

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112018\
 Data File : BG038084.D
 Acq On : 20 Nov 2018 14:09
 Operator : JU/SJ
 Sample : PB114959BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114959BL

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	5.170	314	320	327	rBV	34872	57491	2.07%	0.394%
2	5.628	392	398	408	rBV	681140	1141903	41.14%	7.830%
3	7.232	664	671	685	rBV	718959	1272477	45.84%	8.725%
4	7.614	728	736	749	rVB	672458	1198267	43.17%	8.217%
5	8.084	809	816	824	rVB	141832	247314	8.91%	1.696%
6	8.407	864	871	881	rVB	490181	898569	32.37%	6.161%
7	9.265	1009	1017	1028	rBV	419422	791684	28.52%	5.429%
8	10.904	1289	1296	1304	rBV	203697	367759	13.25%	2.522%
9	13.343	1704	1711	1719	rBV	1163063	1900294	68.46%	13.030%
10	14.723	1938	1946	1954	rBV2	321010	541301	19.50%	3.712%
11	16.210	2192	2199	2213	rBV3	820510	1302907	46.94%	8.934%
12	17.467	2405	2413	2421	rBV2	422083	661474	23.83%	4.536%
13	20.070	2850	2856	2863	rBV	2080642	2775817	100.00%	19.034%
14	21.750	3135	3142	3154	rBV2	430946	737710	26.58%	5.058%
15	25.076	3694	3708	3723	rBV2	206703	688674	24.81%	4.722%

