

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111017\
 Data File : BG030919.D
 Acq On : 10 Nov 2017 22:59
 Operator : SJ/JU
 Sample : PB104078BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB104078BL

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	5.218	323	328	340	rVB	98346	146939	4.10%	1.121%
2	5.705	403	411	424	rBV	379151	643535	17.95%	4.908%
3	7.303	676	683	699	rBV	363292	675168	18.83%	5.149%
4	7.679	739	747	759	rBV	380292	680300	18.98%	5.188%
5	8.144	817	826	834	rBV	115047	200669	5.60%	1.530%
6	8.467	873	881	891	rBV	458232	792797	22.12%	6.046%
7	9.307	1017	1024	1037	rBV	306632	586999	16.38%	4.476%
8	10.958	1298	1305	1326	rBV	132651	274783	7.67%	2.096%
9	13.384	1711	1718	1728	rBV	1000755	1642760	45.83%	12.528%
10	14.765	1946	1953	1968	rVB2	248268	424068	11.83%	3.234%
11	16.246	2198	2205	2221	rBV	598069	982510	27.41%	7.493%
12	17.503	2412	2419	2431	rBV2	393993	634209	17.69%	4.837%
13	20.094	2854	2860	2871	rBV	2664235	3584710	100.00%	27.337%
14	21.775	3139	3146	3158	rBV	532598	913854	25.49%	6.969%
15	25.077	3696	3708	3721	rBV2	301776	929610	25.93%	7.089%

