Investigating physical constraints underlying catalysis and their impact on metabolic systems

Uri Barenholz

CRI Research Symposium

October 11, 2017

https://git.io/vd2x0

Once upon a time...

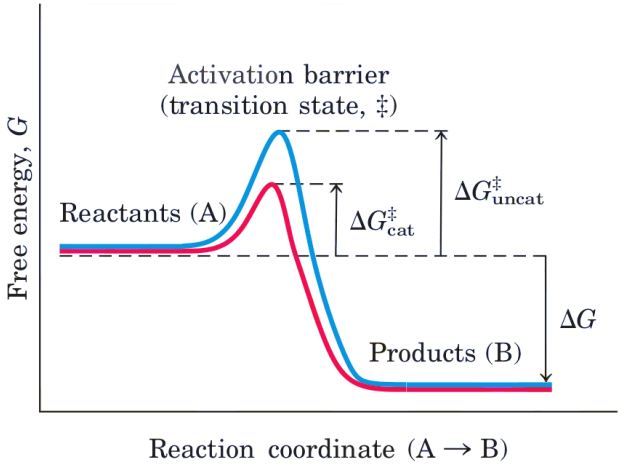
Research questions

- What is the physical limit for lowering the activation energy barrier of a given reaction

Research questions

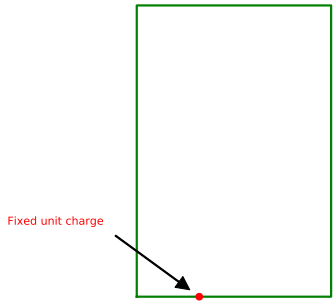
- What is the physical limit for lowering the activation energy barrier of a given reaction
- How is the affinity of an enzyme affected by the requirement to be selective

Textbook illustration

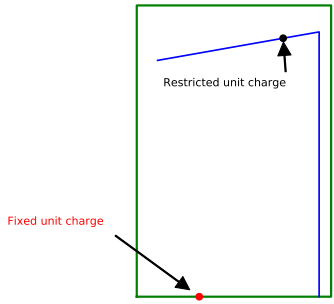


https://git.io/vd2x0

Modeling energy landscape modification in a classical system

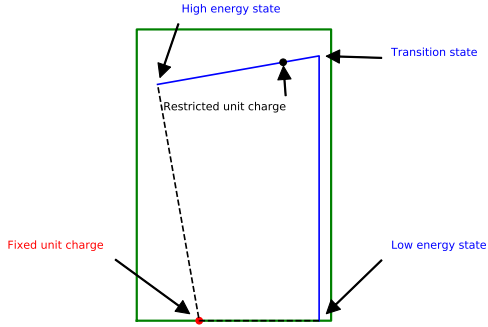


Modeling energy landscape modification in a classical system



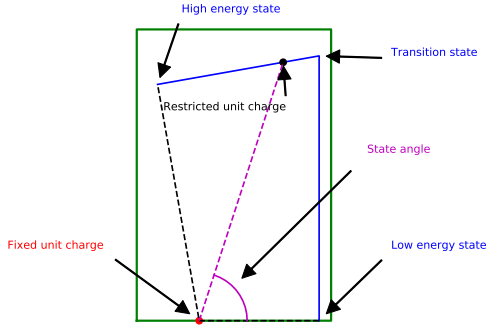
https://git.io/vd2x0

Modeling energy landscape modification in a classical system



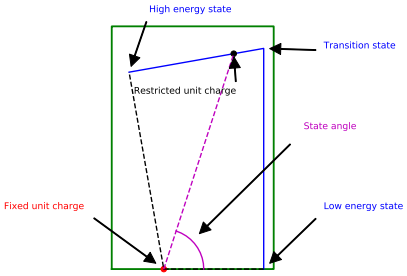
https://git.io/vd2x0

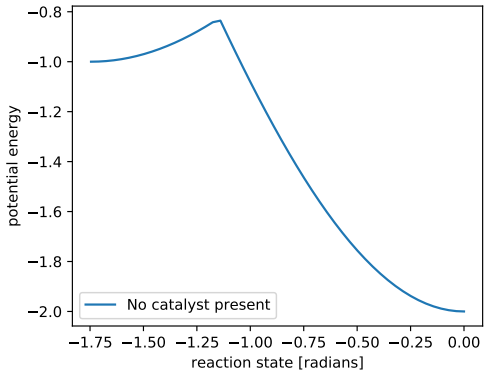
Modeling energy landscape modification in a classical system



