

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110117\
 Data File : BG030646.D
 Acq On : 1 Nov 2017 22:39
 Operator : SJ/JU
 Sample : I6203-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 102717-TIDO-E-U-S

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-BG101617.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	5.240	326	332	341	rVB	85772	134068	2.12%	0.537%
2	5.716	406	413	425	rBV	470667	762559	12.07%	3.057%
3	7.314	678	685	697	rBV	353339	640458	10.13%	2.567%
4	7.690	741	749	762	rBV	783133	1371487	21.70%	5.498%
5	8.160	822	829	836	rBV	226862	384793	6.09%	1.542%
6	8.484	876	884	894	rVB	1021848	1790449	28.33%	7.177%
7	9.324	1018	1027	1036	rBV	833172	1497427	23.70%	6.002%
8	10.975	1300	1308	1320	rBV	333313	615528	9.74%	2.467%
9	13.401	1713	1721	1733	rBV	2659701	4195100	66.38%	16.816%
10	14.776	1946	1955	1964	rBV2	608312	1030194	16.30%	4.129%
11	16.257	2200	2207	2222	rBV	1357512	2108266	33.36%	8.451%
12	17.514	2413	2421	2432	rBV	778597	1212361	19.18%	4.860%
13	20.105	2855	2862	2874	rBV	4656816	6319563	100.00%	25.332%
14	21.786	3140	3148	3158	rBV	879983	1447234	22.90%	5.801%
15	25.094	3700	3711	3731	rBV3	465785	1437954	22.75%	5.764%

