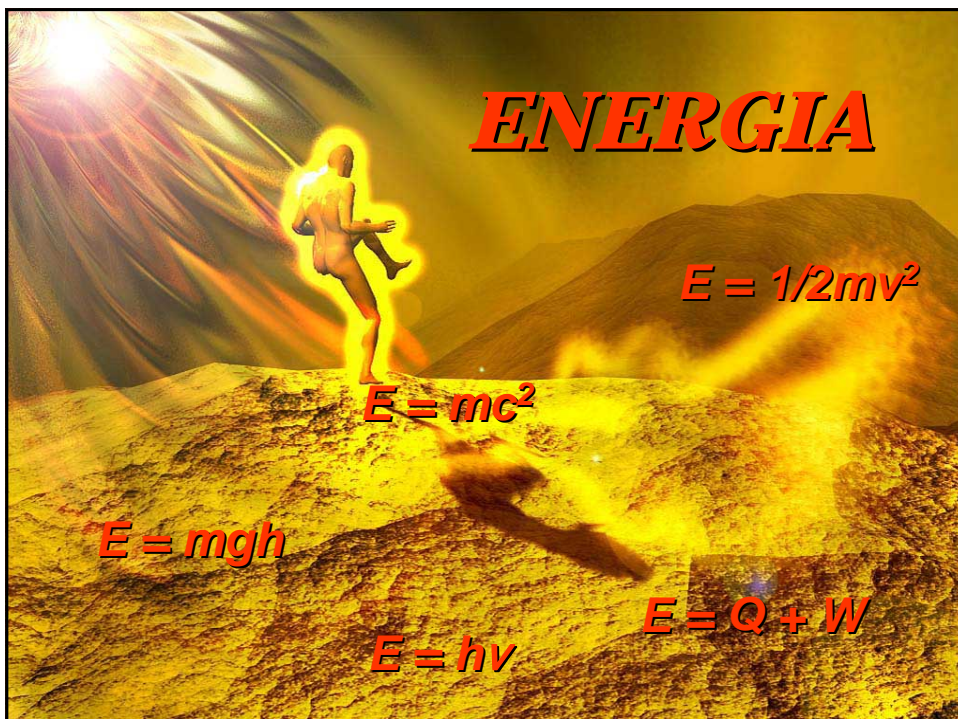


Programa Académico de Bachillerato
Universidad de Chile
Curso Biología



**Clase: Conceptos generales de termodinámica.
Bioenergética.**

*Dr. Enrique Castellón
Programa de Fisiología y Biofísica
Instituto de Ciencias Biomédicas
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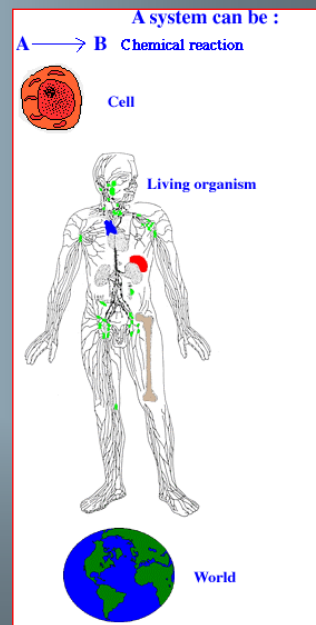


Sistemas y entorno

Hermético (No hay intercambio de energía ni materia)

Cerrado (Hay intercambio de energía pero no de materia)

Abierto (Hay intercambio de energía y materia)



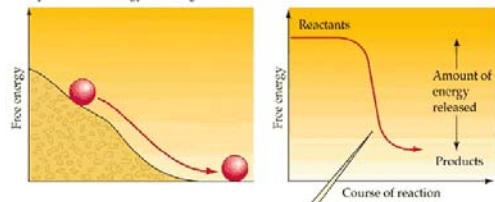
$$\Delta G = \Delta H - T\Delta S$$

$\Delta G < 0$ Exergónica / Espontánea

$\Delta G > 0$ Endergónica / No Espont.

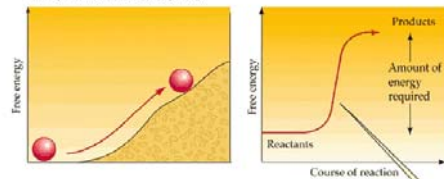
$\Delta G = 0$ Equilibrio termodinámico

(a) Exergonic reaction
(spontaneous; energy-releasing)



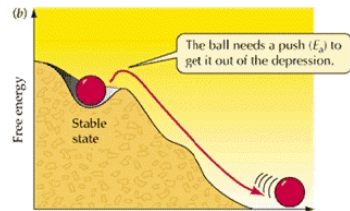
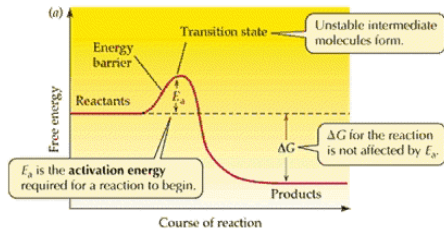
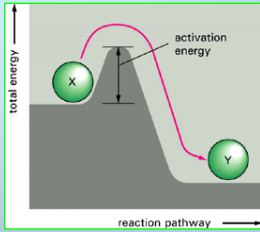
In an exergonic reaction, energy is released as reactants form products.

