

# A MECHANISM FOR FREE-ENERGY TRANSDUCTION BY MOLECULES

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Wednesday, February 16, 2011

4:00 p.m. NSH 118

(Refreshments at 3:30 p.m. NSH 202)

Proteins and other macromolecules can act as molecular machines that convert energy from one form to another through cycles of conformational transitions. In a macroscopically fluctuating environment or at the single-molecule level, the probability for a molecule to be in any state  $j$  fluctuates, and the probability current from any other state  $i$  to state  $j$  is given as the sum of a steady-state current and a pumped current,  $J_{ij} = J_{ij}^{ss} + J_{ij}^{p}$ , where  $J_{ij}^{ss}$  is the fraction of the fluctuating current into and out of state  $j$  coming directly from state  $i$ , and  $J_{ij}^{p}$  is the rate of change of the probability for the molecule to be in state  $j$ . If the fluctuations arise from an equilibrium source, microscopic reversibility guarantees that the time average of the pumped current is zero. If, however, the fluctuations arise due to the action of a nonequilibrium source, the time average of the pumped current is not in general zero and can be opposite in sign to the steady-state current. The pumped current provides a mechanism by which fluctuations, whether generated externally or arising from an internal nonequilibrium chemical reaction, can do electrical, mechanical, or chemical work on a system by coupling into the equilibrium conformational transitions of a protein. In this review I examine work elaborating the mechanism of stochastic pumping and also discuss a thermodynamically consistent approach for modeling the effects of dynamic disorder on enzymes and other proteins.

Colloquium

All interested  
persons are  
cordially  
invited to  
attend.