

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG051315\
 Data File : BG017131.D
 Acq On : 14 May 2015 5:36
 Operator : TP/IZ
 Sample : PB83339BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB83339BL

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-BG051315.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	4.815	354	362	368	rBV	137970	192140	5.50%	1.098%
2	5.303	438	445	451	rBV	936087	1326763	37.99%	7.581%
3	6.907	711	718	727	rBV	979665	1470989	42.12%	8.405%
4	7.248	769	776	784	rBV	925173	1412979	40.46%	8.074%
5	7.706	846	854	862	rBV2	214598	334193	9.57%	1.910%
6	8.023	901	908	916	rBV	618965	994547	28.48%	5.683%
7	8.869	1044	1052	1059	rBV	516628	863389	24.72%	4.933%
8	10.503	1323	1330	1338	rVB	269704	452181	12.95%	2.584%
9	12.982	1745	1752	1759	rBV	1265312	1848081	52.92%	10.560%
10	14.363	1981	1987	1995	rVB3	409339	629485	18.03%	3.597%
11	15.855	2234	2241	2251	rBV	1295911	1729138	49.51%	9.880%
12	17.107	2448	2454	2460	rBV2	608103	854447	24.47%	4.882%
13	19.745	2897	2903	2911	rBV	2858189	3492204	100.00%	19.954%
14	21.290	3161	3166	3172	rBV	764519	967966	27.72%	5.531%
15	23.558	3545	3552	3561	rVB2	496427	932921	26.71%	5.331%

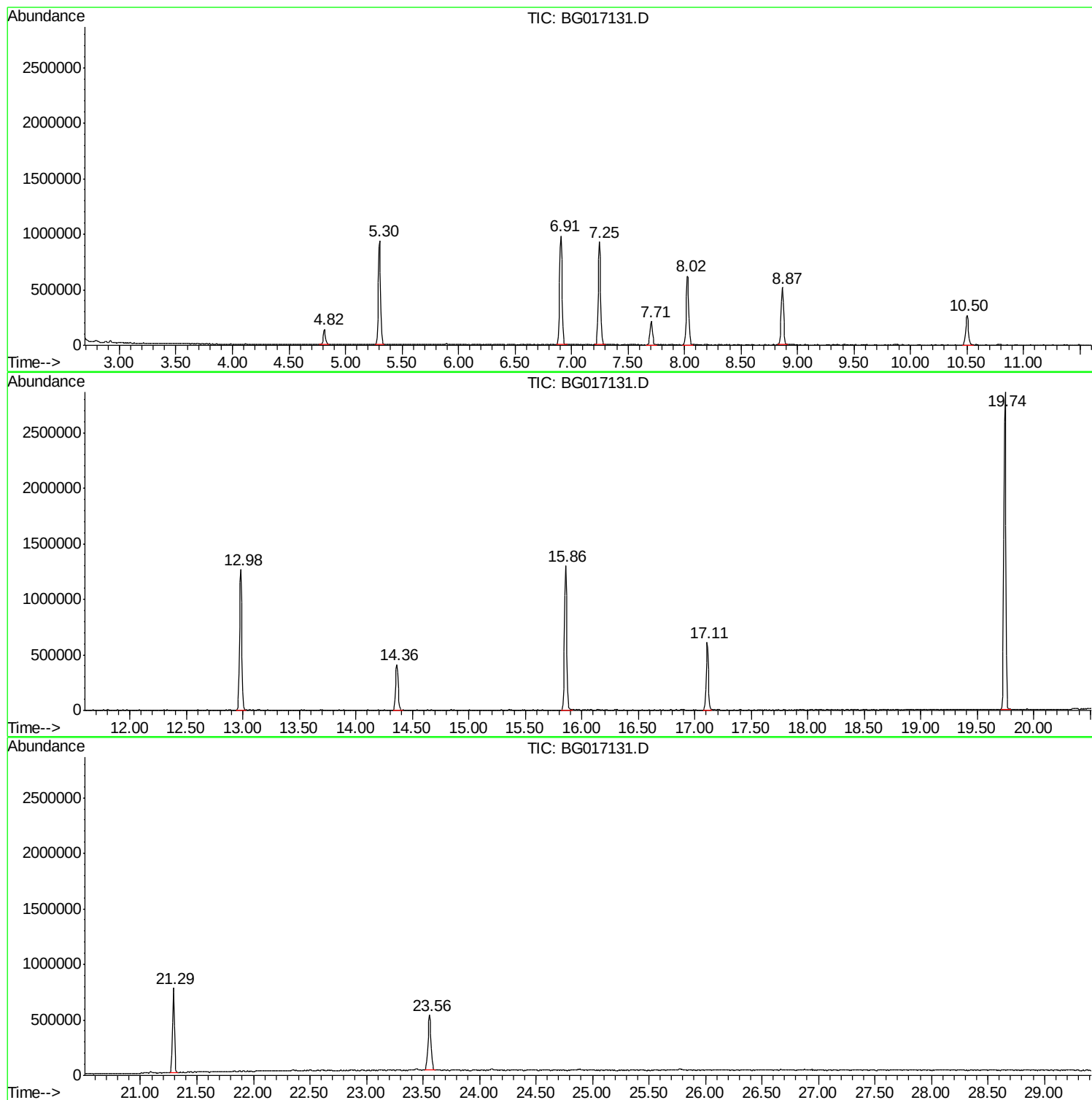
Sum of corrected areas: 17501423

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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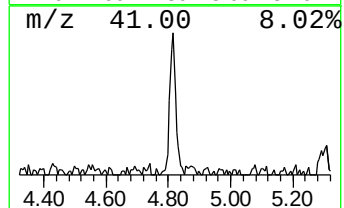
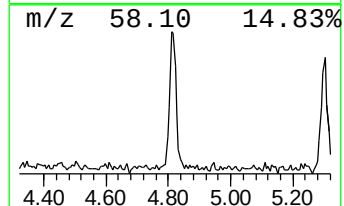
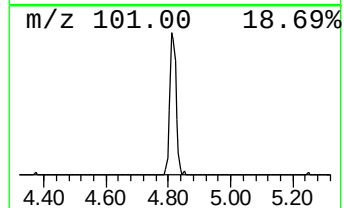
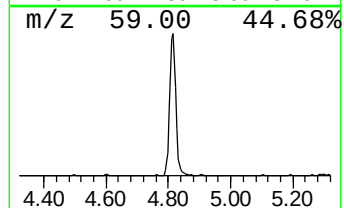
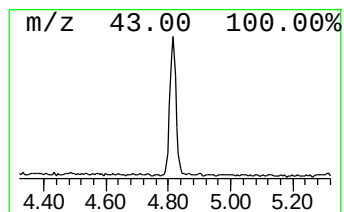
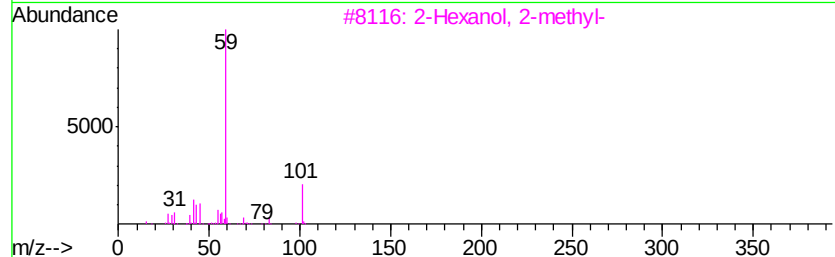
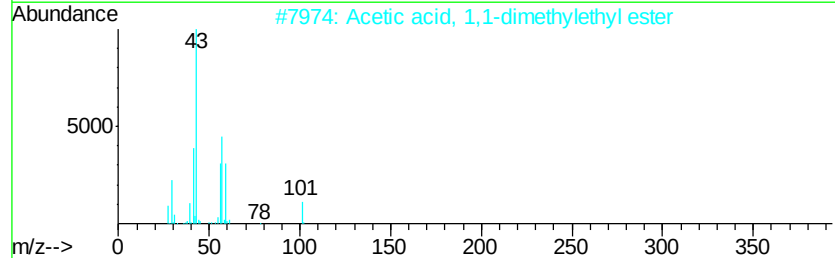
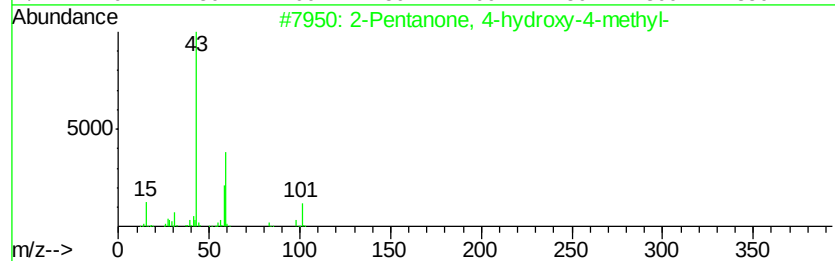
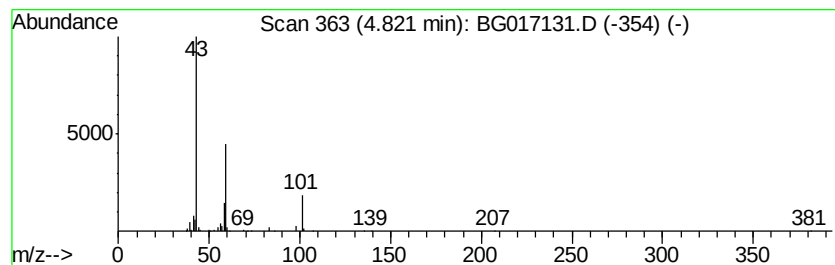
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.82	11.50 ng	192140	1,4-Dichlorobenzene-d4	7.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	25
