

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061515\
 Data File : BF079987.D
 Acq On : 15 Jun 2015 18:44
 Operator : TP/IZ
 Sample : G2639-11
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP15-LF2MW4-9016

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_F\METHODS\8270-BF061115.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	5.567	215	217	222	rVB	73058	79744	3.48%	0.702%
2	5.956	249	251	254	rBV	470806	482235	21.06%	4.248%
3	6.950	335	338	340	rBV	374802	308127	13.46%	2.714%
4	7.121	350	353	355	rBV	784162	1026796	44.85%	9.045%
5	7.350	371	373	375	rVB	331775	268442	11.72%	2.365%
6	7.510	385	387	389	rBV	1430951	1135036	49.57%	9.998%
7	7.910	419	422	424	rBV	786011	801286	35.00%	7.058%
8	8.642	484	486	488	rBV	454734	356179	15.56%	3.138%
9	9.716	578	580	582	rBV	2372205	1845046	80.58%	16.253%
10	10.413	639	641	644	rVB	479371	400397	17.49%	3.527%
11	11.202	708	710	713	rBV2	1017611	1041066	45.47%	9.171%
12	11.910	770	772	775	rBV	309357	416346	18.18%	3.668%
13	13.579	915	918	920	rBV	2217551	2289578	100.00%	20.169%
14	14.814	1023	1026	1028	rBV	352411	458830	20.04%	4.042%
15	16.654	1183	1187	1194	rBV	255292	443057	19.35%	3.903%

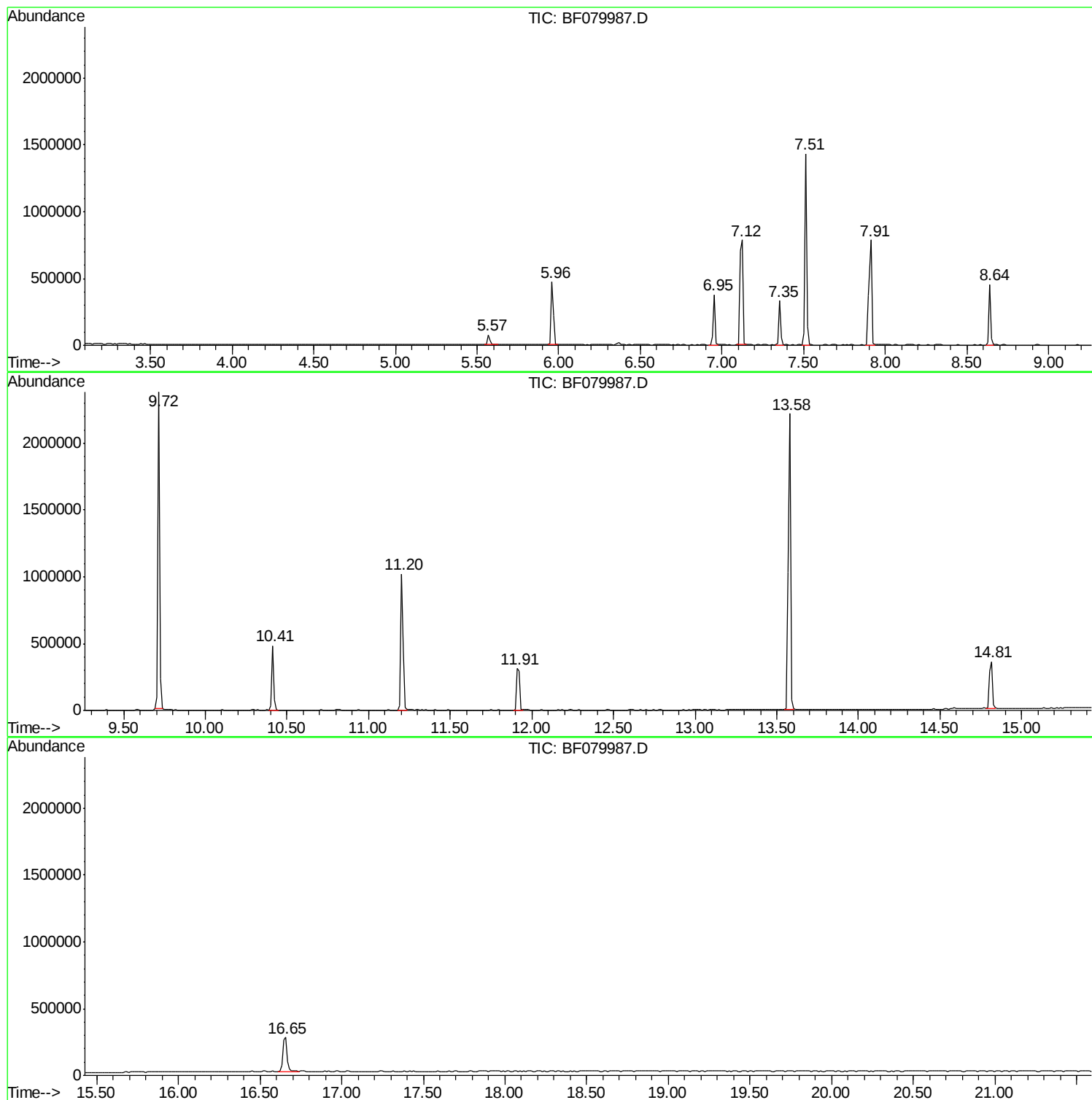
Sum of corrected areas: 11352165

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.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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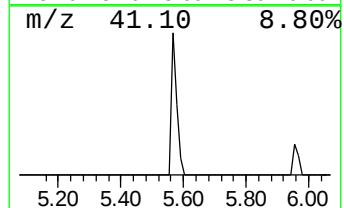
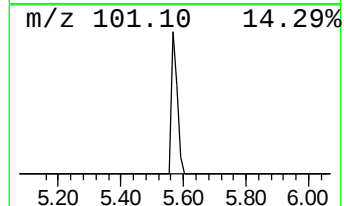
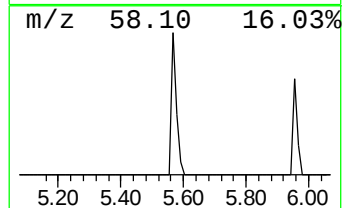
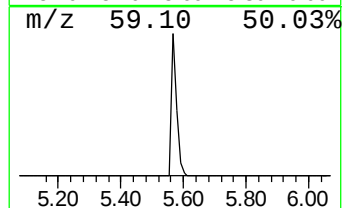
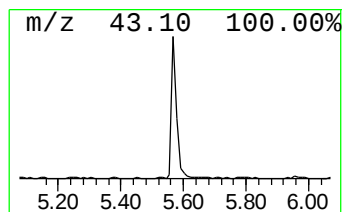
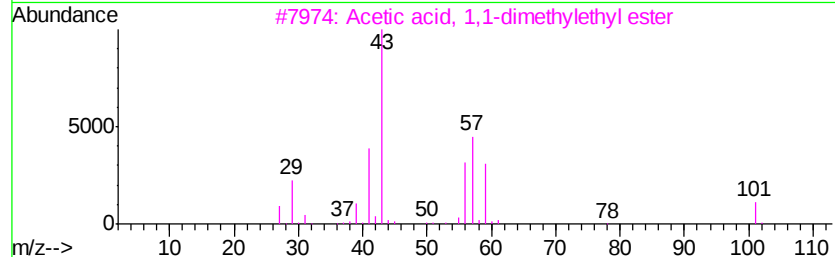
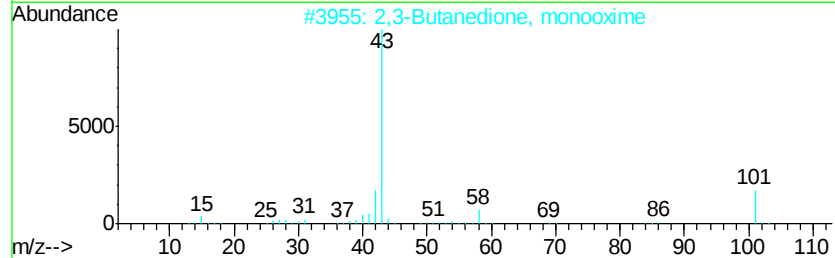
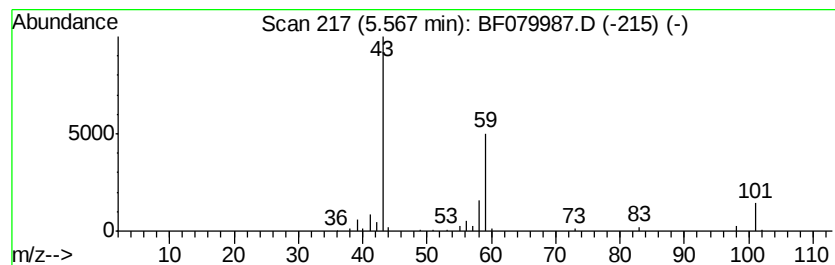
