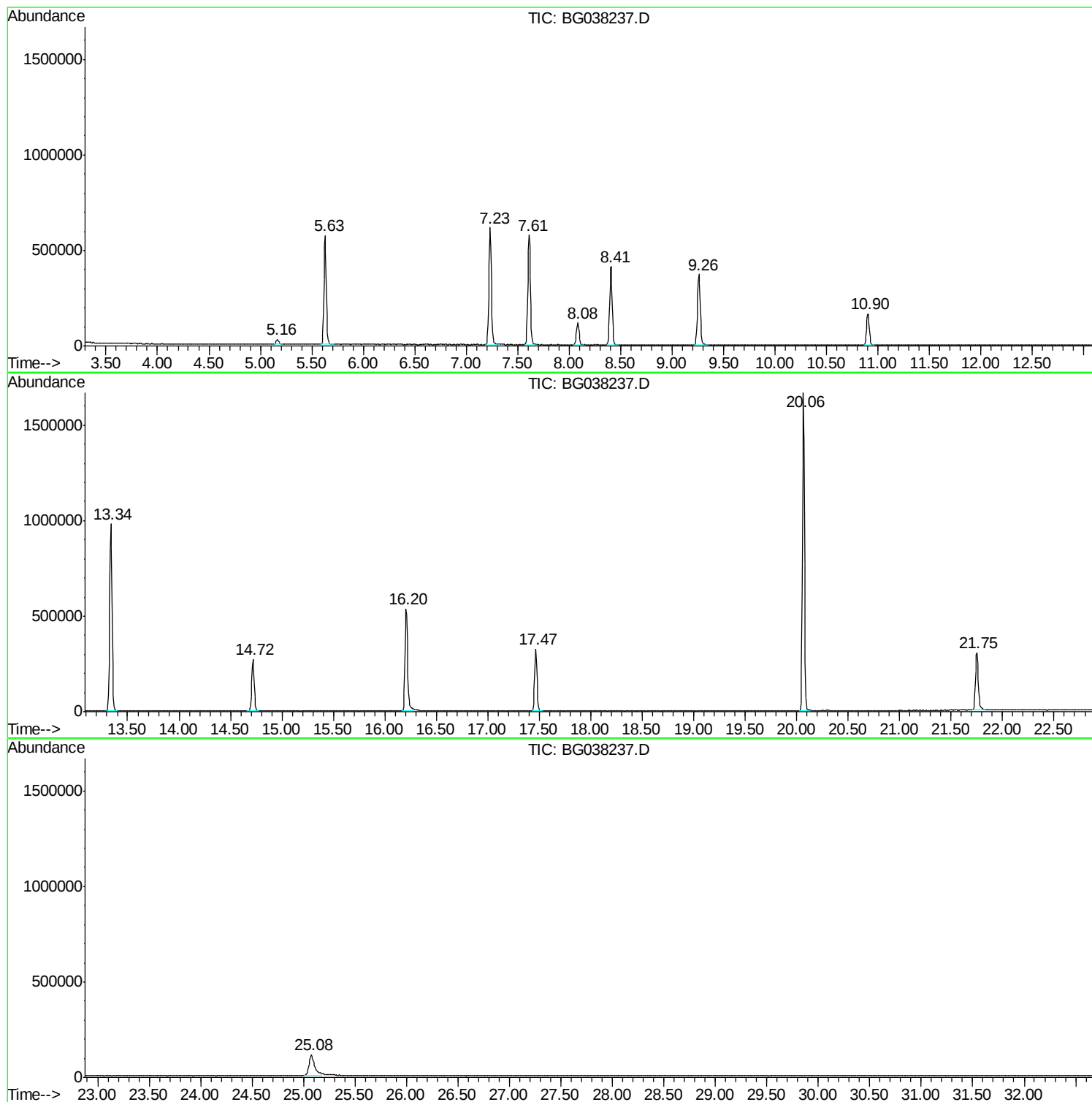
Sum of corrected areas: 11770710

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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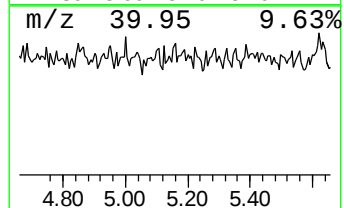
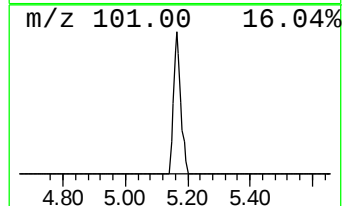
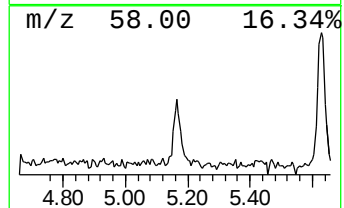
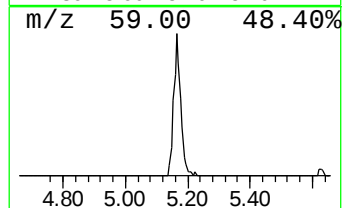
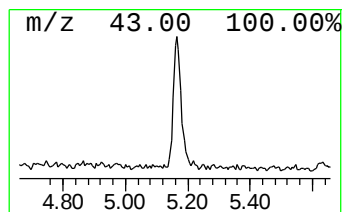
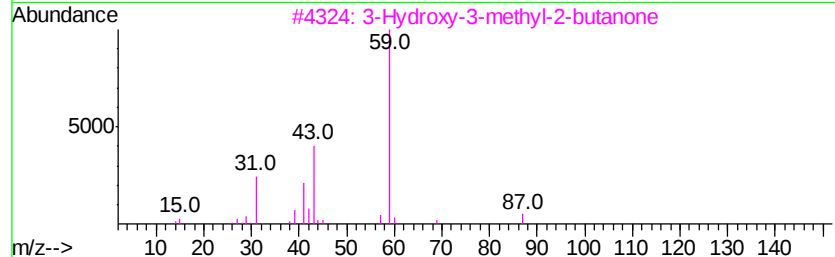
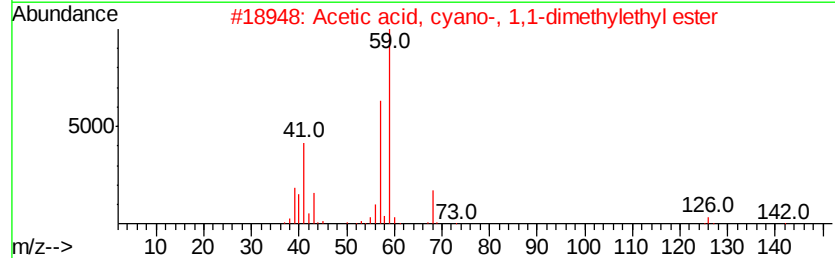
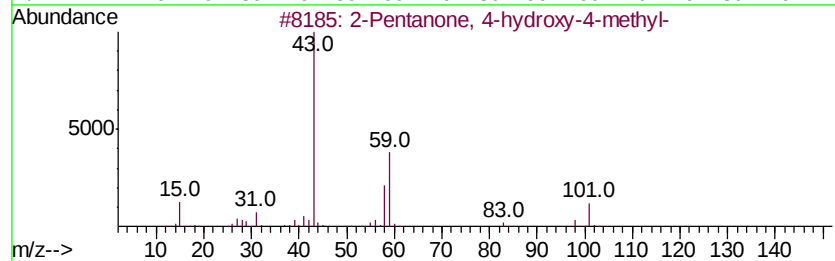
