

Data Path : Z:\HPCHEM1\BNA G\DATA\BG113017\
 Data File : BG031309.D
 Acq On : 1 Dec 2017 4:36
 Operator : SJ/JU
 Sample : I6635-07
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 112817-TIDO-E-U-B

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:\HPCHEM1\BNA_G\METHODS\8270-BG111017.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	5.203	318	326	340	rBV	60450	101599	1.68%	0.472%
2	5.684	402	408	421	rBV	303568	553854	9.14%	2.572%
3	7.294	675	682	695	rBV	178809	362134	5.97%	1.682%
4	7.659	736	744	756	rBV	561762	1054798	17.40%	4.899%
5	8.129	816	824	833	rBV	207337	366086	6.04%	1.700%
6	8.452	871	879	892	rBV	835827	1526666	25.19%	7.090%
7	9.298	1015	1023	1042	rBV	573105	1139457	18.80%	5.292%
8	10.949	1295	1304	1323	rBV	238609	521305	8.60%	2.421%
9	13.375	1710	1717	1730	rBV	1920997	3209664	52.96%	14.907%
10	14.756	1943	1952	1970	rBV2	455989	838399	13.83%	3.894%
11	16.237	2198	2204	2222	rBV	1004708	1726705	28.49%	8.019%
12	17.494	2411	2418	2432	rBV2	713479	1138618	18.79%	5.288%
13	20.085	2853	2859	2870	rBV	4546786	6061014	100.00%	28.149%
14	21.378	3074	3079	3093	rBV9	34696	92412	1.52%	0.429%
15	21.766	3138	3145	3161	rVB	832430	1461802	24.12%	6.789%
16	25.068	3696	3707	3722	rBV2	465962	1377423	22.73%	6.397%

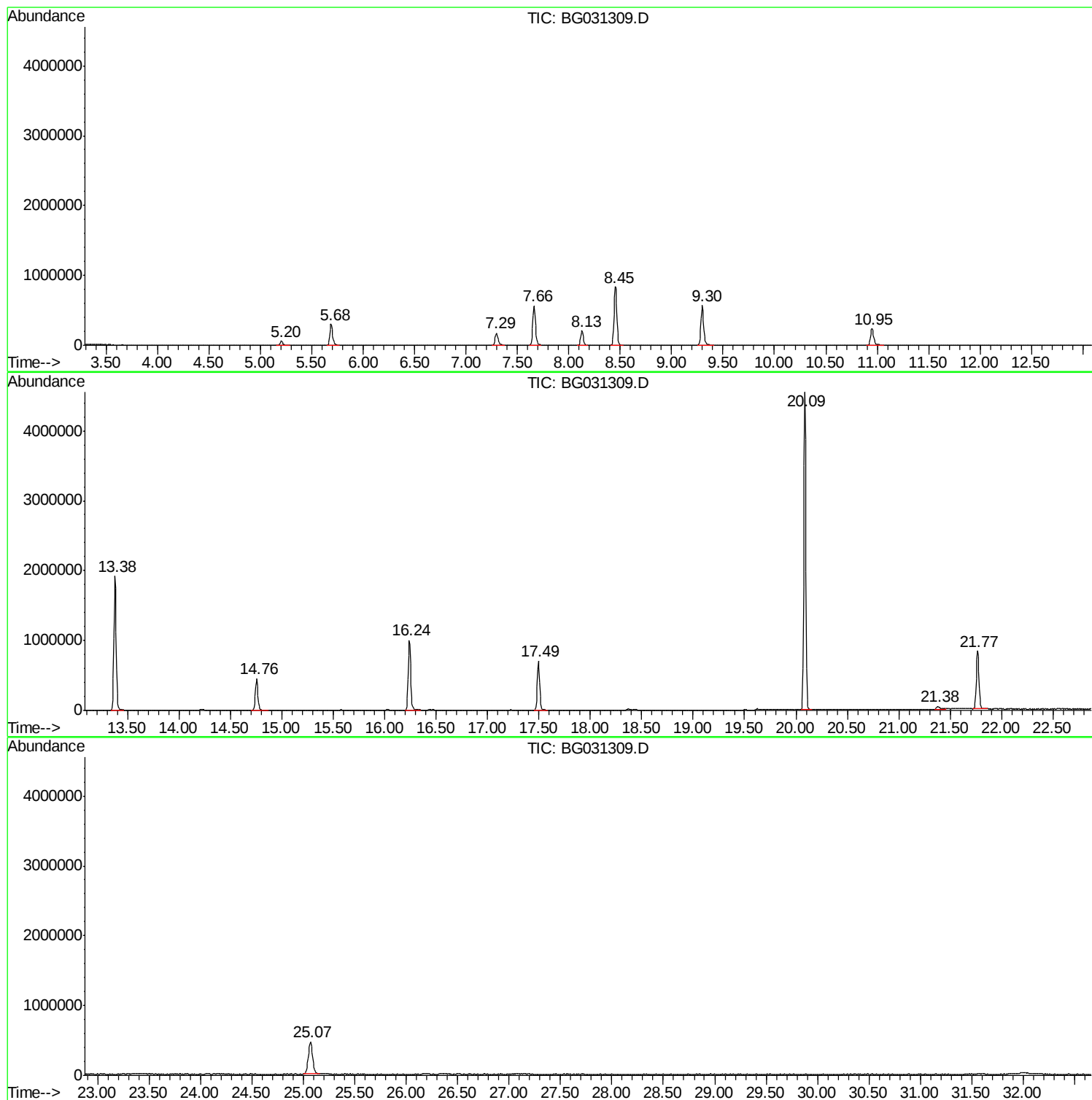
Sum of corrected areas: 21531936

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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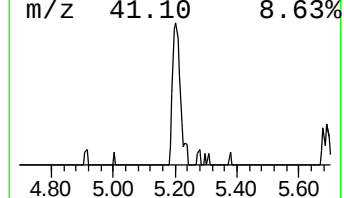
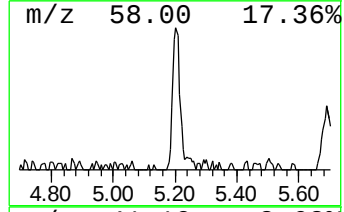
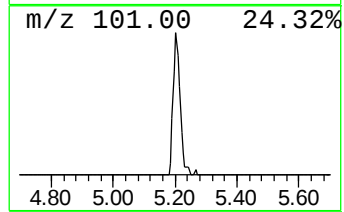
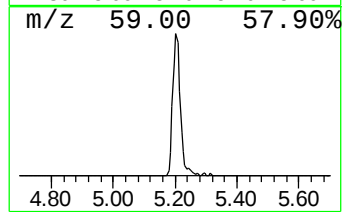
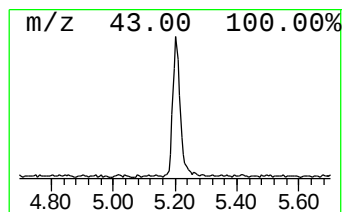
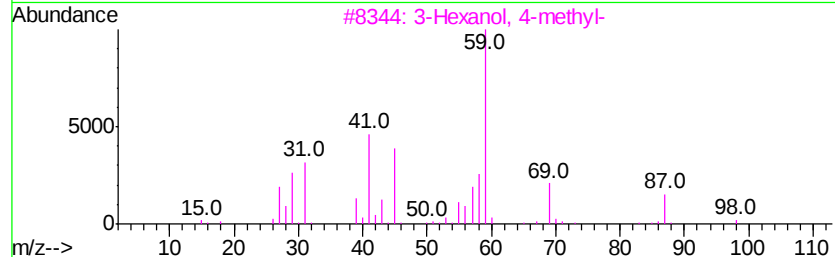
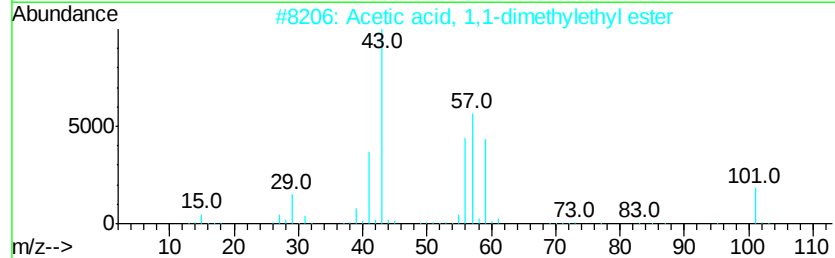
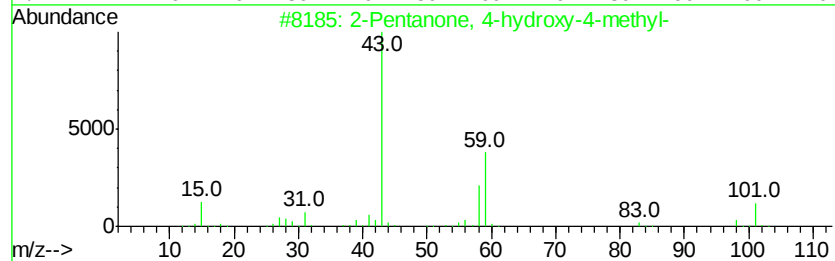
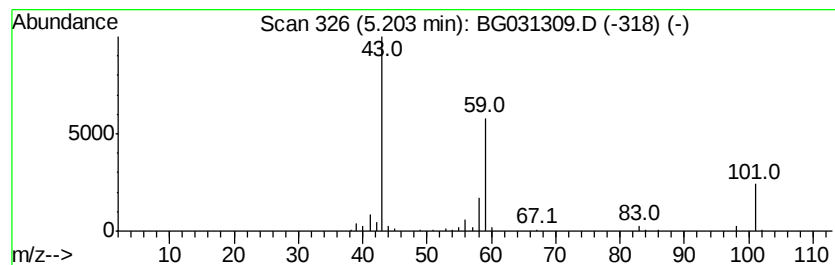
