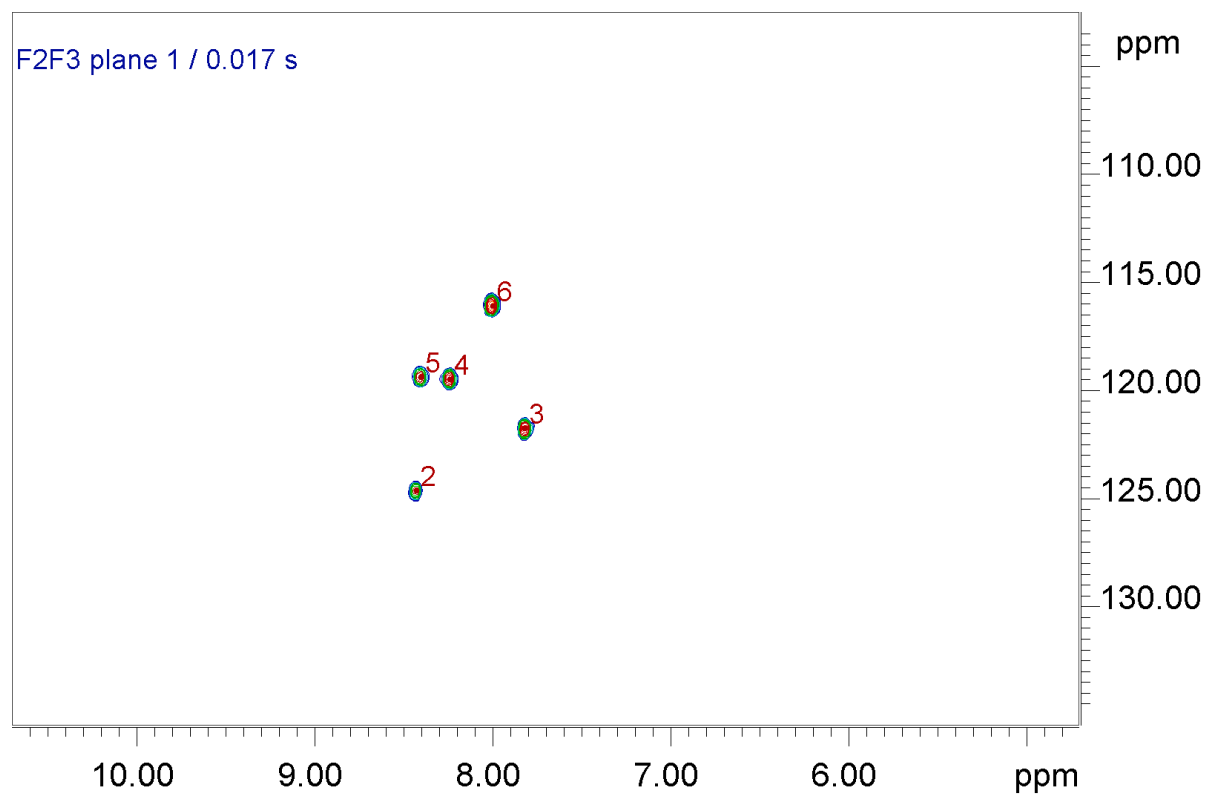


## ● T2 Analysis



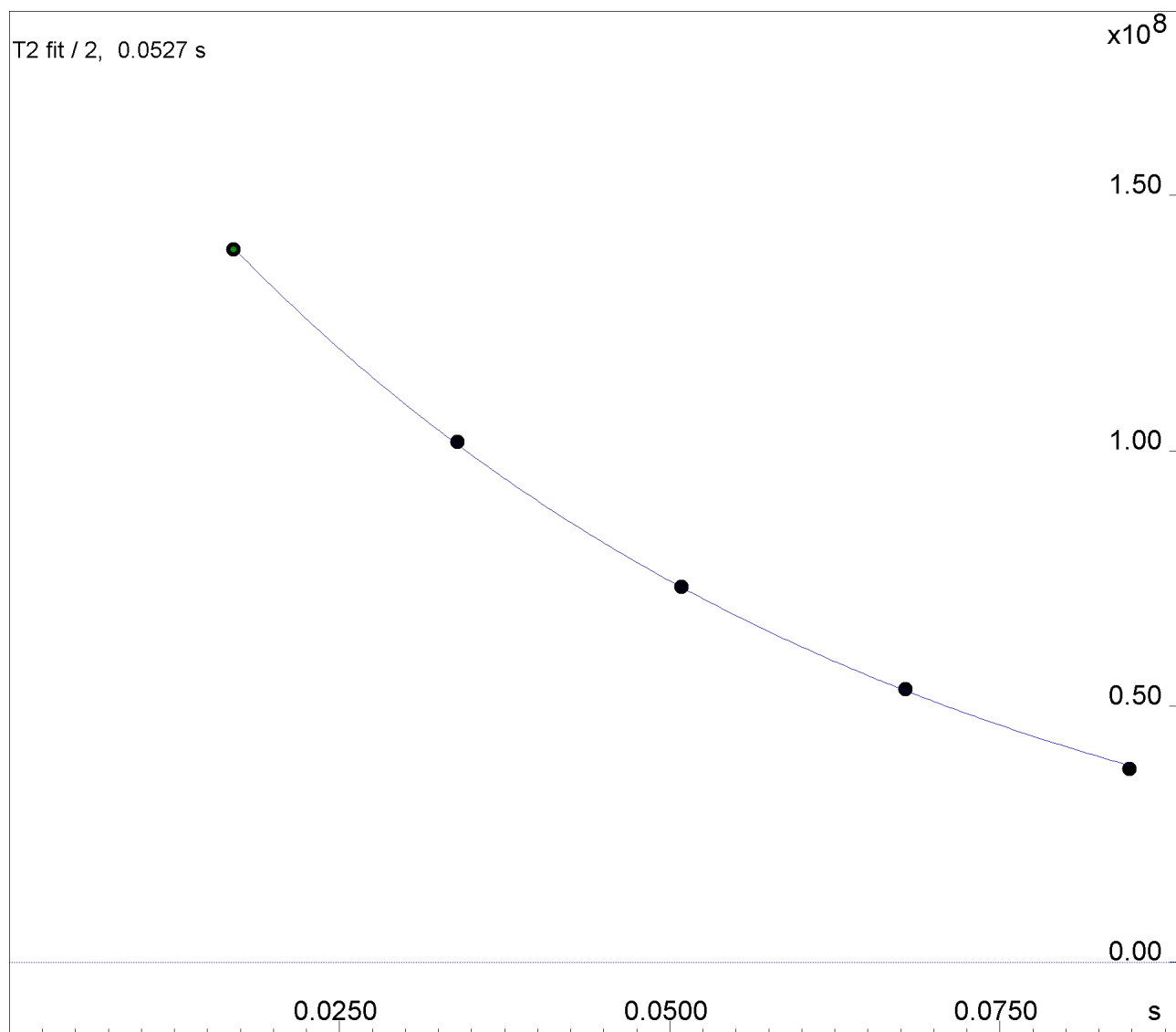
sample name:	ubiquitine
Description/Title:	standard demo sample
Origin:	in-house
Date of preparation:	06 / 2005
Solvent:	D2O
pH:	7.0
Sample tube:	normal
Tube diameter (mm):	3
Concentration (mM):	10.00
Temperature (K):	308.0
Weight (Daltons):	8900
Correlation time (ns):	12.40
Labelling:	15N
AminoAcid storage:	SEQ file
AminoAcid file:	/opt/data/nmrsu/nmr/sample/ubiquitine.seq
PDB file:	/opt/data/nmrsu/nmr/sample/ubiquitine.pdb



Fitted function:	$f(t) = I_0 * \exp(-t/T)$
Random error estimation of data:	RMS per spectrum (or trace/plane)
Systematic error estimation of data:	worst case per peak scenario
Fit parameter Error estimation method:	from fit using arbitrary y uncertainties
Confidence level:	95%
Used peaks:	automatically picked peaks
Used integrals:	peak intensities
Used Time:	all values (including replicates) used

Peak name	F1 [ppm]	F2 [ppm]	T2 [s]	error	fitInfo
2	124.625	8.425	0.0527	0.001285	Done
3	121.756	7.816	0.0622	0.008432	Done
4	119.467	8.238	0.0527	0.01155	Done
5	119.371	8.402	0.0507	0.008522	Done
6	116.083	7.998	0.0624	0.01382	Done

## Current fit display



# T2 Analysis

/home/nmrsu/Samuli/Cory\_Peptides\_2022/magainine\_POPC\_bic\_200mM\_37deg/120/pdata/1/3rrr