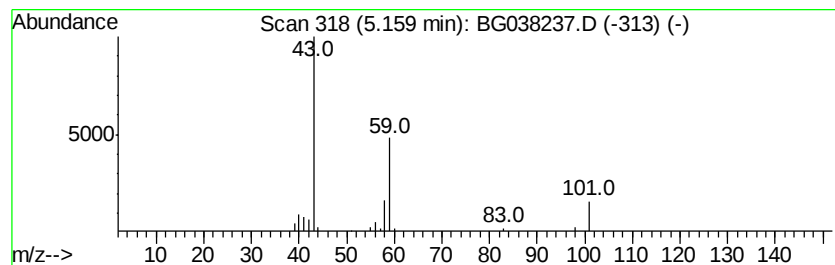
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	4.47 ng	46020	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	23
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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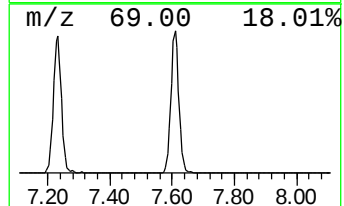
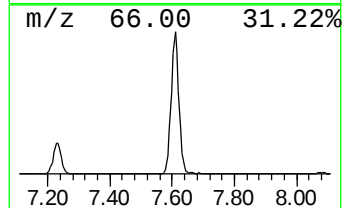
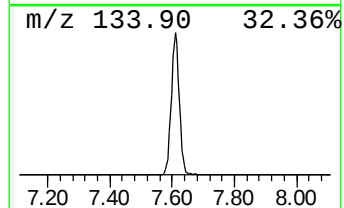
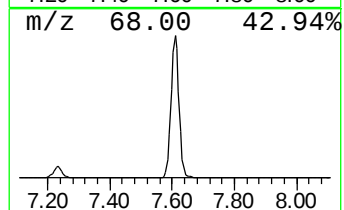
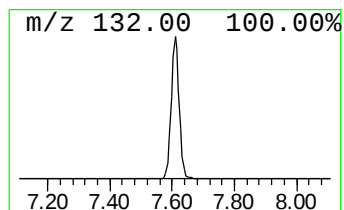
