

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112018\
 Data File : BG038097.D
 Acq On : 20 Nov 2018 22:44
 Operator : JU/SJ
 Sample : PB114966BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114966BL

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	5.163	314	319	329	rBV2	29416	50254	2.04%	0.382%
2	5.633	391	399	410	rBV	619590	1047375	42.56%	7.962%
3	7.231	663	671	680	rBV	653382	1146633	46.59%	8.717%
4	7.613	728	736	745	rBV	610762	1082572	43.99%	8.230%
5	8.083	810	816	827	rVB	134628	238466	9.69%	1.813%
6	8.406	864	871	879	rBV	458177	823927	33.48%	6.263%
7	9.264	1009	1017	1028	rBV	386603	713173	28.98%	5.421%
8	10.903	1289	1296	1304	rVB	190577	354979	14.42%	2.698%
9	13.342	1703	1711	1718	rBV	1068620	1675404	68.08%	12.736%
10	14.723	1938	1946	1957	rBV2	295471	500480	20.34%	3.805%
11	16.209	2192	2199	2216	rBV	685898	1105754	44.93%	8.406%
12	17.466	2405	2413	2424	rBV2	388872	616018	25.03%	4.683%
13	20.069	2850	2856	2864	rBV	1913909	2460881	100.00%	18.707%
14	21.750	3135	3142	3152	rBV2	399607	690072	28.04%	5.246%
15	25.075	3694	3708	3724	rBV	196078	648700	26.36%	4.931%