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	5.55 ng	101599	1,4-Dichlorobenzene-d4	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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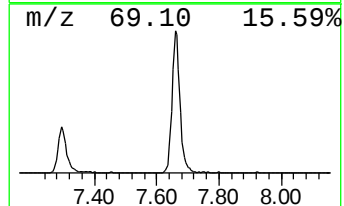
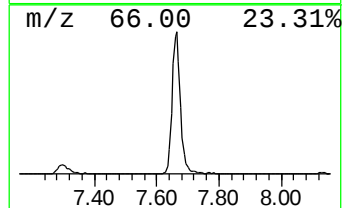
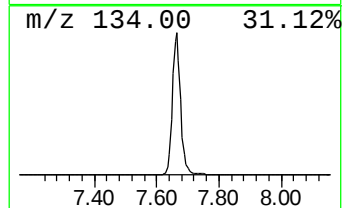
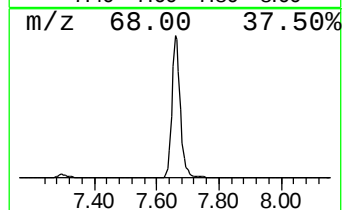
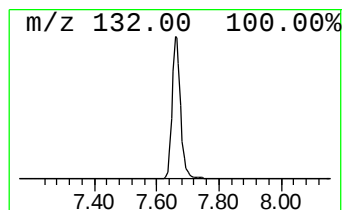
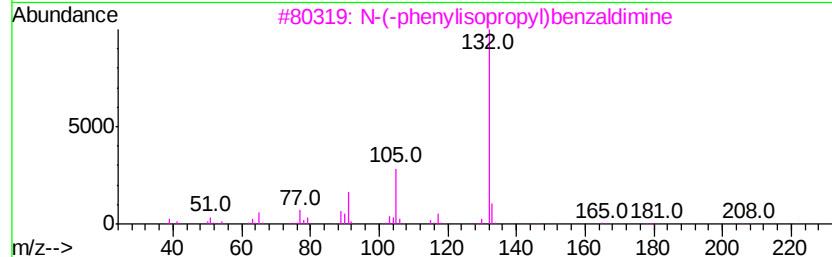
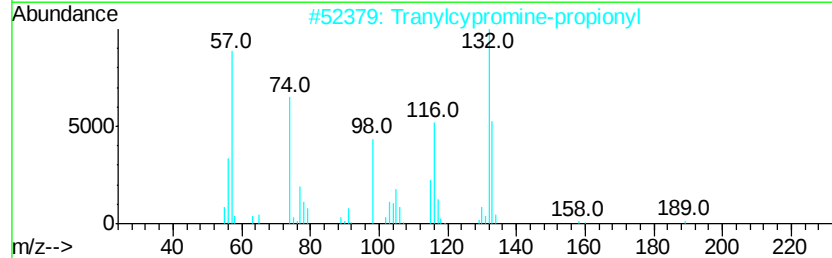
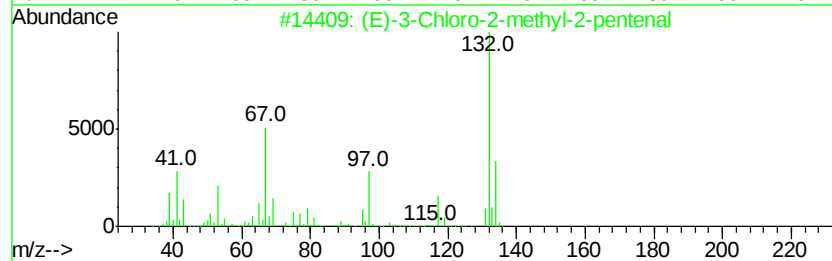
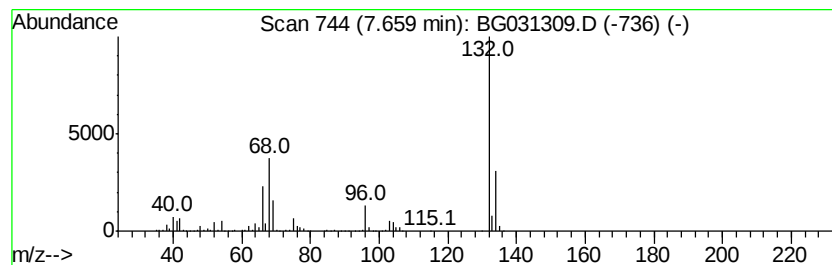
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Peak Number 2 unknown7.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	57.63 ng	1054800	1,4-Dichlorobenzene-d4	8.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3	N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	12
4	5-Aminoindole	132	C8H8N2	005192-03-0	12
5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	5.5	ng	101599	1	8.13	366086	20.0
unknown7.66	7.66	57.6	ng	1054800	1	8.13	366086	20.0