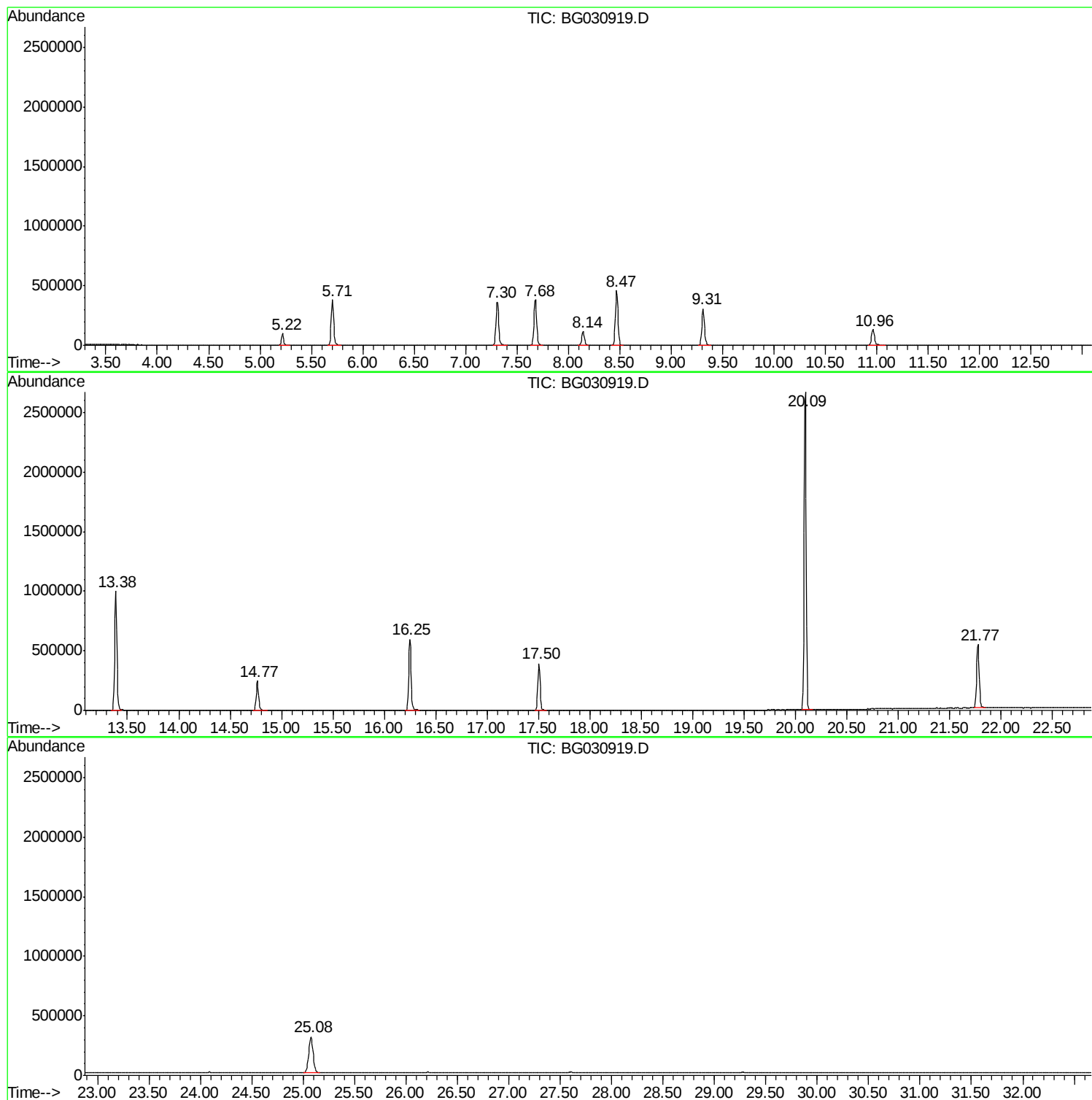
Sum of corrected areas: 13112911

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



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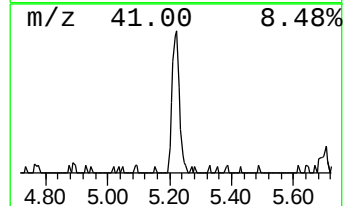
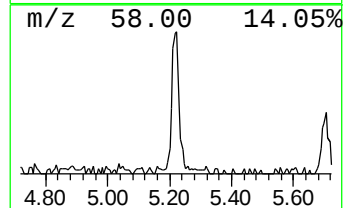
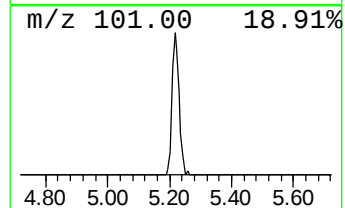
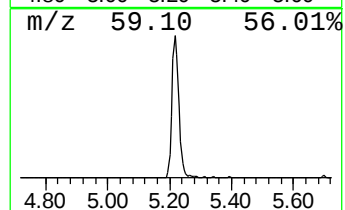
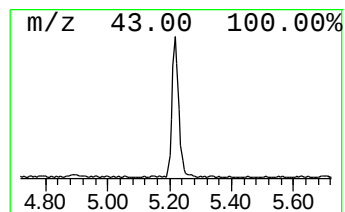
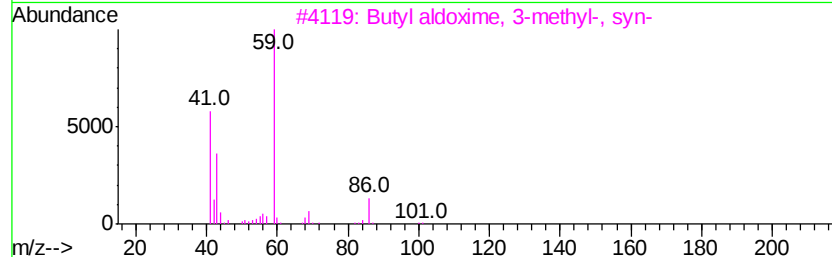
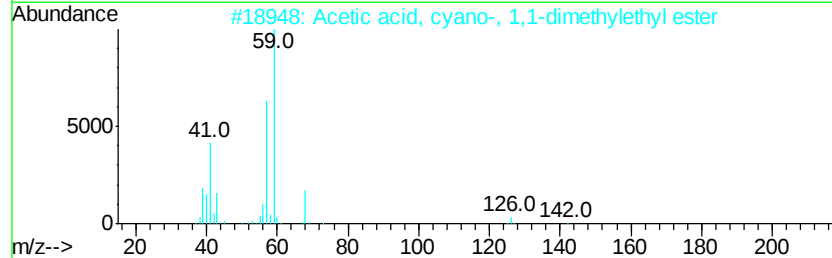
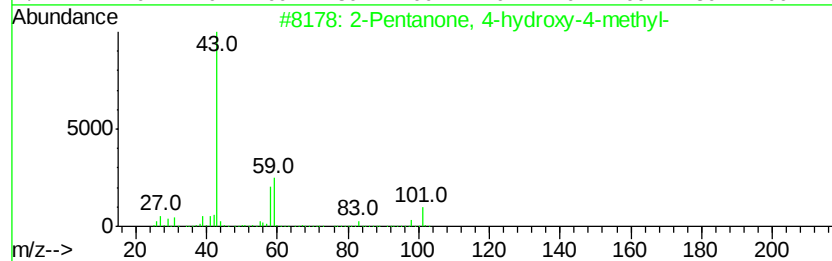
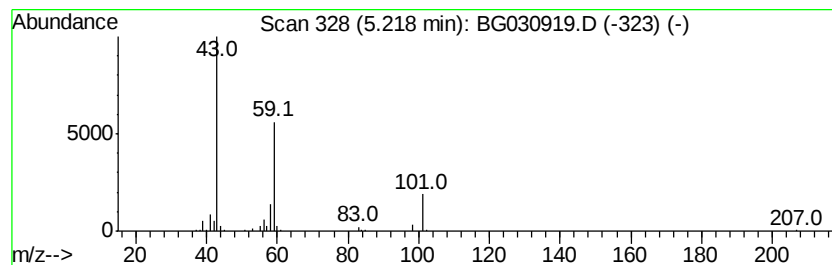
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	14.64 ng	146939	1,4-Dichlorobenzene-d4	8.14

