

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061615\
 Data File : BF079990.D
 Acq On : 16 Jun 2015 15:02
 Operator : TP/IZ
 Sample : G2638-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP15-MMW0002

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	221	rVB	102833	99397	4.25%	0.831%
2	5.956	249	251	254	rBV	554122	561483	23.98%	4.693%
3	6.950	336	338	340	rBV	475606	373193	15.94%	3.119%
4	7.122	350	353	355	rBV	1102030	1153546	49.27%	9.641%
5	7.350	371	373	375	rVB	319395	278369	11.89%	2.327%
6	7.510	385	387	390	rBV	1370143	1156146	49.38%	9.663%
7	7.910	420	422	424	rBV	1057552	906707	38.73%	7.578%
8	8.642	484	486	489	rBV	423013	366322	15.65%	3.062%
9	9.716	578	580	583	rVB	1883041	1810043	77.31%	15.128%
10	10.413	639	641	644	rBV	421506	404409	17.27%	3.380%
11	11.213	708	711	713	rBV	1042641	1098793	46.93%	9.184%
12	11.934	772	774	776	rBV	540142	453453	19.37%	3.790%
13	13.682	924	927	929	rBV	2511758	2341220	100.00%	19.568%
14	15.008	1038	1043	1046	rBV	423724	500606	21.38%	4.184%
15	16.963	1210	1214	1219	rBV	265844	461125	19.70%	3.854%

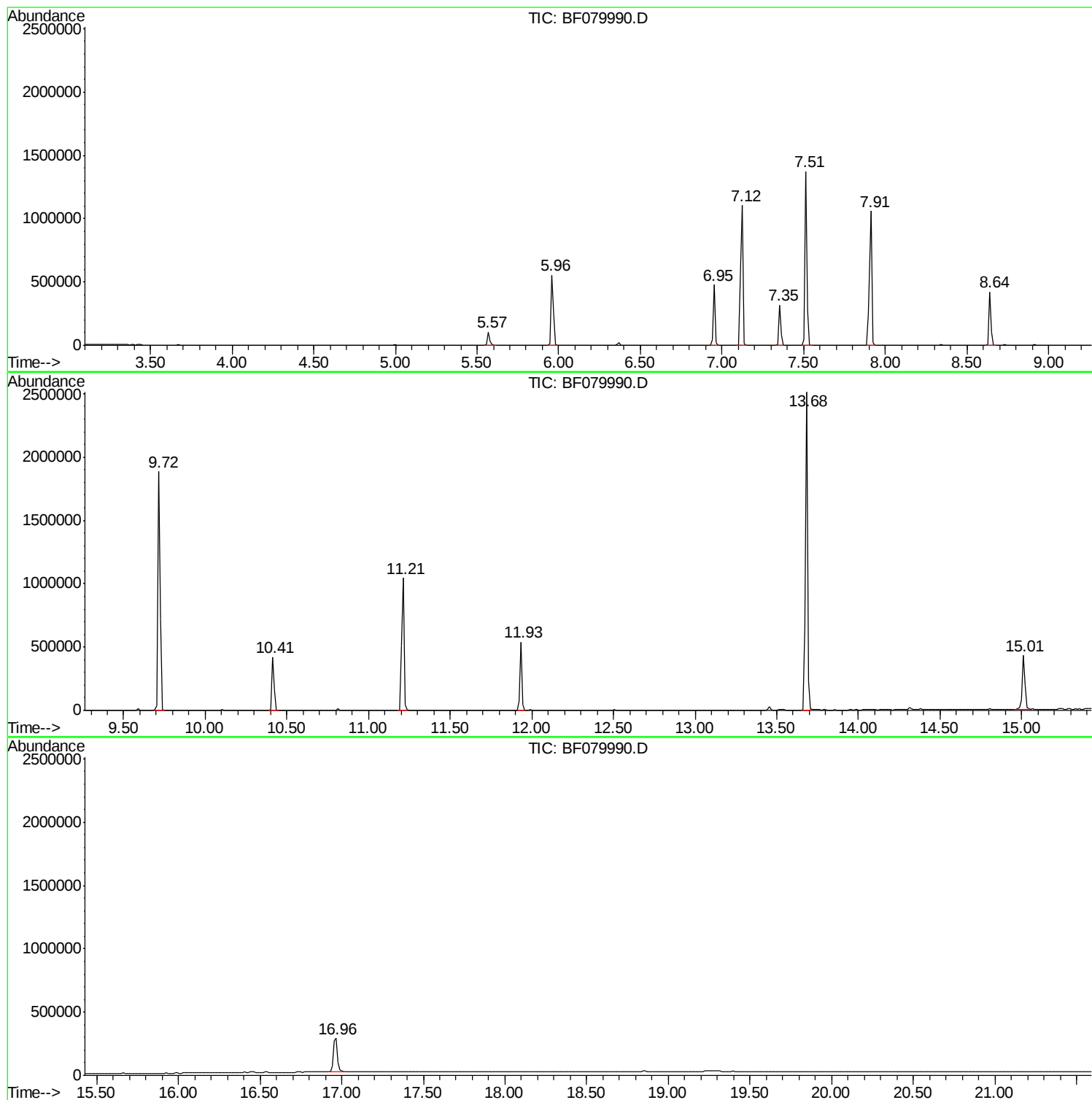
Sum of corrected areas: 11964812

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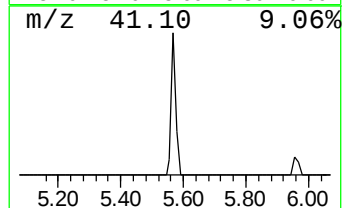
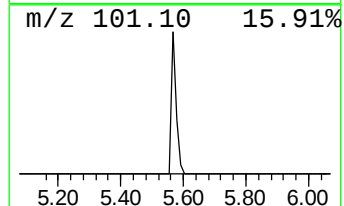
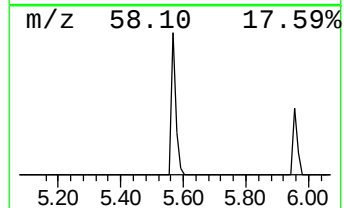
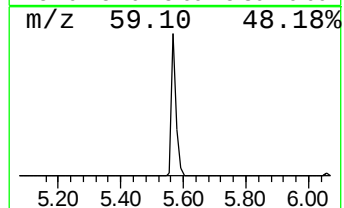
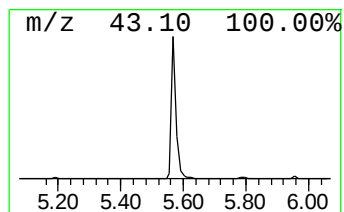
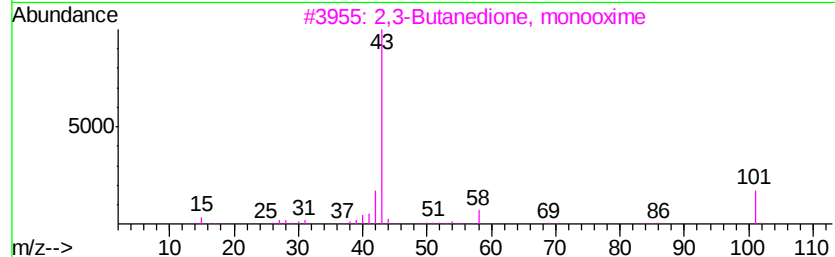
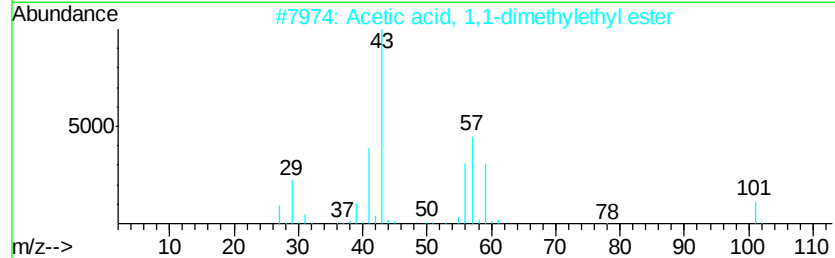
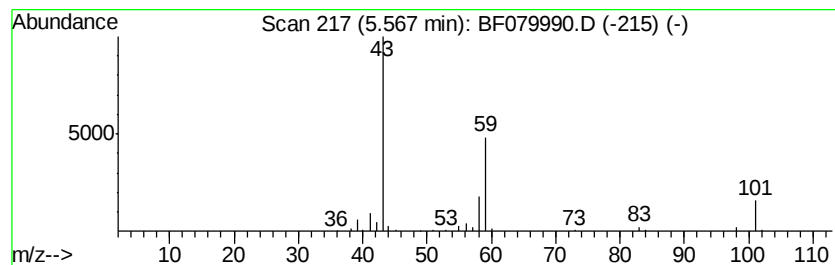
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.57	7.14 ng	99397	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			