

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111317\
 Data File : BG030973.D
 Acq On : 13 Nov 2017 15:09
 Operator : SJ/JU
 Sample : PB104186BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB104186BL

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	5.221	322	329	338	rBV	109157	172436	4.13%	1.018%
2	5.703	404	411	428	rBV	682094	1100960	26.35%	6.501%
3	7.307	677	684	694	rBV	642631	1126089	26.95%	6.650%
4	7.677	737	747	757	rBV	657017	1160016	27.76%	6.850%
5	8.141	819	826	835	rVB	142983	240036	5.74%	1.417%
6	8.465	873	881	890	rBV	523249	944759	22.61%	5.579%
7	9.311	1016	1025	1036	rBV	356392	684330	16.38%	4.041%
8	10.962	1298	1306	1314	rBV	161135	320702	7.67%	1.894%
9	13.382	1711	1718	1728	rBV	1219866	1912857	45.78%	11.295%
10	14.757	1945	1952	1962	rBV2	286634	489938	11.72%	2.893%
11	16.244	2198	2205	2223	rBV	1140262	1834515	43.90%	10.833%
12	17.495	2408	2418	2430	rBV	458033	734481	17.58%	4.337%
13	20.086	2853	2859	2868	rBV	3104262	4178805	100.00%	24.676%
14	21.767	3138	3145	3159	rVB	601658	1041834	24.93%	6.152%
15	25.063	3695	3706	3721	rBV	338197	993112	23.77%	5.864%

