

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121117\
 Data File : BG031462.D
 Acq On : 11 Dec 2017 19:14
 Operator : SJ/JU
 Sample : PB104880BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB104880BL

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-BG121117.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	5.180	317	322	336	rVB	113902	185024	3.72%	1.096%
2	5.667	399	405	421	rBV	437472	787270	15.84%	4.666%
3	7.271	671	678	693	rBV	431630	833465	16.77%	4.939%
4	7.642	734	741	752	rBV	471562	838237	16.86%	4.968%
5	8.106	813	820	831	rBV	139335	255044	5.13%	1.511%
6	8.435	868	876	887	rBV	521879	971263	19.54%	5.756%
7	9.275	1011	1019	1034	rBV	344028	721898	14.52%	4.278%
8	10.926	1293	1300	1310	rBV	143714	337755	6.79%	2.002%
9	13.352	1706	1713	1725	rBV	1231820	2072434	41.69%	12.282%
10	14.733	1941	1948	1963	rBV2	293033	559252	11.25%	3.314%
11	16.220	2194	2201	2216	rBV	571283	1067193	21.47%	6.324%
12	17.477	2408	2415	2432	rBV2	479455	882572	17.75%	5.230%
13	20.062	2849	2855	2873	rBV	3580751	4971410	100.00%	29.462%
14	21.743	3134	3141	3158	rVB2	687852	1248369	25.11%	7.398%
15	25.021	3687	3699	3716	rBV2	367438	1143046	22.99%	6.774%