https://git.io/vd2x0

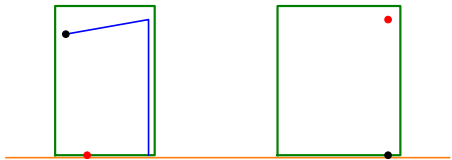
Reaction energy landscape of model substrate





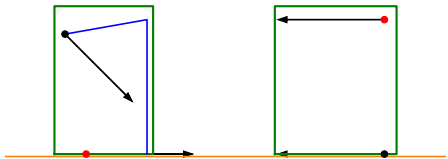
https://git.io/vd2x0

Introducing a model catalyst

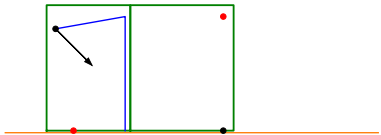


https://git.io/vd2x0

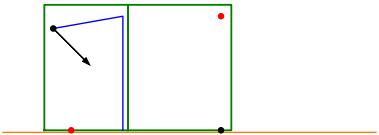
Introducing a model catalyst

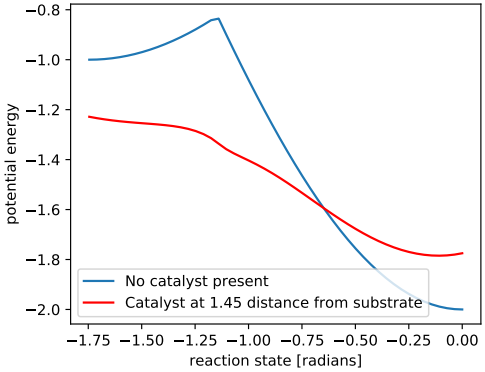


Reaction energy landscape of bound substrate



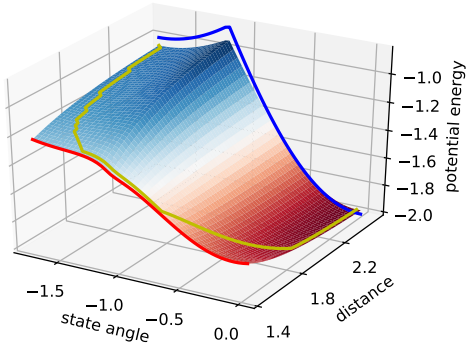
Reaction energy landscape of bound substrate





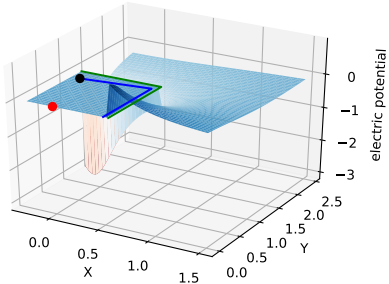
https://git.io/vd2x0

The catalyst creates a bypass to the energy barrier at the transition state



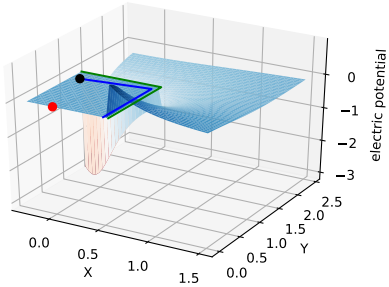
https://git.io/vd2x0

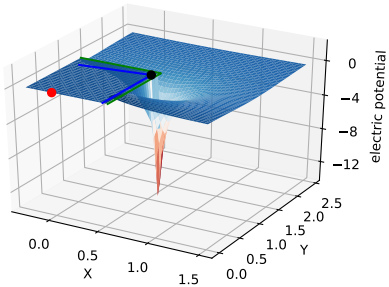
Subtracting the potential field at the transition state from the initial state produces an energy barrier reduction landscape