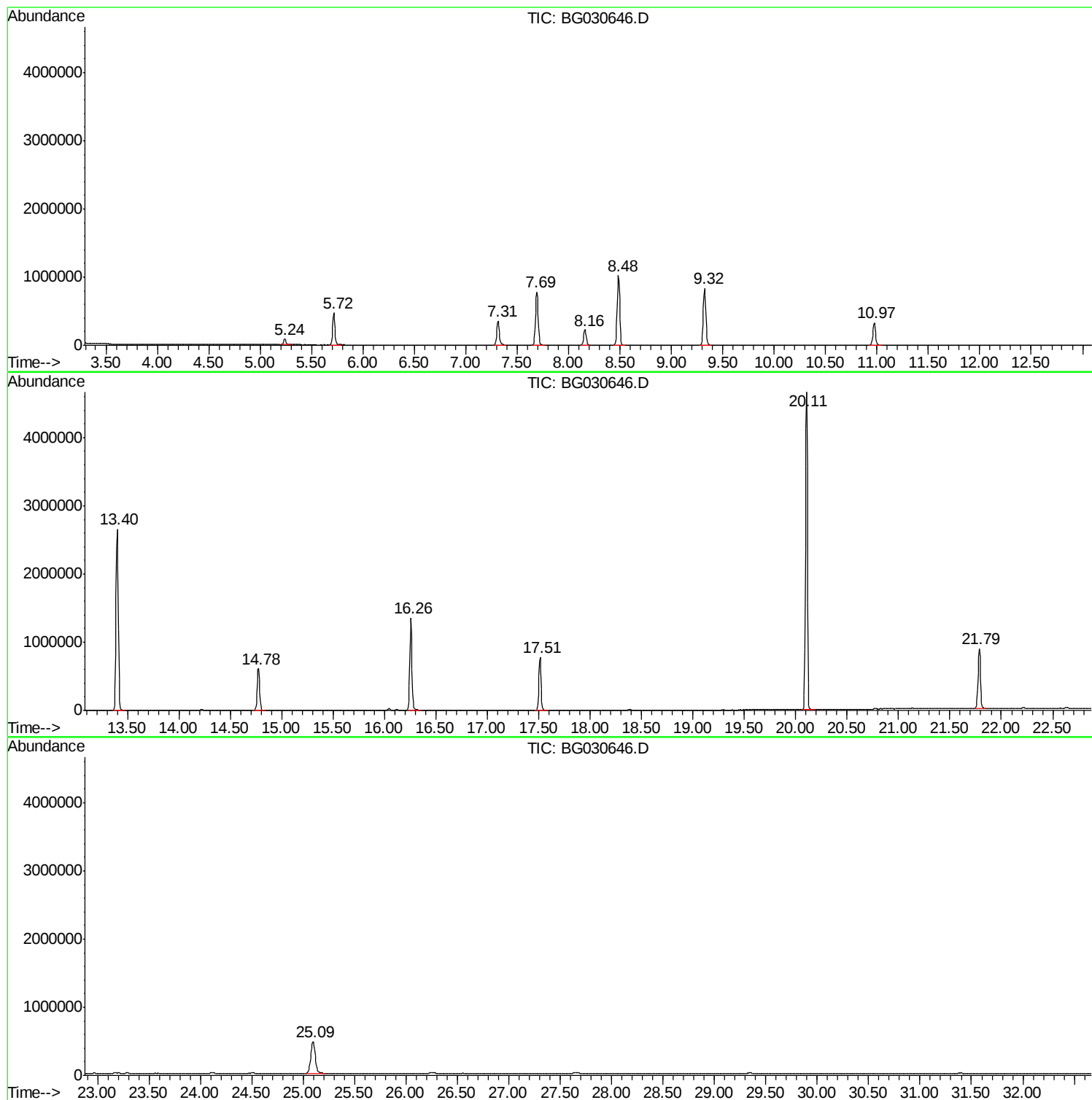
Sum of corrected areas: 24947441

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.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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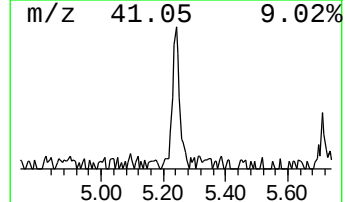
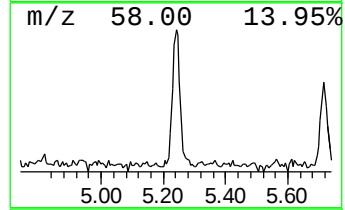
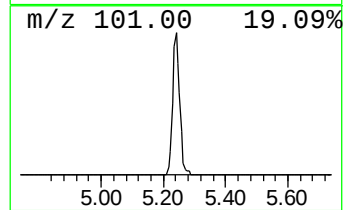
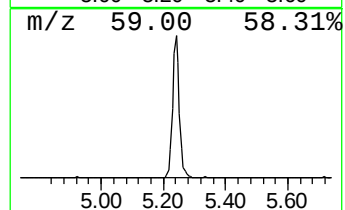
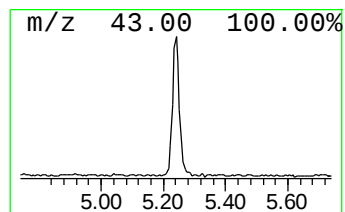
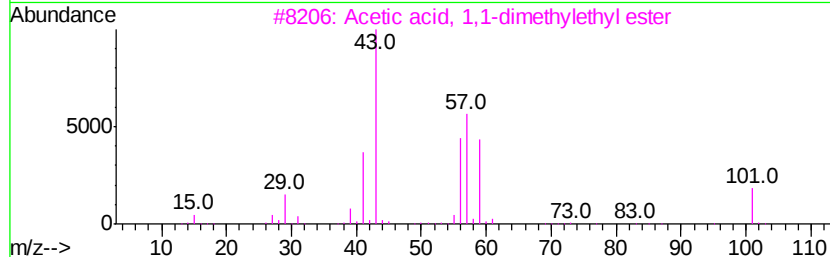
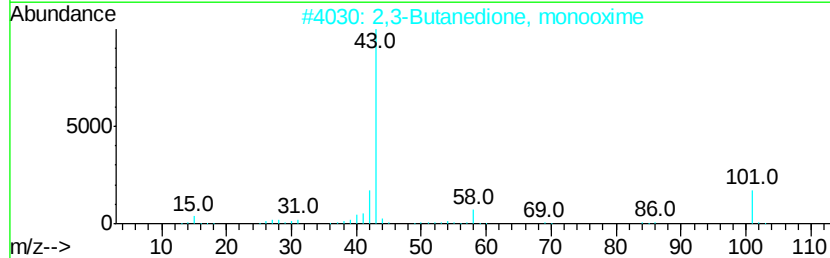
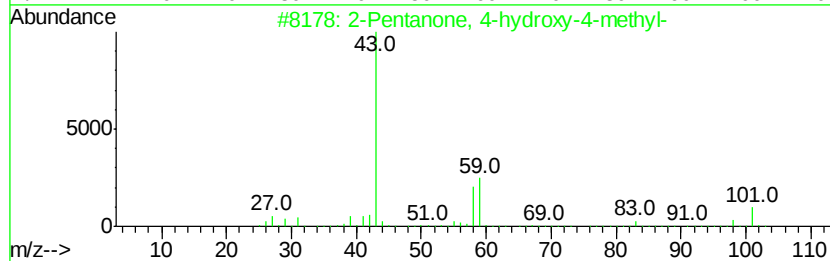
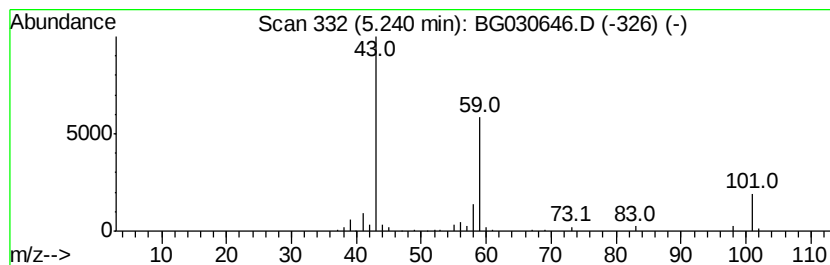
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	6.97 ng	134068	1,4-Dichlorobenzene-d4	8.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
4			3-Pentanol	88	C5H12O	000584-02-1	25
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25