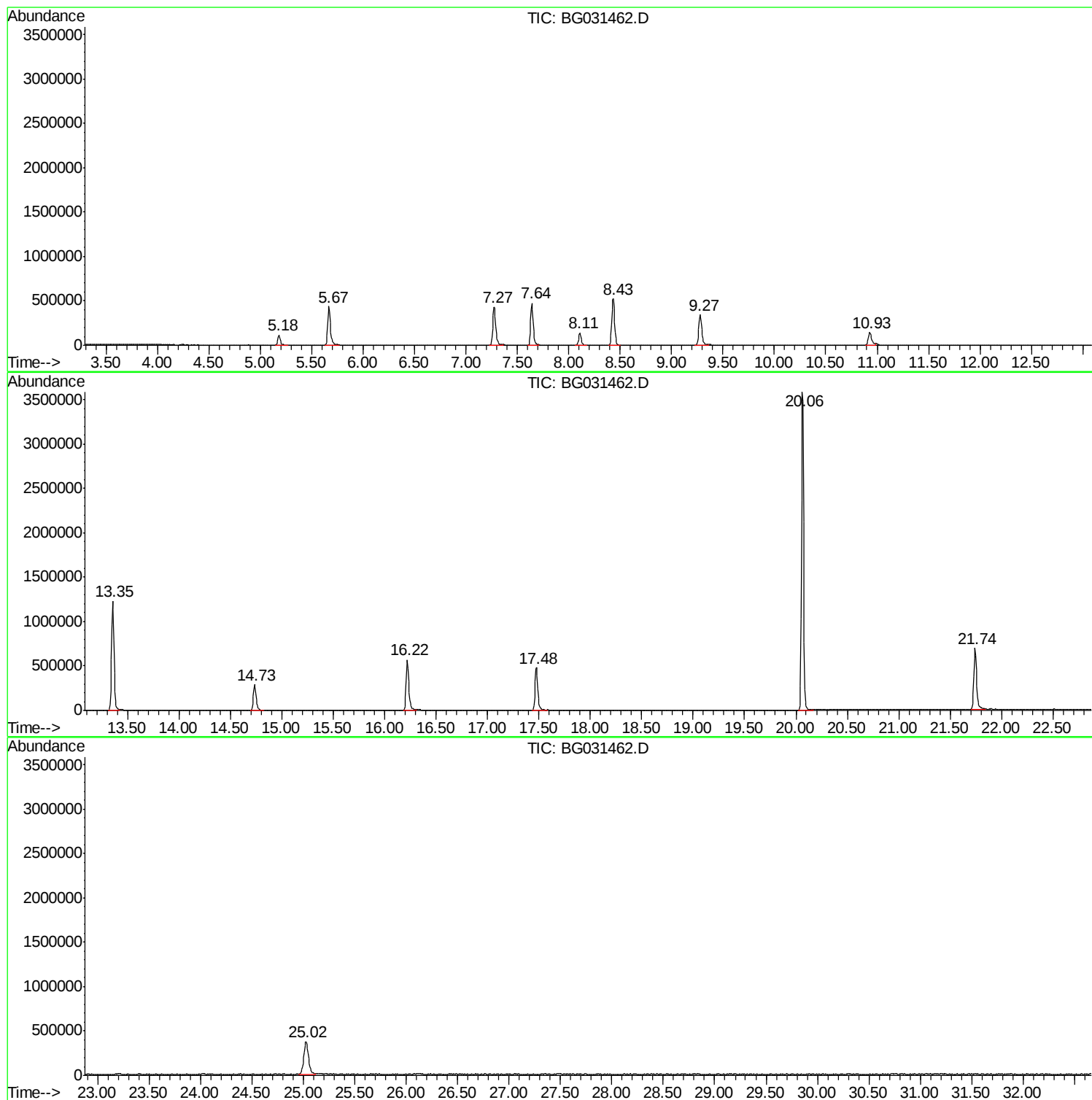
Sum of corrected areas: 16874232

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



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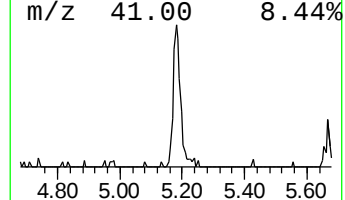
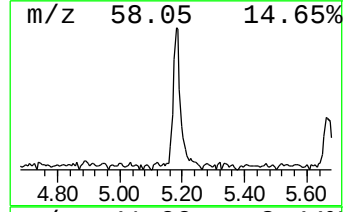
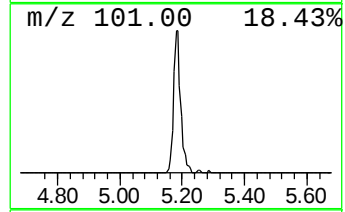
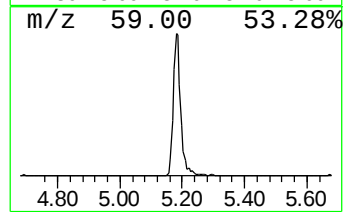
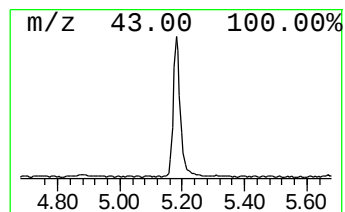
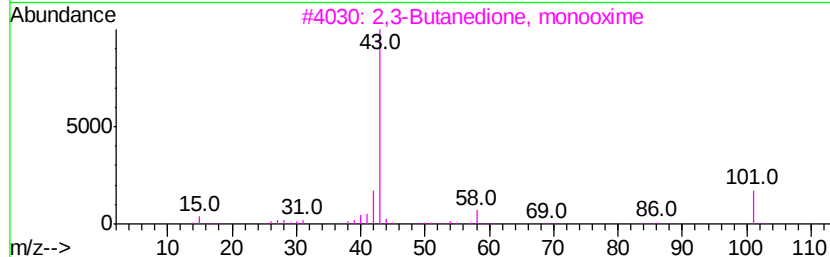
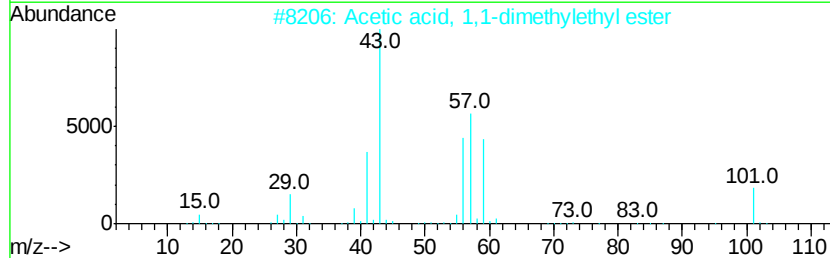
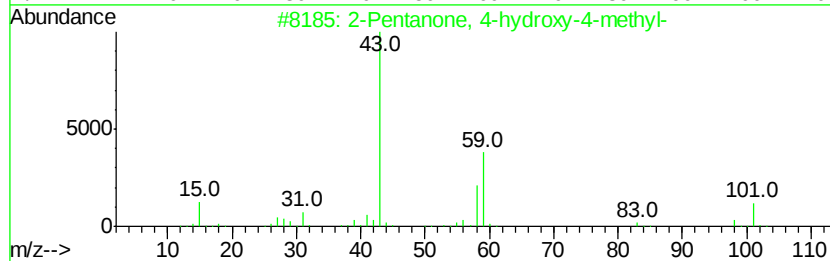
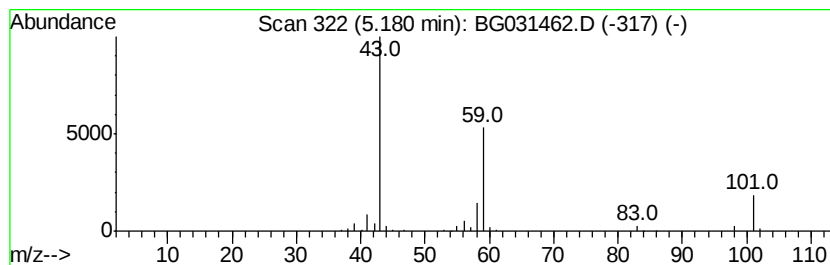
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	14.51 ng	185024	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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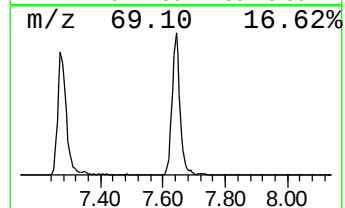