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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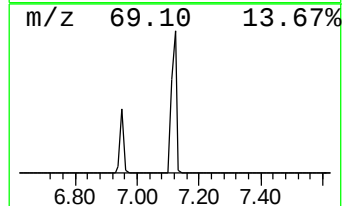
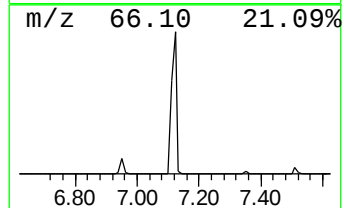
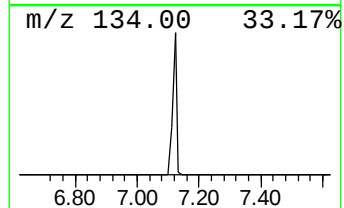
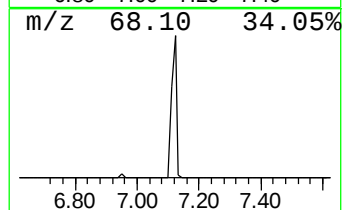
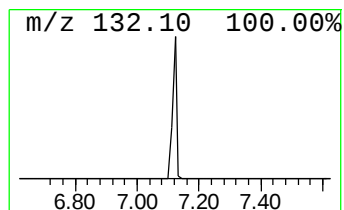
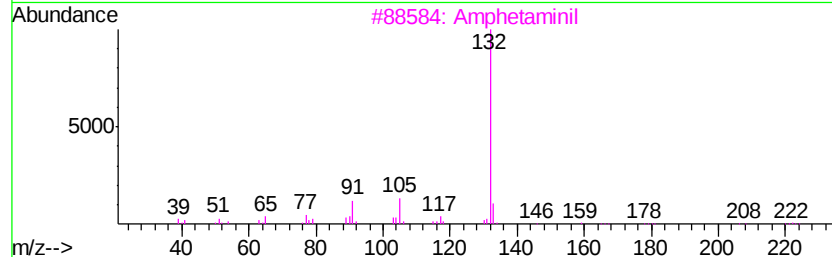
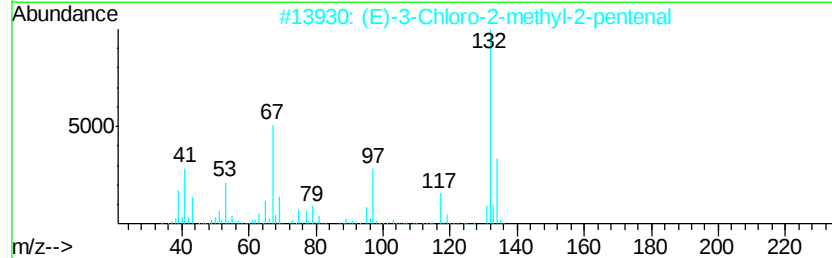
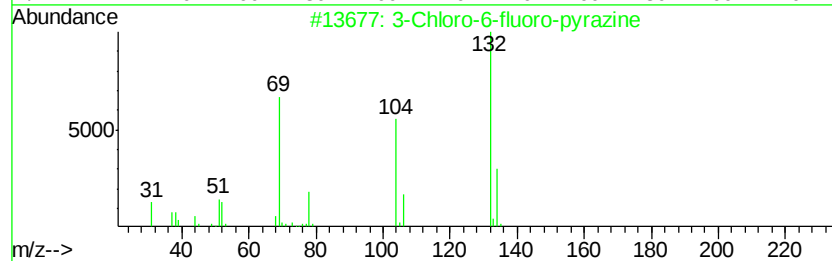
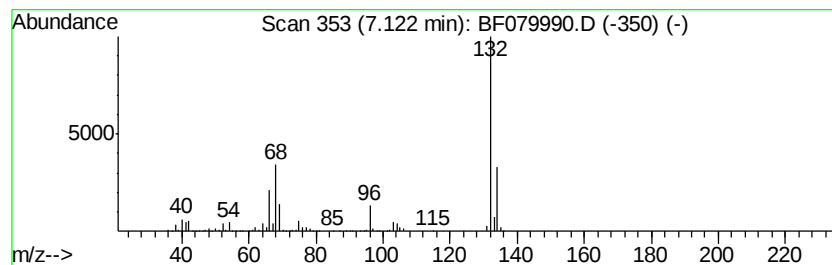
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Peak Number 2 unknown7.12 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	82.88 ng	1153550	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		Xenon	132	Xe	007440-63-3	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.57	7.1	ng	99397	1	7.35	278369	20.0
unknown7.12	7.12	82.9	ng	1153550	1	7.35	278369	20.0