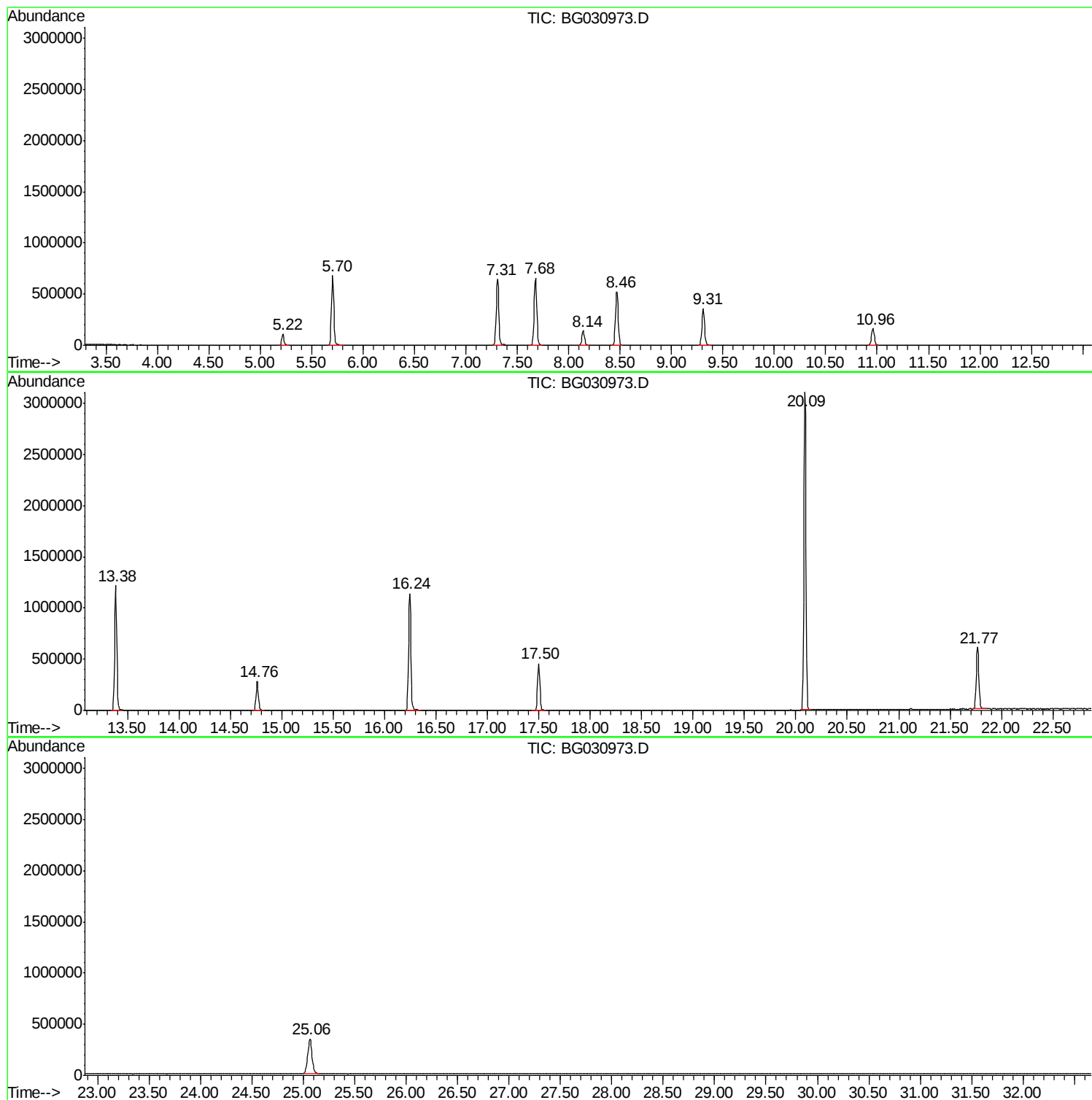
Sum of corrected areas: 16934870

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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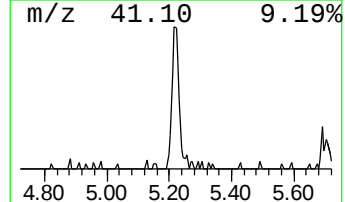
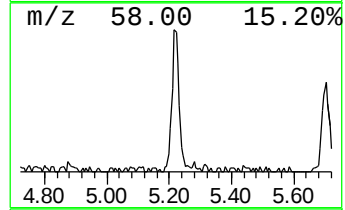
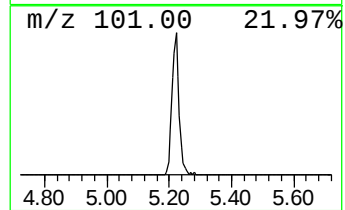
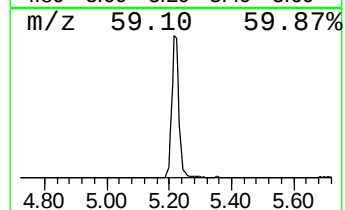
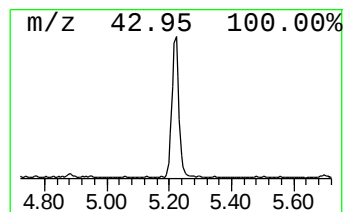
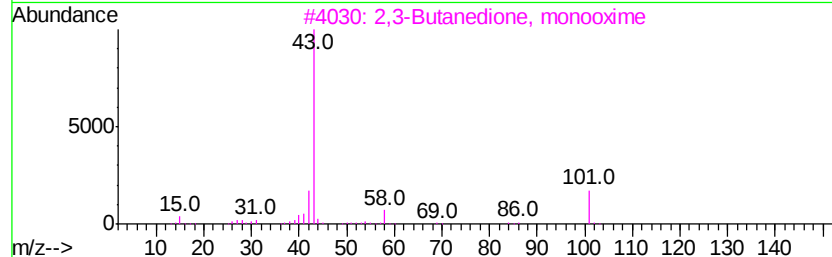
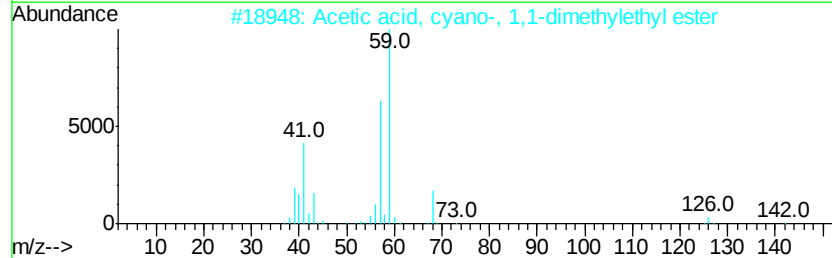
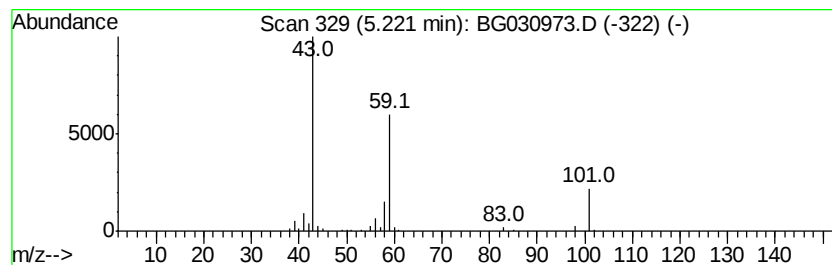
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	14.37 ng	172436	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5		Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9



