Measuring the roughness of protein molecules using the single molecule unfolding method

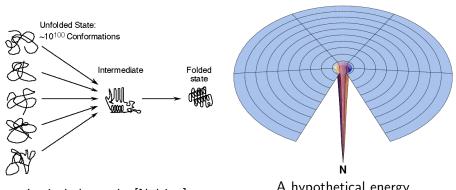
Trevor King

Drexel University

June 7, 2007

Finding stable protein conformations

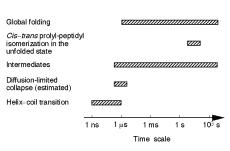
Exhaustive searches are impossible.



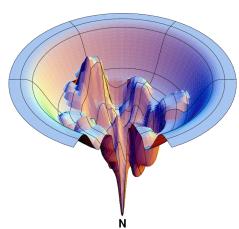
Levinthal paradox[Nolting].

A hypothetical energy landscape[Dill].

But natural proteins do fold

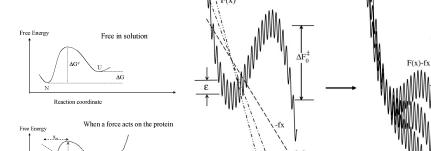


Protein folding timescales at 25°C and pH 7[Nolting].



How do we measure this energy landscape?

Measuring the energy landscape roughness ε



Proposed energy roughness mode [Hyeon]. $F(x) = F_0(x) + F_1(x).$

Reaction coordinate

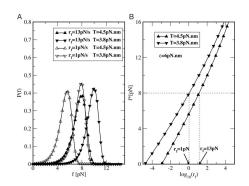
x=D

Kramers' problem and diffusion in rough potentials

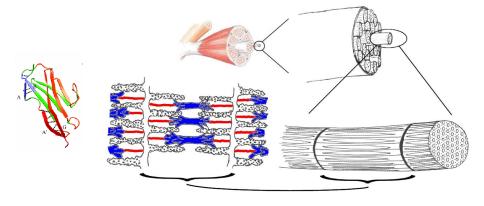
$$\varepsilon^2 \approx \frac{k_B^2 T_1 T_2}{T_2 - T_1} \cdot \left[T_1 \log \left(\frac{r_f(T_1) \Delta x(f^*)}{k_0(f^*) k_B T_1} \right) - T_2 \log \left(\frac{r_f(T_2) \Delta x(f^*)}{k_0(f^*) k_B T_2} \right) \right],$$

where

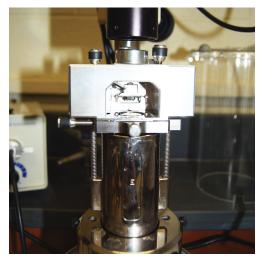
- $\Delta x(f^*)$ is the unfolding distance from the bound state to the transition state for the smooth potential $F_0(x) f^*(x)$.
- $k_0(f^*)$ is the escape rate from the smooth potential.



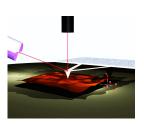
Titin and temperature dependence in muscle



Atomic force microscopy



MultiMode head.

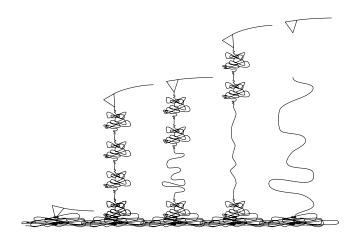


Operation[Warr].



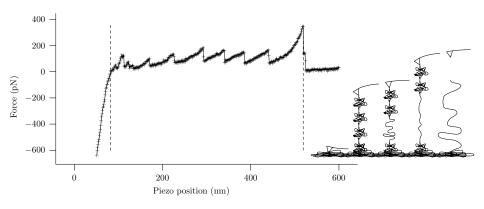
Fluid cell.

Protein unfolding, understanding the mechanics





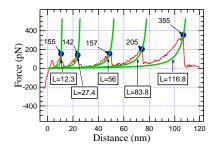
Protein unfolding, understanding the data



The characteristic unfolding 'sawtooth'.

Worm like chains

To convince ourselves that we're measuring what we think we're measuring



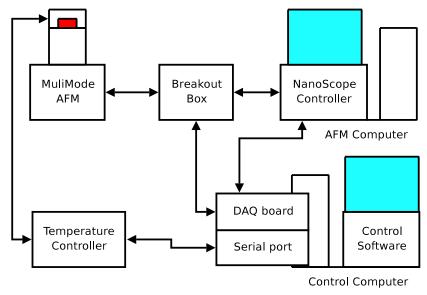


$$F(x) = \frac{k_B T}{p} \left[\frac{1}{4} \left(\frac{1}{(1 - x/L)^2} - 1 \right) + \frac{x}{L} \right],$$

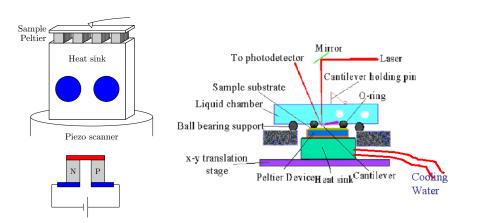
where p is the persistence length and L is the contour length.



Equipment layout

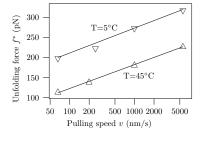


Temperature control



Temperature control module and the Peltier heater/cooler.