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2		000123-42-2	59
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2		001116-98-9	25
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO		005780-40-5	23
4	2,3-Butanedione, monooxime	101	C4H7NO2		000057-71-6	16
5	1-Propen-2-ol, acetate	100	C5H8O2		000108-22-5	10



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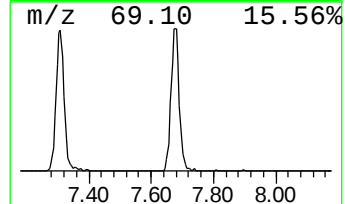
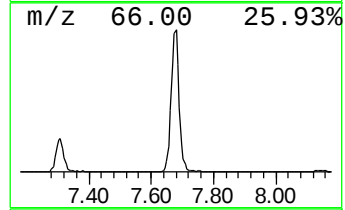
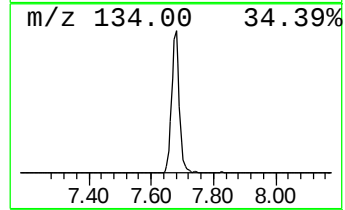
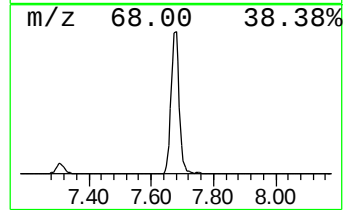
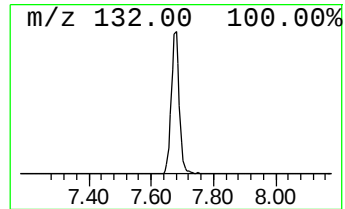
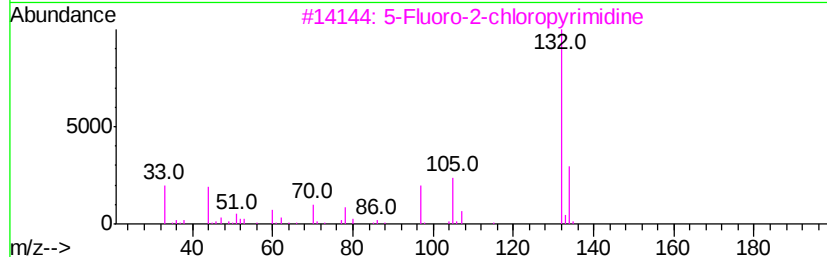
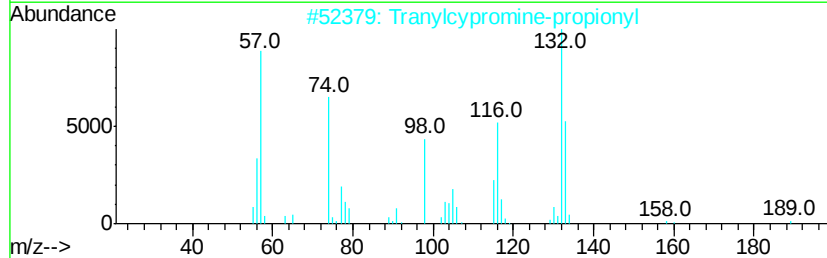
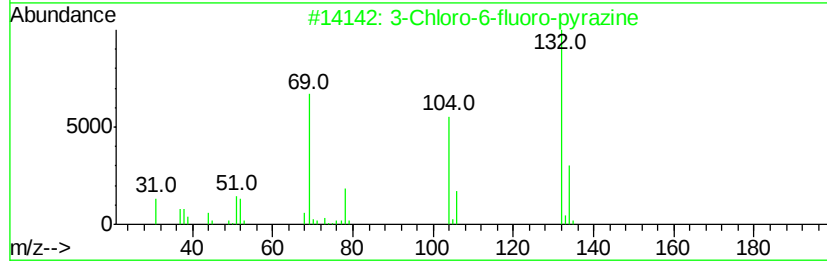
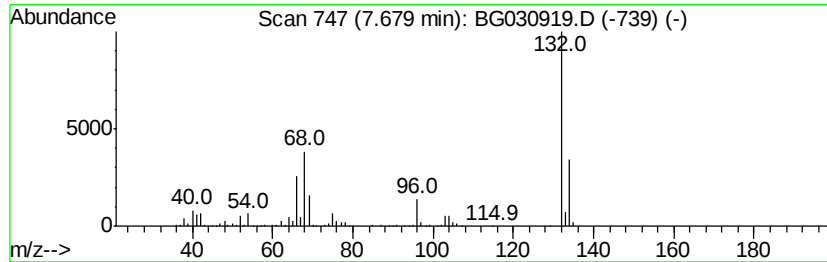
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Peak Number 2 unknown7.68 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	67.80 ng	680300	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	14.6	ng	146939	1	8.14	200669	20.0
unknown7.68	7.68	67.8	ng	680300	1	8.14	200669	20.0