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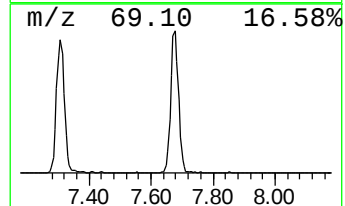
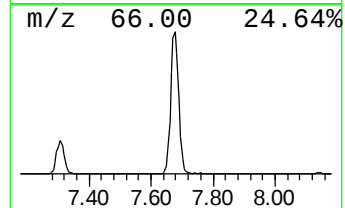
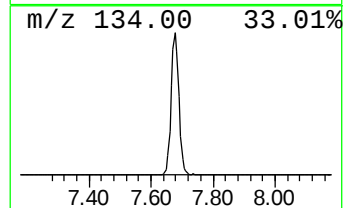
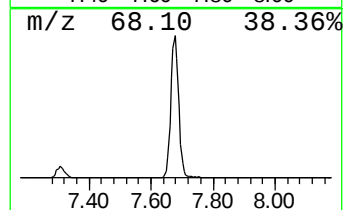
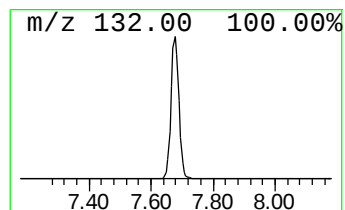
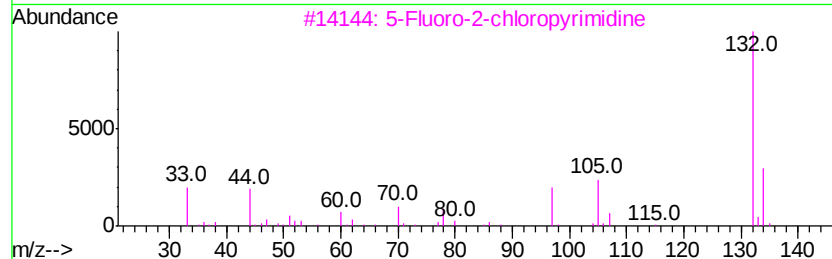
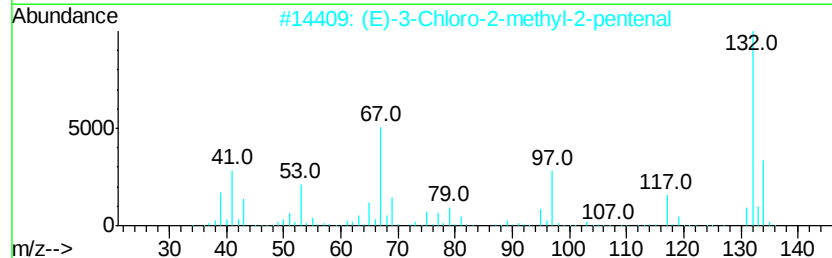
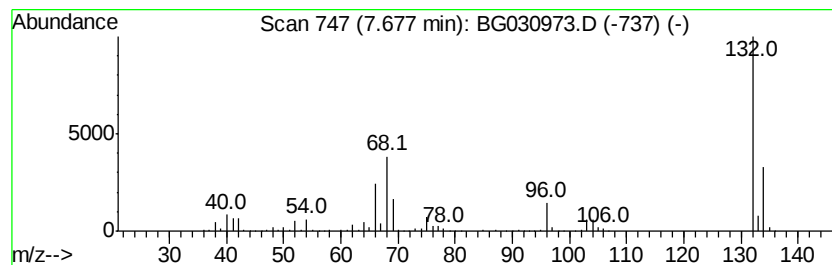
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Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	96.65 ng	1160020	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		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	14.4	ng	172436	1	8.14	240036	20.0
unknown7.68	7.68	96.7	ng	1160020	1	8.14	240036	20.0