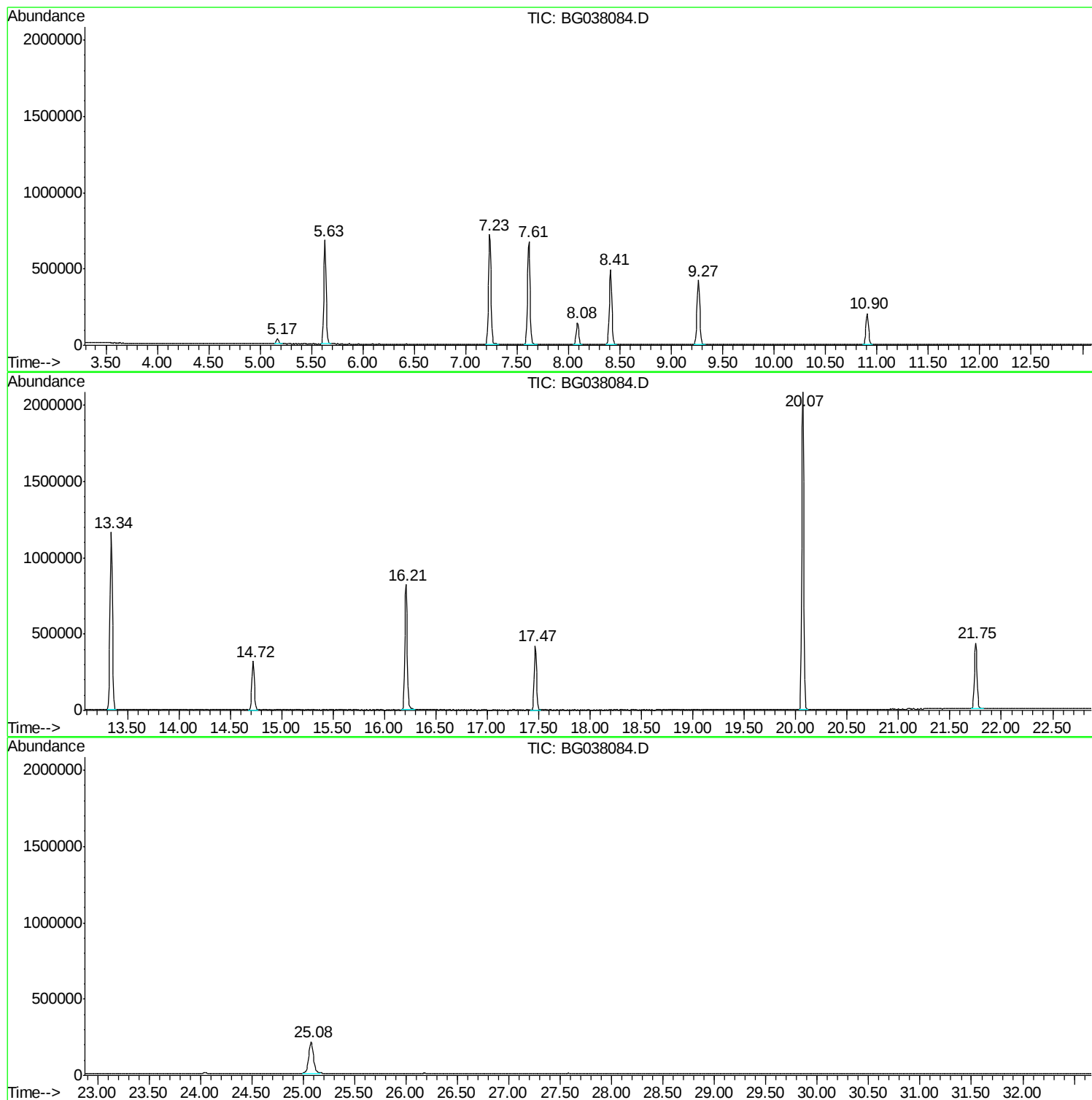
Sum of corrected areas: 14583641

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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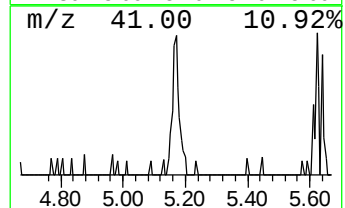
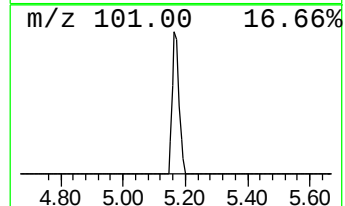
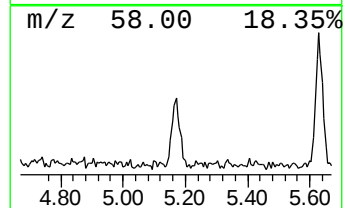
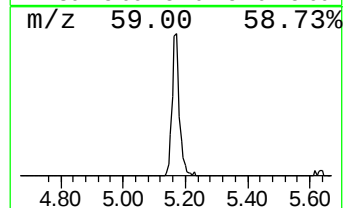
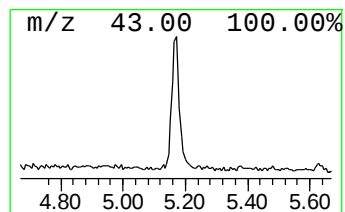
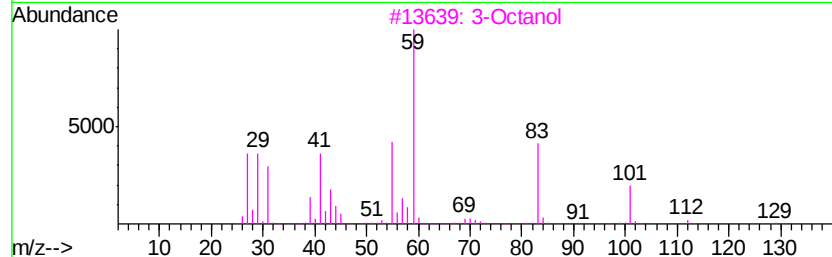
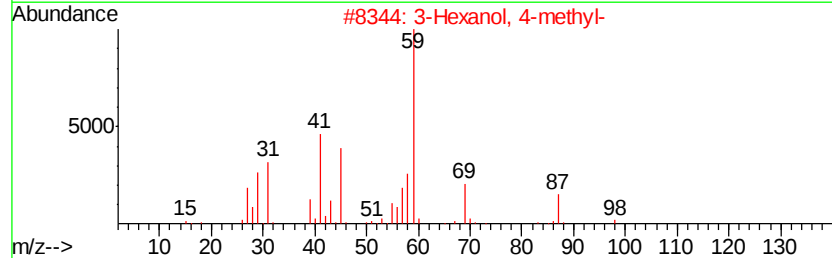
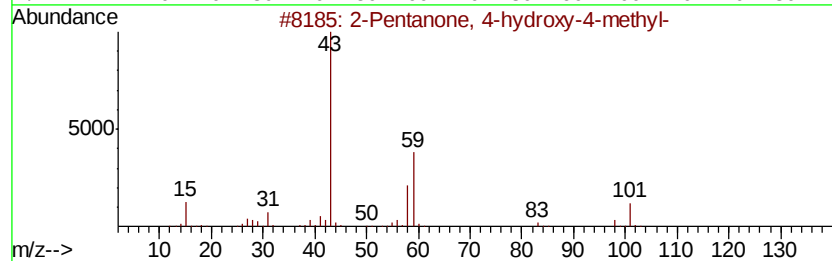
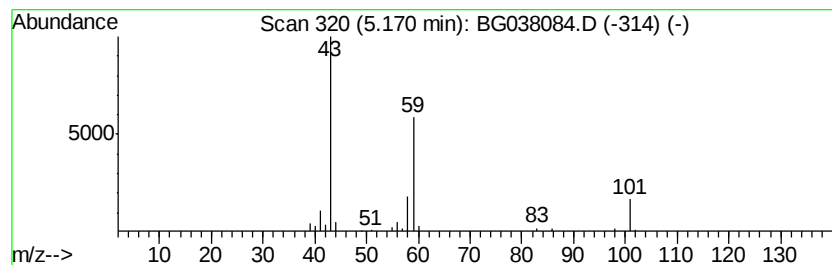
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	4.65 ng	57491	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	64
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	40
3			3-Octanol	130	C8H18O	000589-98-0	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5			3-Heptanol	116	C7H16O	000589-82-2	25



