

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061615\
 Data File : BF079996.D
 Acq On : 16 Jun 2015 17:55
 Operator : TP/IZ
 Sample : PB84029BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84029BL

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	220	rBV	573373	523602	26.22%	3.743%
2	5.967	249	252	254	rBV	938513	1275775	63.89%	9.119%
3	6.367	285	287	289	rBV	34056	39131	1.96%	0.280%
4	6.950	336	338	340	rBV	1754913	1378434	69.03%	9.853%
5	7.121	350	353	355	rBV	1189739	1350628	67.64%	9.654%
6	7.350	371	373	376	rBV	400859	331828	16.62%	2.372%
7	7.510	385	387	389	rBV	1384538	1101240	55.15%	7.872%
8	7.910	419	422	424	rBV	749568	760895	38.11%	5.439%
9	8.642	484	486	488	rBV	560688	444389	22.26%	3.176%
10	9.716	578	580	583	rBV	2071562	1676376	83.95%	11.983%
11	10.413	639	641	644	rVB	602198	502069	25.14%	3.589%
12	11.202	708	710	713	rBV2	773691	947785	47.47%	6.775%
13	11.922	770	773	775	rBV	558640	530733	26.58%	3.794%
14	13.591	917	919	922	rBV	2187400	1996805	100.00%	14.273%
15	14.848	1026	1029	1032	rBV	574627	576728	28.88%	4.122%
16	16.711	1189	1192	1199	rBV	385430	553560	27.72%	3.957%

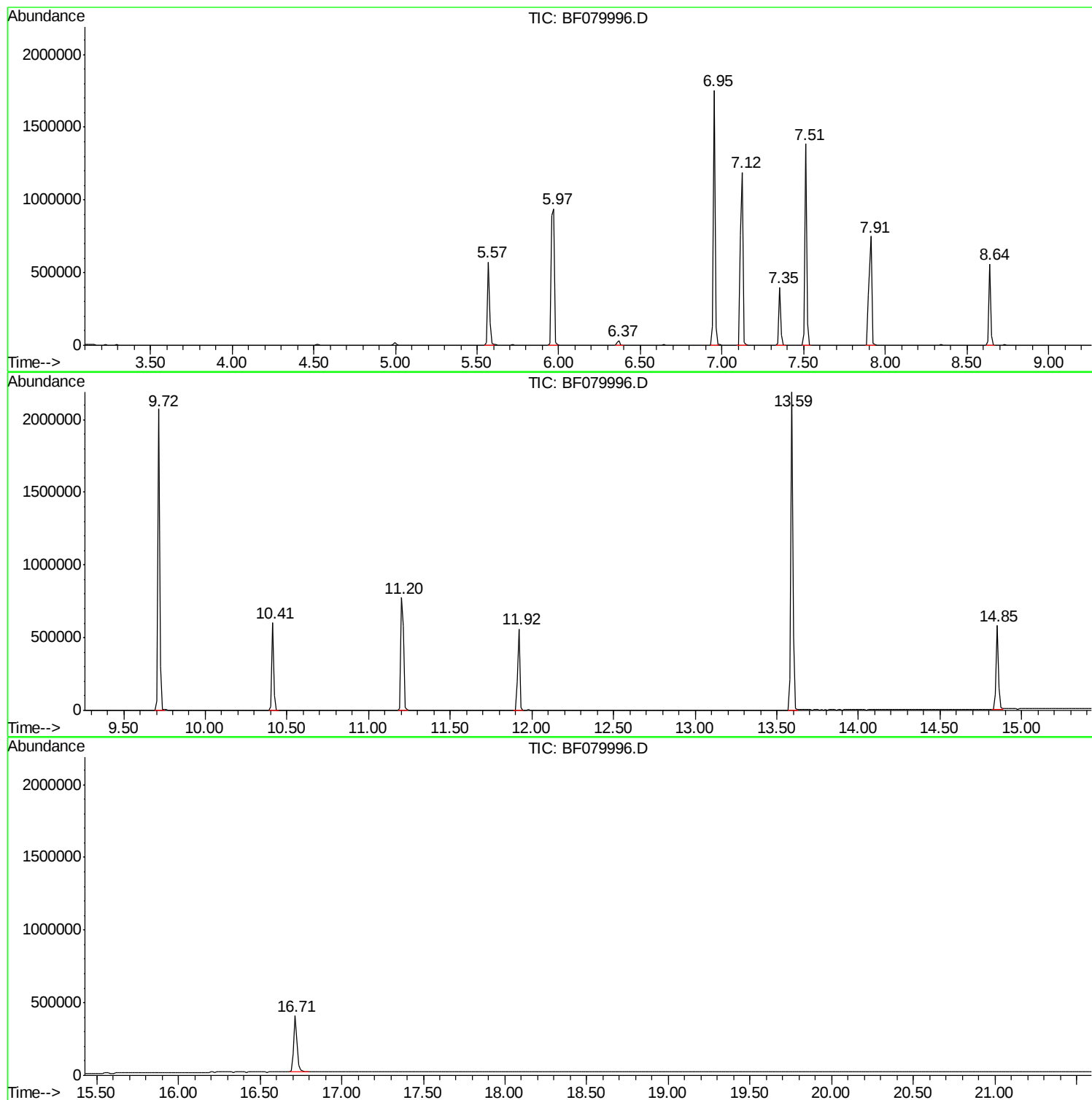
Sum of corrected areas: 13989978

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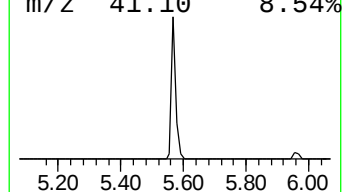
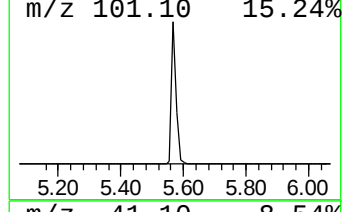
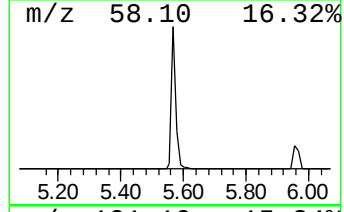
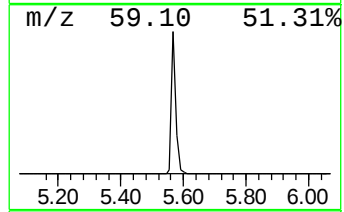
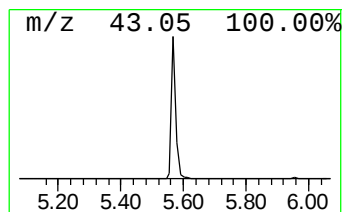
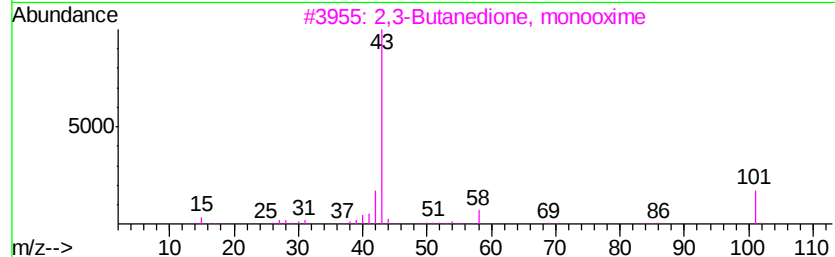
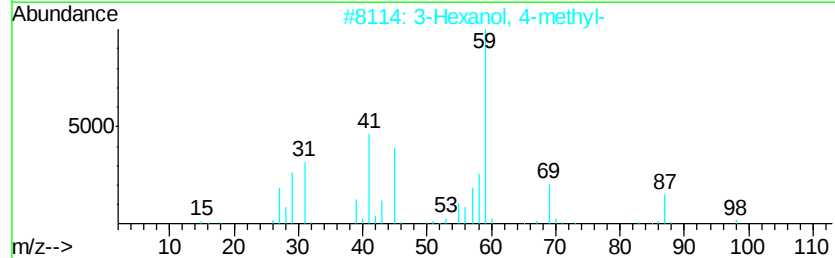
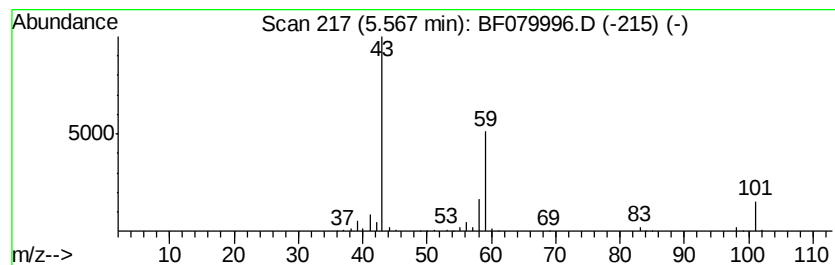
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.57	31.56 ng	523602	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		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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ClientSampled :