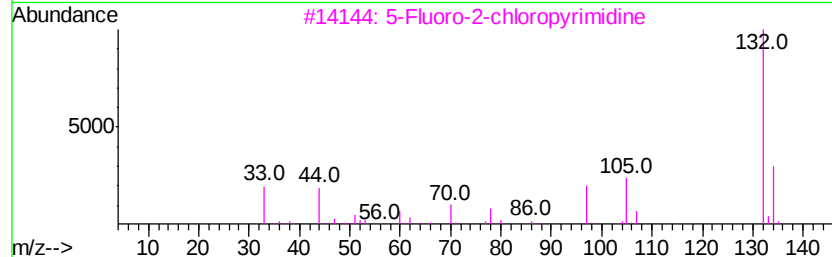
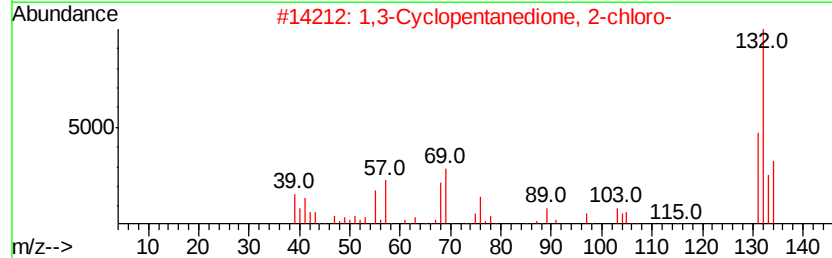
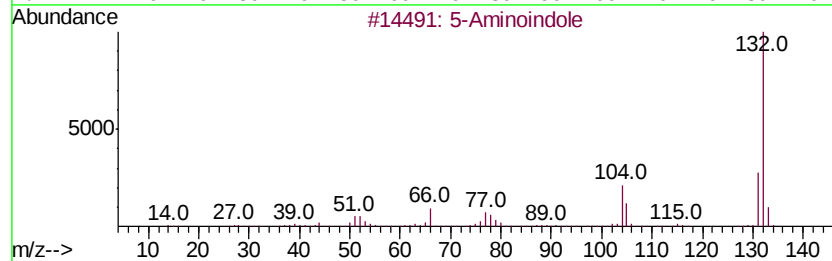
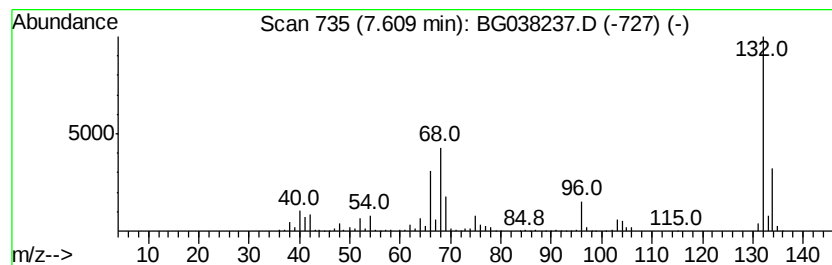
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	98.81 ng	1017190	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	22
2		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	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...	5.16	4.5	ng	46020	1	8.08	205881	20.0
unknown7.61	7.61	98.8	ng	1017190	1	8.08	205881	20.0