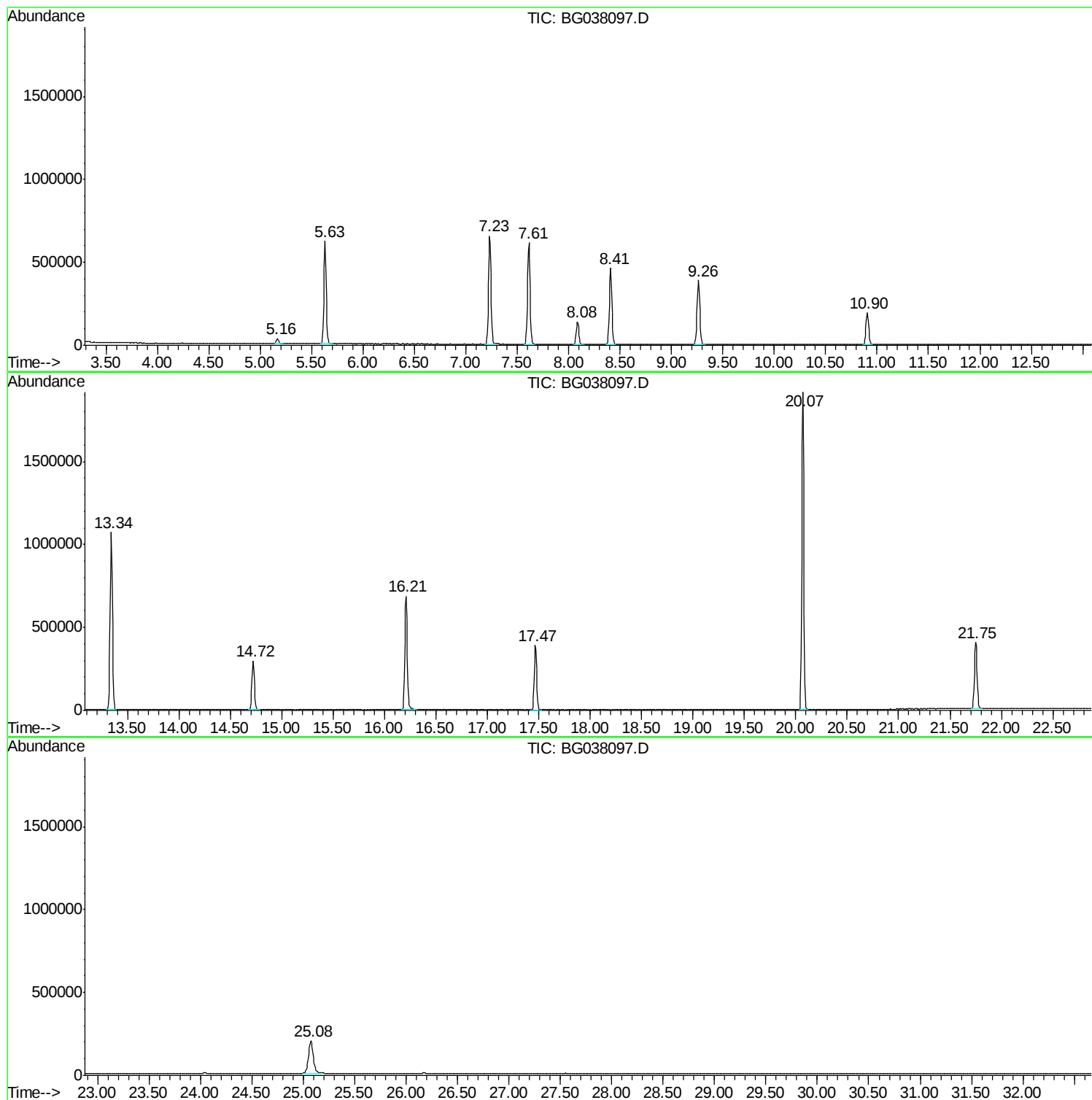
Sum of corrected areas: 13154688

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



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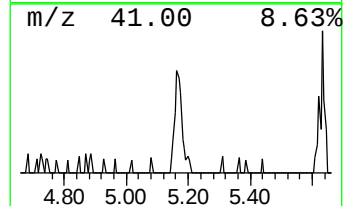
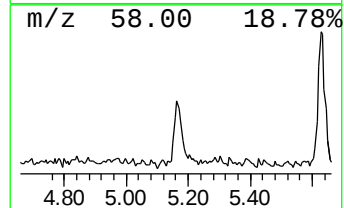
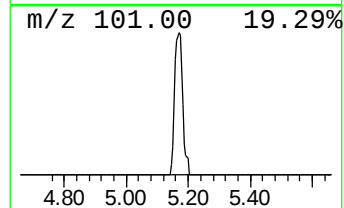
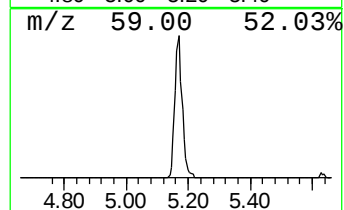
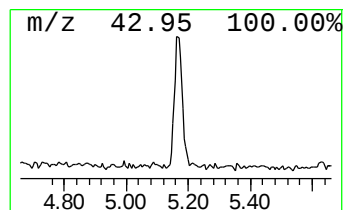
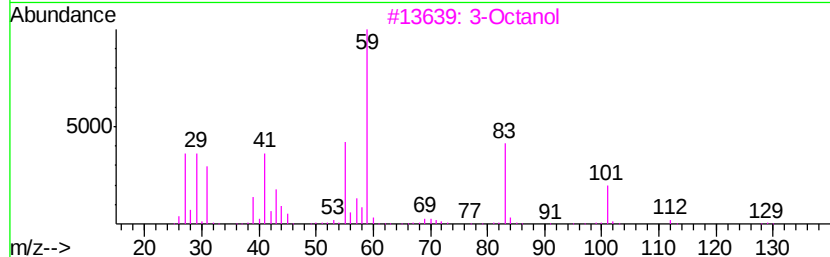
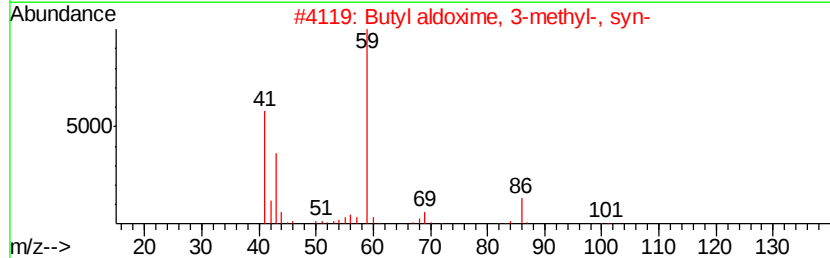
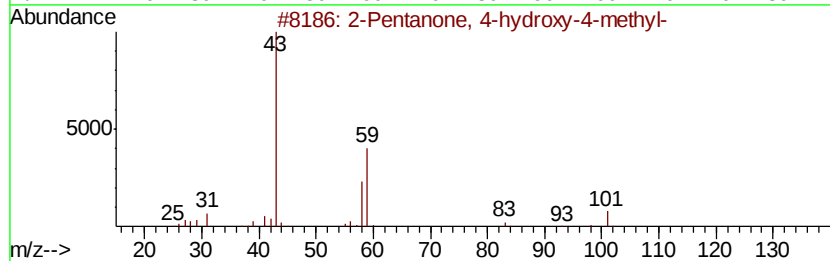
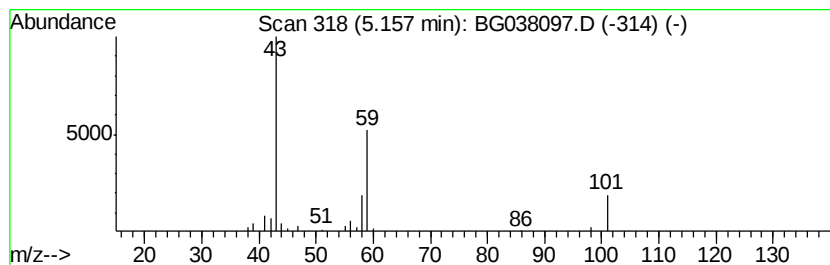
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	4.21 ng	50254	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	37
3			3-Octanol	130	C8H18O	000589-98-0	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
5			Propylamine	59	C3H9N	000107-10-8	9