https://git.io/vd2x0

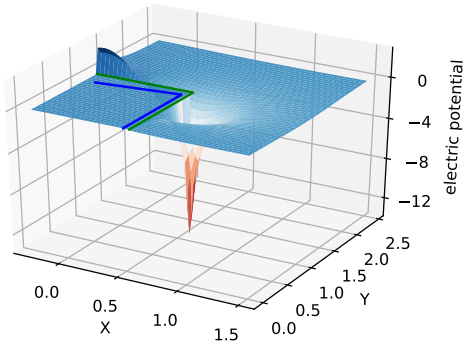
Subtracting the potential field at the transition state from the initial state produces an energy barrier reduction landscape





https://git.io/vd2x0

Subtracting the potential field at the transition state from the initial state produces an energy barrier reduction landscape



https://git.io/vd2x0

Subtracting the potential field at the transition state from the initial state produces an energy barrier reduction landscape

- The resulting function quantifies the barrier reduction when positioning a positive point charge at any coordinate in space

Subtracting the potential field at the transition state from the initial state produces an energy barrier reduction landscape

- The resulting function quantifies the barrier reduction when positioning a positive point charge at any coordinate in space
- Placing charges at extremum points of this function achieves maximal barrier reduction

Methodological approach for investigating catalytic constraints

Methodological approach for investigating catalytic constraints

- Crowd-sourcing platform
  - Challenge existing assumptions
  - Reveal potential catalytic mechanisms

Methodological approach for investigating catalytic constraints

- Crowd-sourcing platform
  - Challenge existing assumptions
  - Reveal potential catalytic mechanisms
- Handheld, functional models
  - Demonstrate and communicate catalytic principles

Methodological approach for investigating catalytic constraints

- Crowd-sourcing platform
  - Challenge existing assumptions
  - Reveal potential catalytic mechanisms
- Handheld, functional models
  - Demonstrate and communicate catalytic principles
- Apply theoretical framework to molecular domain

Methodological approach for investigating catalytic constraints

- Crowd-sourcing platform
  - Challenge existing assumptions
  - Reveal potential catalytic mechanisms
- Handheld, functional models
  - Demonstrate and communicate catalytic principles
- Apply theoretical framework to molecular domain
- Investigate metabolic network design implications
  - Synthetic biology applications
  - Origins of life metabolism

Does structural similarity limit affinity in metabolic networks?

Does structural similarity limit affinity in metabolic networks?

https://git.io/vd2x0

Does structural similarity limit affinity in metabolic networks?

- Most enzymes are substrate-specific

Does structural similarity limit affinity in metabolic networks?

- Most enzymes are substrate-specific
- Structural similarity is used for drug discovery and promiscuous activity tests

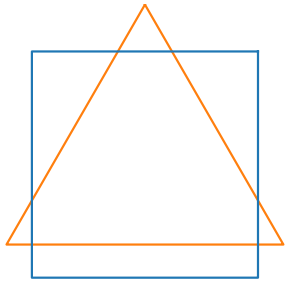
Does structural similarity limit affinity in metabolic networks?

- Most enzymes are substrate-specific
- Structural similarity is used for drug discovery and promiscuous activity tests
- Metabolic networks must contain structurally similar metabolites
  - But can potentially reduce similarities at critical points

Does structural similarity limit affinity in metabolic networks?

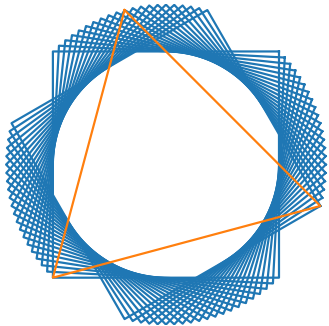
- Most enzymes are substrate-specific
- Structural similarity is used for drug discovery and promiscuous activity tests
- Metabolic networks must contain structurally similar metabolites
  - But can potentially reduce similarities at critical points
- Numerous examples for specificity tradeoffs in the literature

Why do we expect selectivity to decrease affinity?



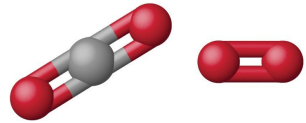
https://git.io/vd2x0

Why do we expect selectivity to decrease affinity?