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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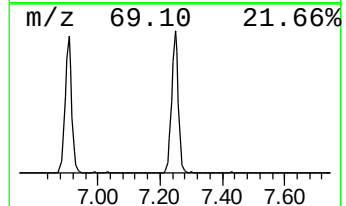
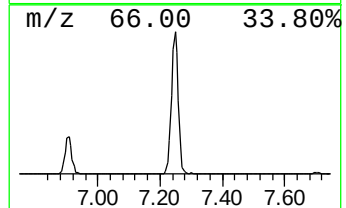
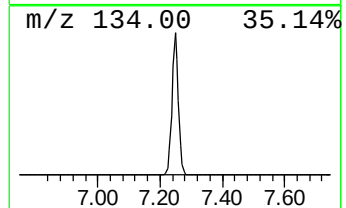
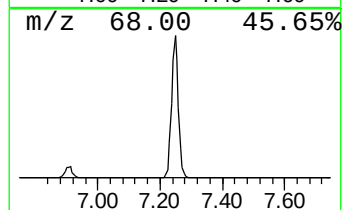
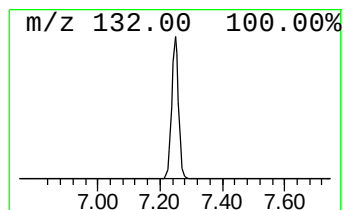
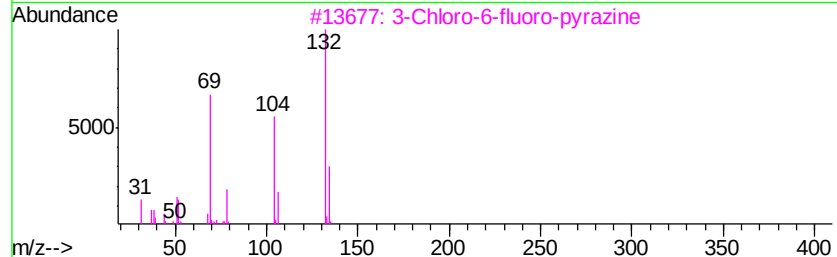
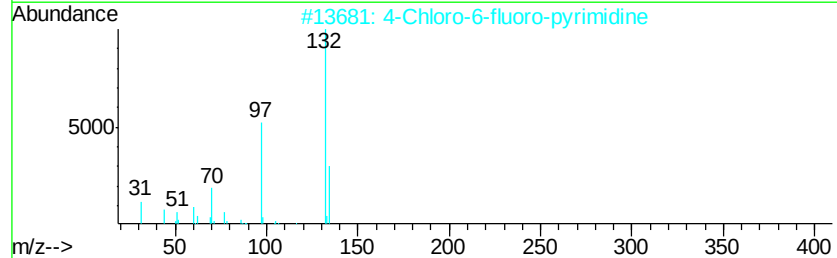
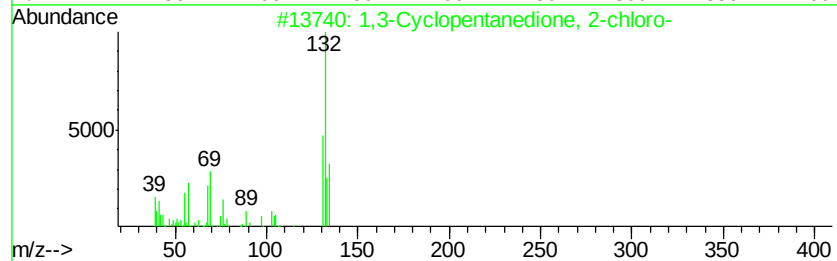
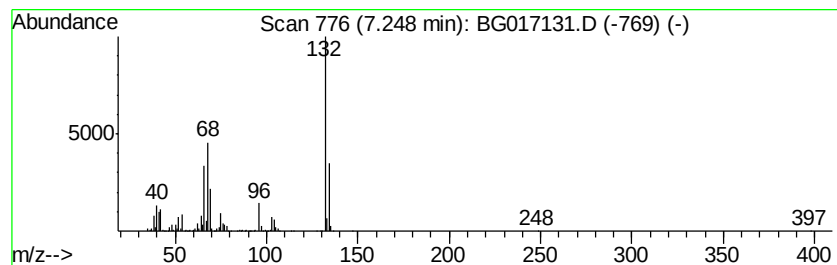
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Peak Number 2 unknown7.25 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.25	84.56 ng	1412980	1,4-Dichlorobenzene-d4	7.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	32
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.82	11.5	ng	192140	1	7.71	334193	20.0
unknown7.25	7.25	84.6	ng	1412980	1	7.71	334193	20.0