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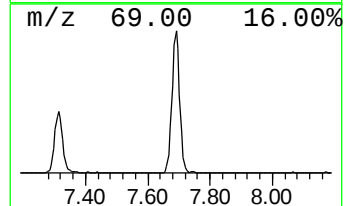
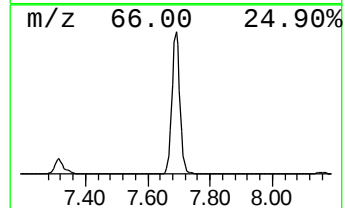
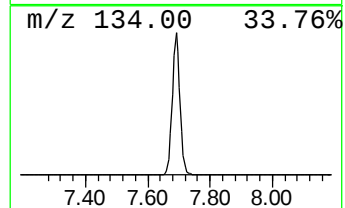
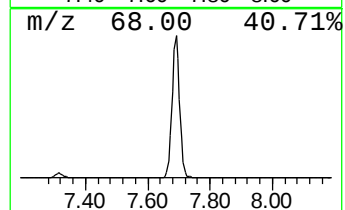
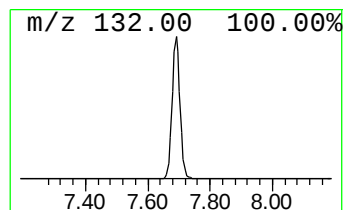
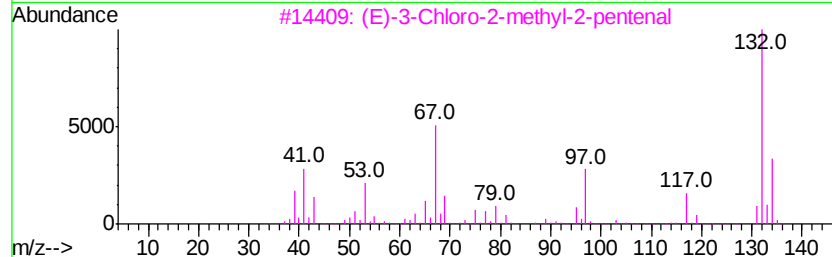
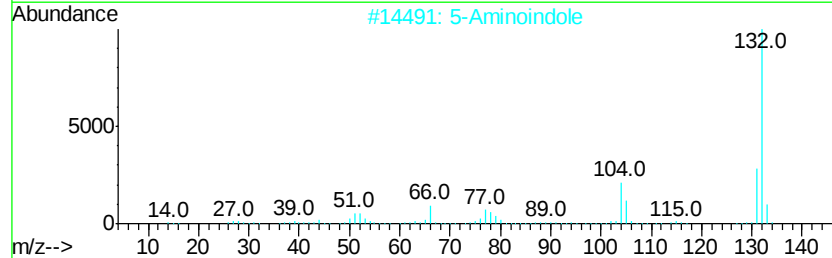
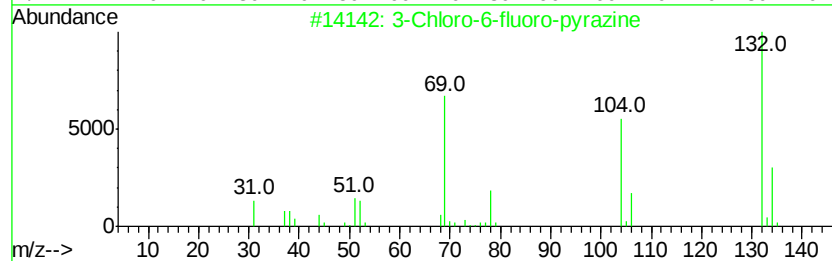
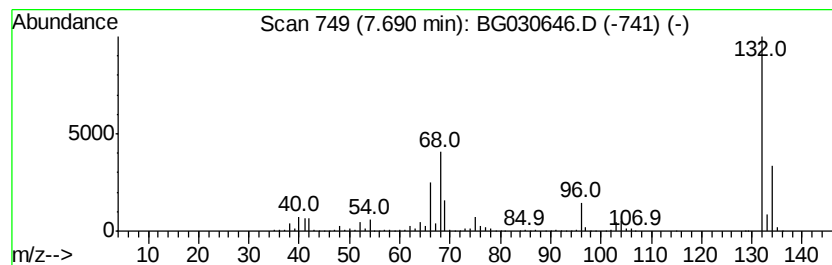
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Peak Number 2 unknown7.69 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	71.28 ng	1371490	1,4-Dichlorobenzene-d4	8.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.24	7.0	ng	134068	1	8.16	384793	20.0
unknown7.69	7.69	71.3	ng	1371490	1	8.16	384793	20.0