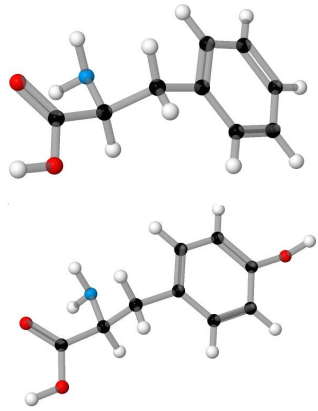
Examples of specificity-affinity challenges

- RuBisCo
  - CO_2 versus O_2



Examples of specificity-affinity challenges

- RuBisCo
  - CO_2 versus O_2
- Tyrosine ammonia lyase
  - Tyr versus Phe



https://git.io/vd2x0

Examples of specificity-affinity challenges

- RuBisCo
  - CO_2 versus O_2
- Tyrosine ammonia lyase
  - Tyr versus Phe
- Bacterial DNA methyltransferase
  - Relaxing sequence specificity accelerates rate
- Bacterial hexose phosphate transporter

Can we formulate a quantitative evaluation of the selectivity challenge?

- Given metabolites concentration data
  - Identify challenging reactions
  - Quantify expected cost

https://git.io/vd2x0

Can we formulate a quantitative evaluation of the selectivity challenge?

- Given metabolites concentration data
  - Identify challenging reactions
  - Quantify expected cost
- Given reaction possibilities
  - Find biases in metabolic network structure maximizing structural differences

Methodological approach for investigating selectivity tradeoffs

- Impact on metabolites concentrations and enzymes
  - BRENDA - identifying weak affinity enzymes
  - Promiscuous activity data from Sauer lab
  - Structural similarity metrics comparison with measured metabolites concentrations

Methodological approach for investigating selectivity tradeoffs

- Impact on metabolites concentrations and enzymes
  - BRENDA - identifying weak affinity enzymes
  - Promiscuous activity data from Sauer lab
  - Structural similarity metrics comparison with measured metabolites concentrations
- Impact on network structure
  - Project metabolic networks to chemical space
  - Implement selectivity in constraint based modeling of metabolic networks

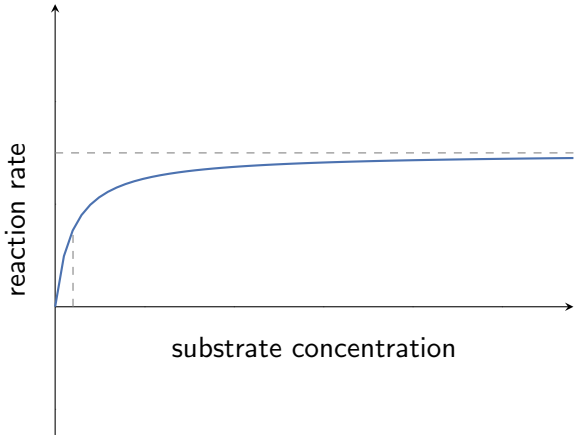
Summary

- Basic challenges of biological systems are rarely investigated theoretically
- Transforming key problems to simplified models in accessible platforms can leverage innovation of wider audience and reveal novel principles
- Recently available datasets allow evaluation of hypotheses
- Mapping metabolic networks into the chemical space can highlight metabolic network motifs

Thank You!

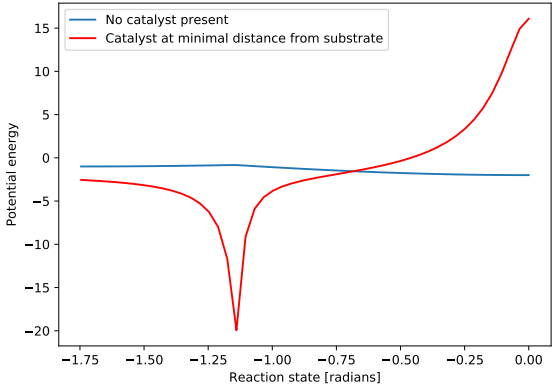
https://git.io/vd2x0

The Michaelis-Menten model for catalyzed chemical reaction rate



$$V = \frac{k_{\text{cat}}[X]}{k_M + [X]}$$

Catalyst design must track the entire reaction pathway



https://git.io/vd2x0

Structural similarity inhibits enzymes due to finite structural diversity