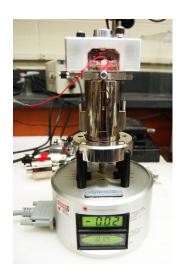
Temperature dependent unfolding in Ubiquitin



The unfolding energy according to Hyeon and Thirmulai's equation is

$$\varepsilon \approx 1.98 \cdot 10^{-20} \text{J} \approx 4.8 k_B T.$$

Streamlining the procedure



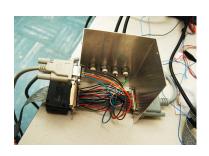
Difficulties with temperature dependent measurement

- Thermal drift and limited piezo range.
- Procedure not very reliable, needs hand tweaking.
- Many parameters to optimize once it is working.
- Lots of data to search through to find our sawtooth.

Goals

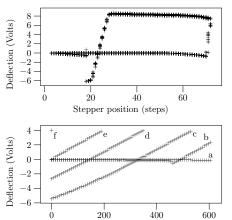
- Automate stepping motor control.
- Re-architect experiment software.
- Automate force-curve sorting.

Breaking into AFM communication

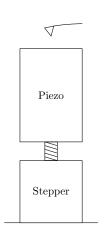




Testing our motor control



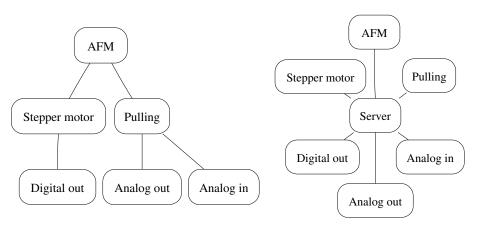
Piezo position (nm)



Measuring backlash (\sim 46 steps) and stepsize (\sim 170nm).

Architecture comparison

Choose for flexibility with minimal programmer time.



Tree vs. network.



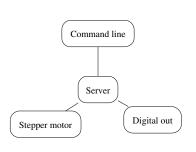
The goal: Flexible control scripts in action

```
> approach_surface
Lost signal!
> approach_surface
> piezo_measure_distance
372 nm
> pull_protein --save 'D:\data\' &
saving data to 'D:\data\070510\'
> monitor_unfolding
> sort data --dir 'D:\data\070510\',
. . .
> 1s
pulls_good
pulls_bad_structure
. . .
```

Example script: approach_surface

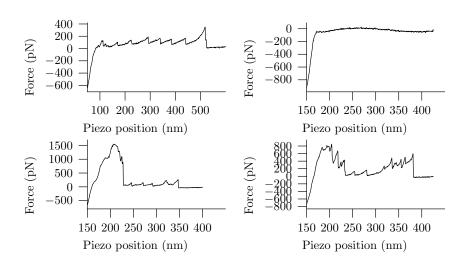
```
#!/bin/sh
                                    Minimize programmer time
TARG DEF=1 # Volts
MIN TOT=3 # Volts
get_feedback() {
  DEF='cmmdline -c':AFM|Get:Deflection'
  TOT='cmmdline -c':AFM|Get:TotalPhotodiode'
step_closer() {
  ANS='cmmdline -c 'Stepper|Set:RelPos 1'
get_feedback
while [ $DEF -lt $TARG_DEF ] && [ $TOT -gt $MIN_TOT ]; do
  step_closer
  get_feedback
done
```

Tracing commands



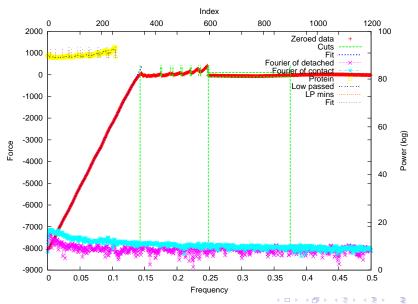
| From | То | Message | |
|----------|----------|--------------|--|
| Cmmdline | Stepper | Set:AbsPos 3 | |
| Stepper | Dig. Out | Set:Output 1 | |
| Dig. Out | Stepper | DONE | |
| Stepper | Dig. Out | Set:Output 2 | |
| Dig. Out | Stepper | DONE | |
| Stepper | Dig. Out | Set:Output 3 | |
| Dig. Out | Stepper | DONE | |
| Stepper | Cmmdline | DONE | |

Automatically sorted pulling curves

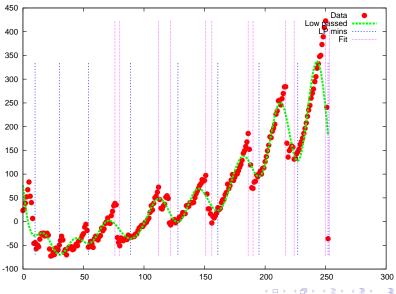




Sorting example



Sorting example



Preliminary results

| % | Good | Bad | % | Good | Bad |
|---|------|-----|---|------|------|
| + | 1.9 | 23 | + | 2.2 | 3.9 |
| _ | 0.3 | 75 | _ | 2.2 | 91.5 |

The sorting algorithm on training and test datasets.

Why are they so different? What do we do about it?

Conclusions

We have

- Automated control of the stepping motor, which will allow longer data runs without interruption.
- Developed a flexible software architecture, which will ease the implementation of future refinements and reduce troubleshooting time during setup.
- Beginning development of an automatic force curve sorter, which will allow automatic experiment parameter optimization and ease the sorting of large datasets.

The next step will be putting these pieces together in order to measure Titin's temperature dependent unfolding properties.