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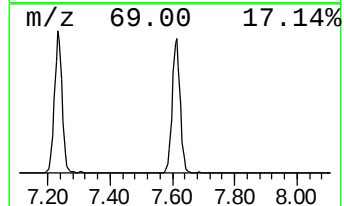
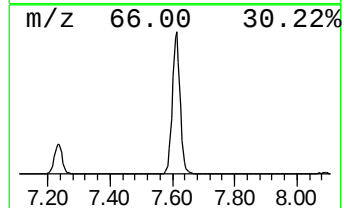
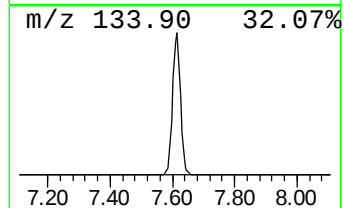
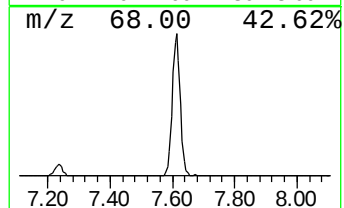
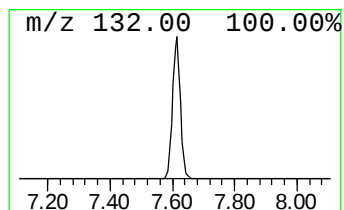
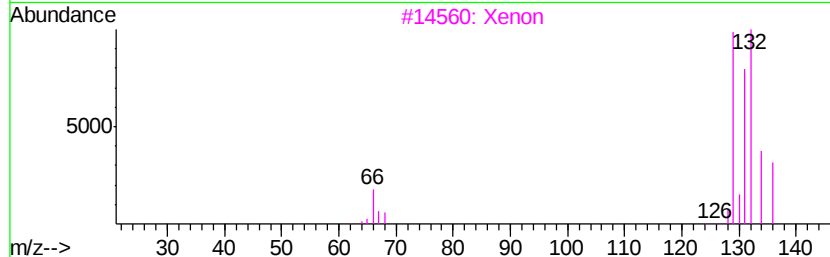
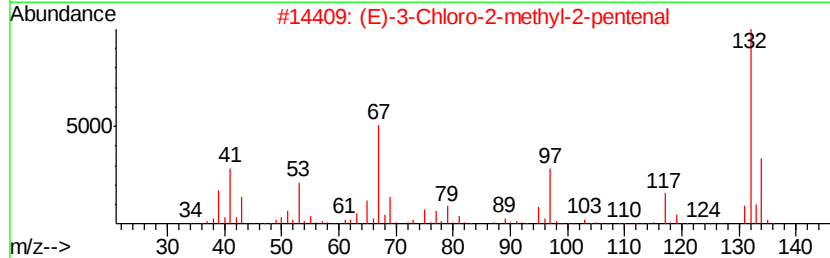
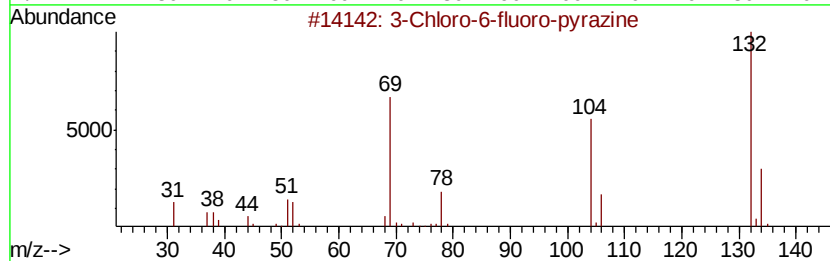
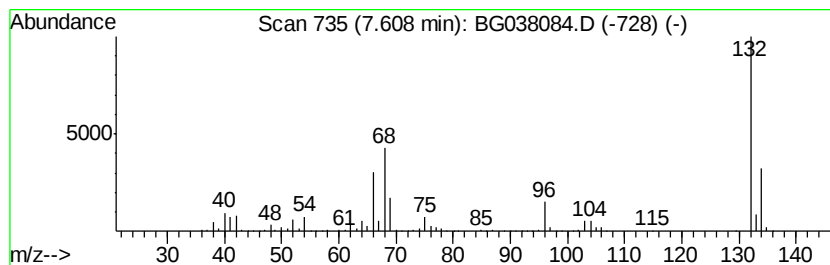
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	96.90 ng	1198270	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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			Xenon	132	Xe	007440-63-3	9
4			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	9
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...	5.17	4.7	ng	57491	1	8.08	247314	20.0
unknown7.61	7.61	96.9	ng	1198270	1	8.08	247314	20.0