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.57	5.94 ng	79744	1,4-Dichlorobenzene-d4	7.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5			2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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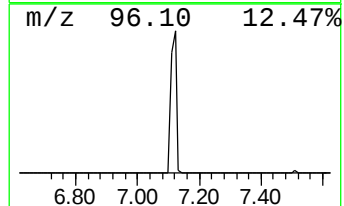
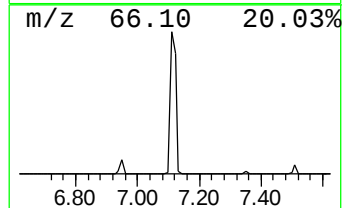
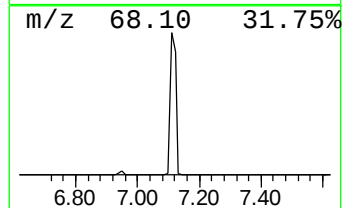
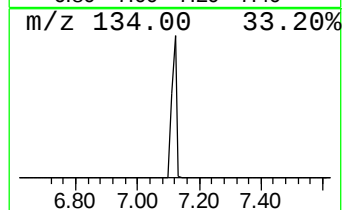
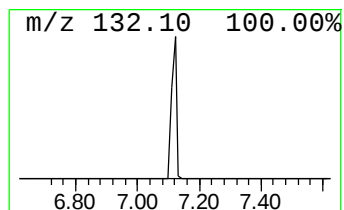
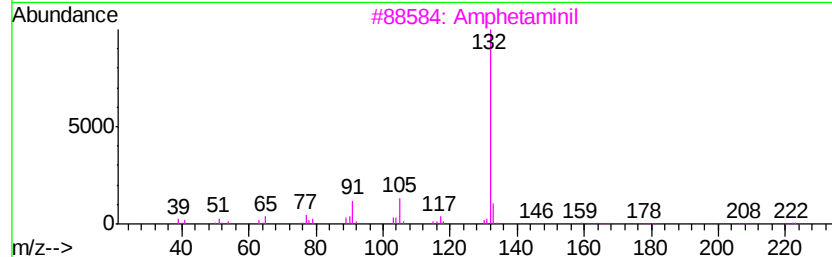
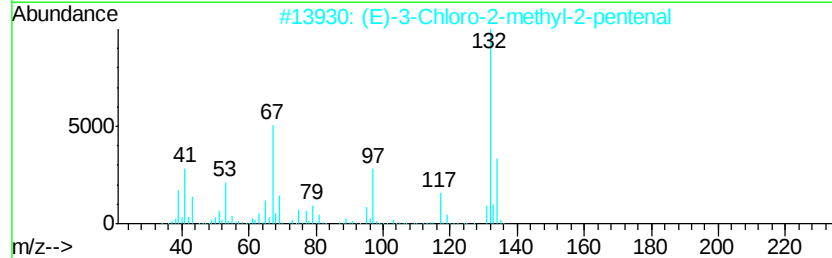
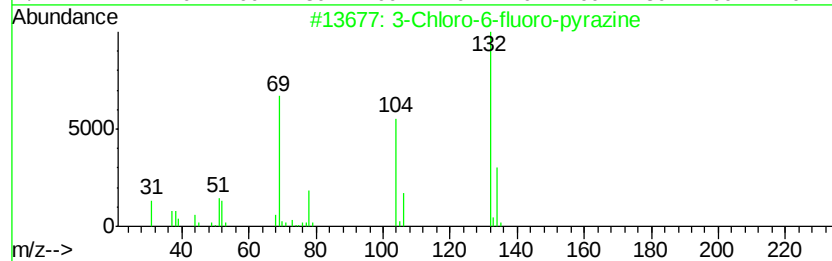
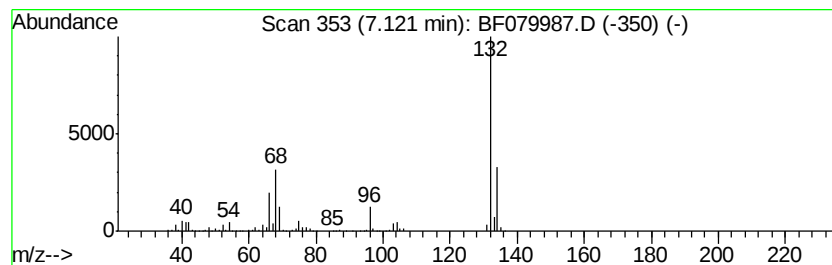
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Peak Number 2 unknown7.12 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	76.50 ng	1026800	1,4-Dichlorobenzene-d4	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.57	5.9	ng	79744	1	7.35	268442	20.0
unknown7.12	7.12	76.5	ng	1026800	1	7.35	268442	20.0