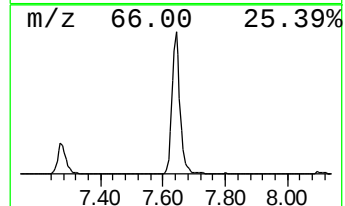
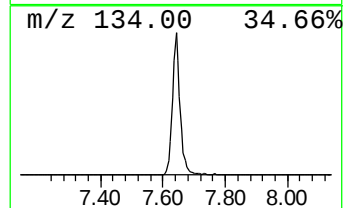
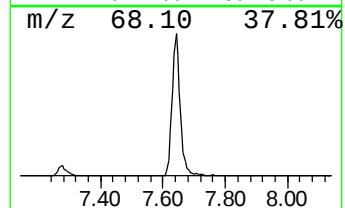
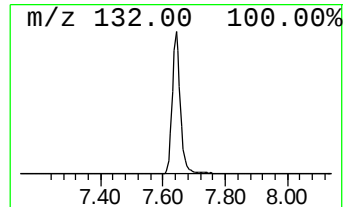
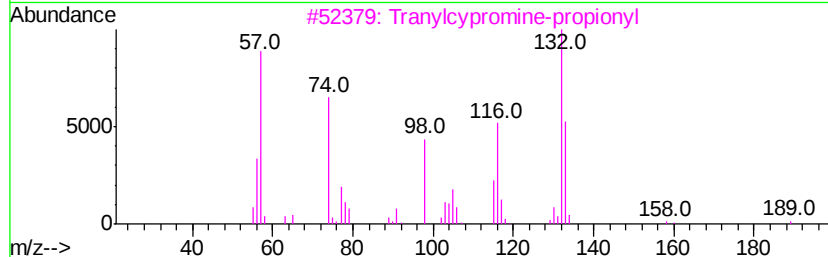
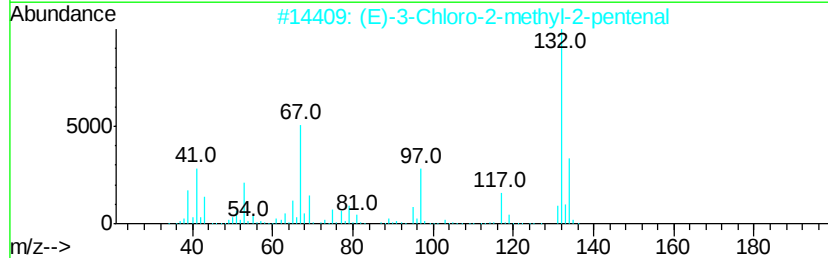
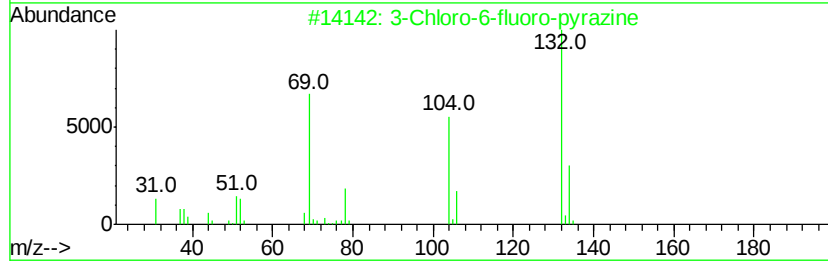
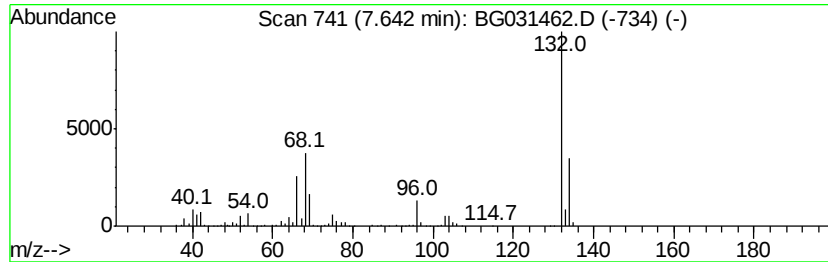
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	65.73 ng	838237	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcypropane-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
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...	5.18	14.5	ng	185024	1	8.11	255044	20.0
unknown7.64	7.64	65.7	ng	838237	1	8.11	255044	20.0