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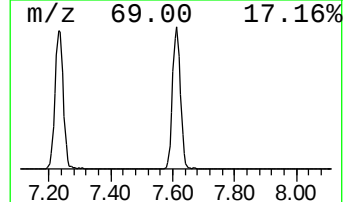
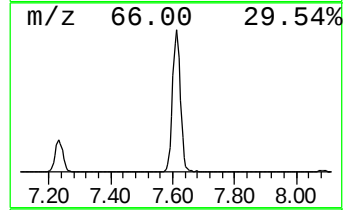
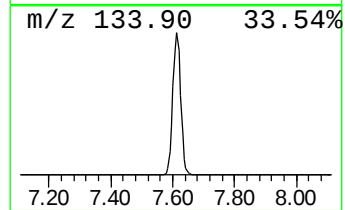
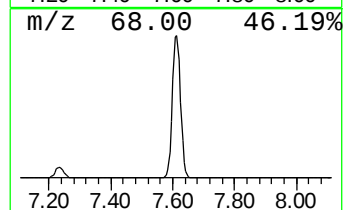
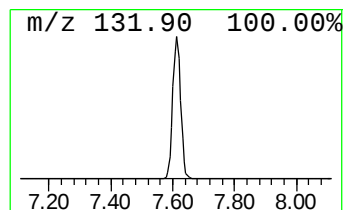
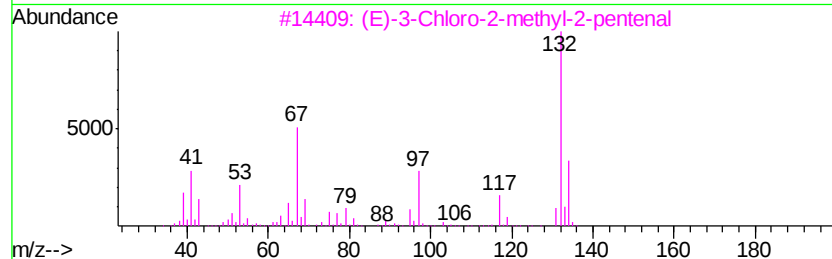
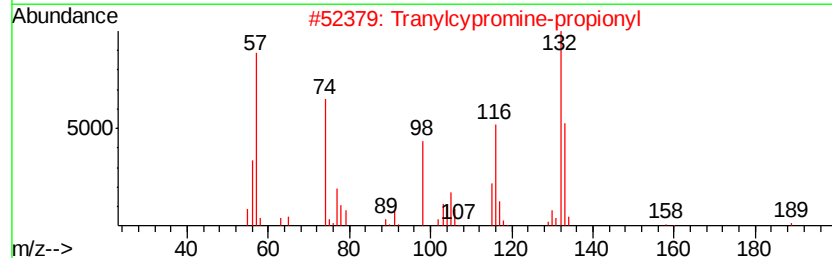
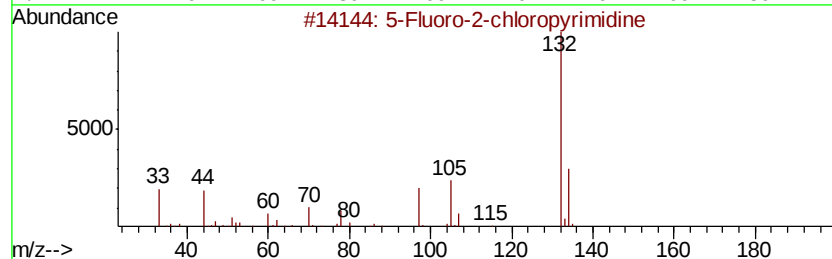
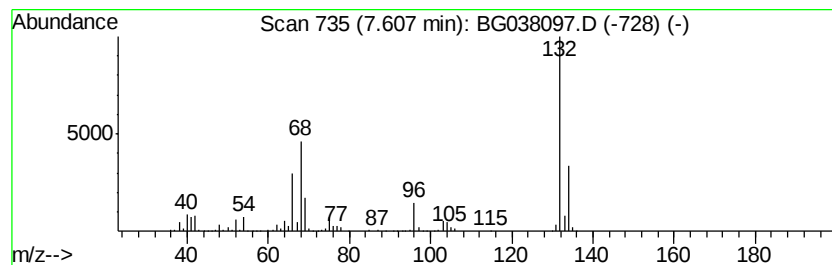
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	90.79 ng	1082570	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		5-Aminoindole	132	C8H8N2	005192-03-0	9
5		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	4.2	ng	50254	1	8.08	238466	20.0
unknown7.61	7.61	90.8	ng	1082570	1	8.08	238466	20.0