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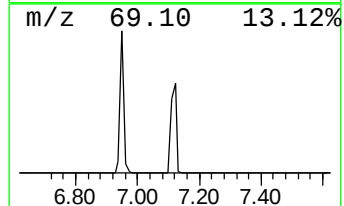
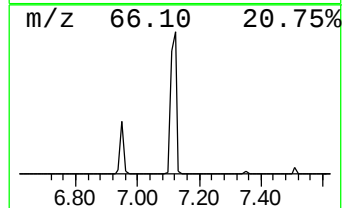
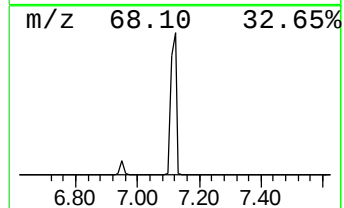
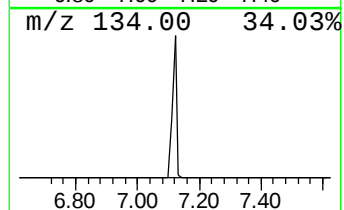
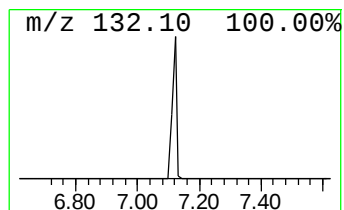
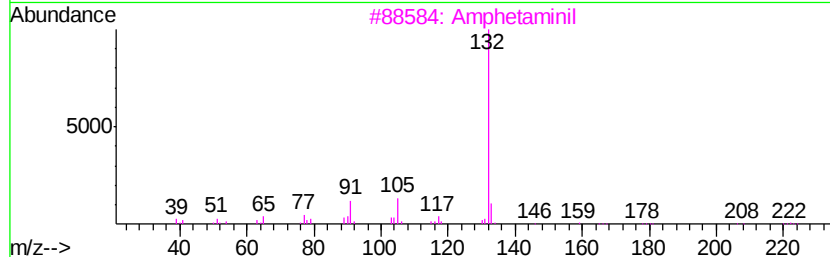
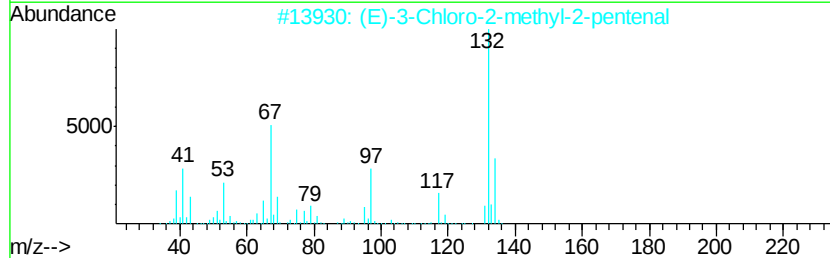
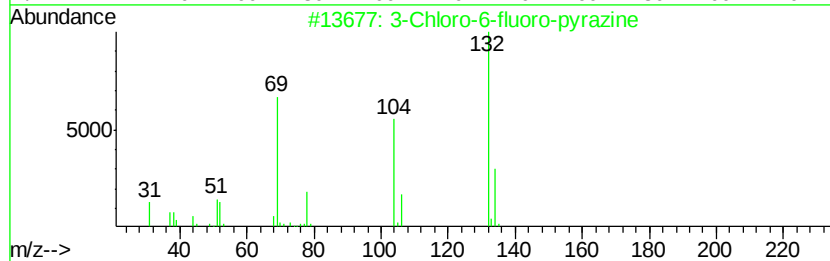
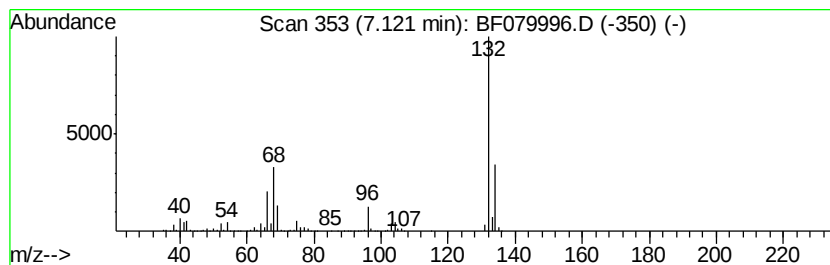
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Peak Number 3 unknown7.12 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	81.41 ng	1350630	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	17
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.57	31.6	ng	523602	1	7.35	331828	20.0
unknown7.12	7.12	81.4	ng	1350630	1	7.35	331828	20.0