(b) Endergonic reaction
(not spontaneous; energy-requiring)

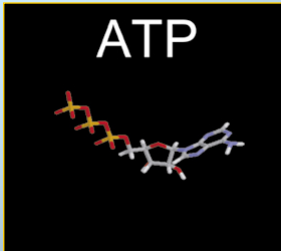


Energy is required for an endergonic reaction, in which reactants with a low energy content are converted to products with a higher energy level.

CATALYSIS



ATP



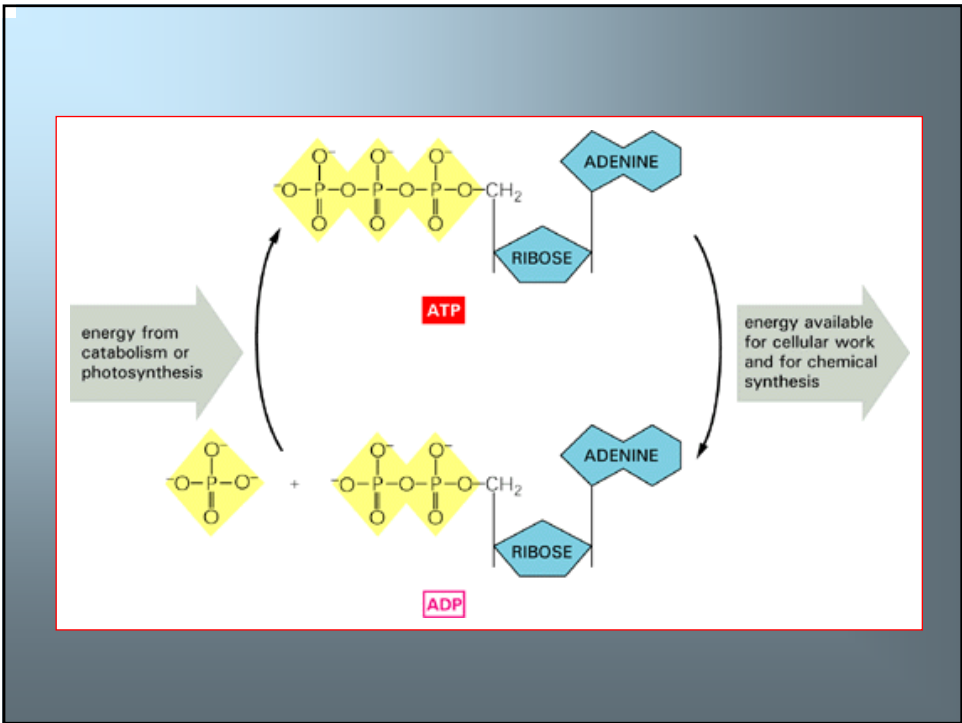
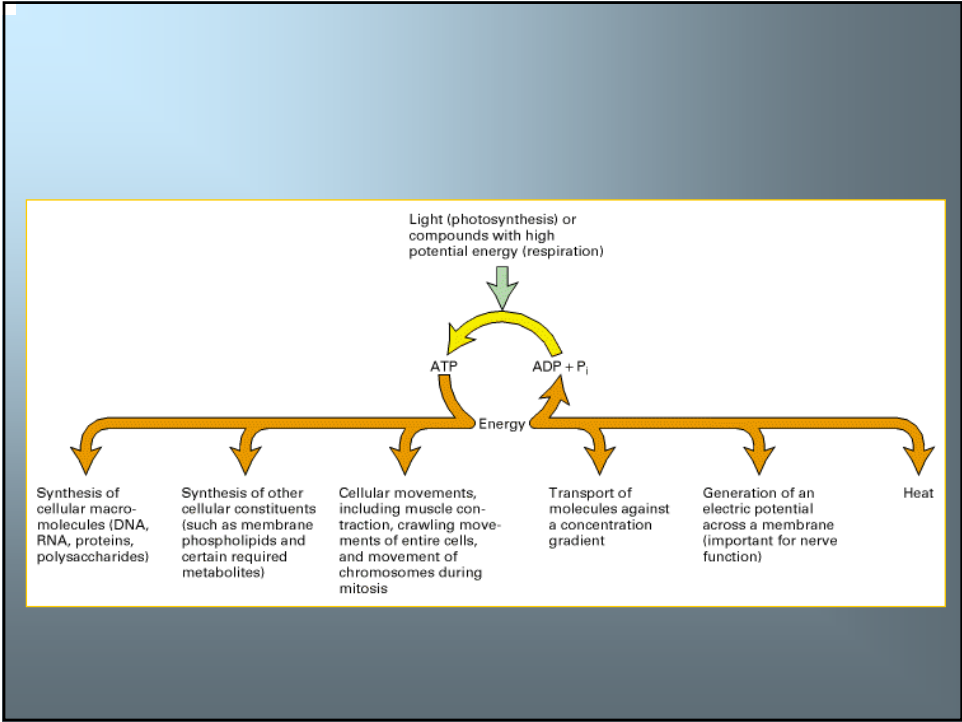
ATP - LIKE A RECHARGEABLE BATTERY

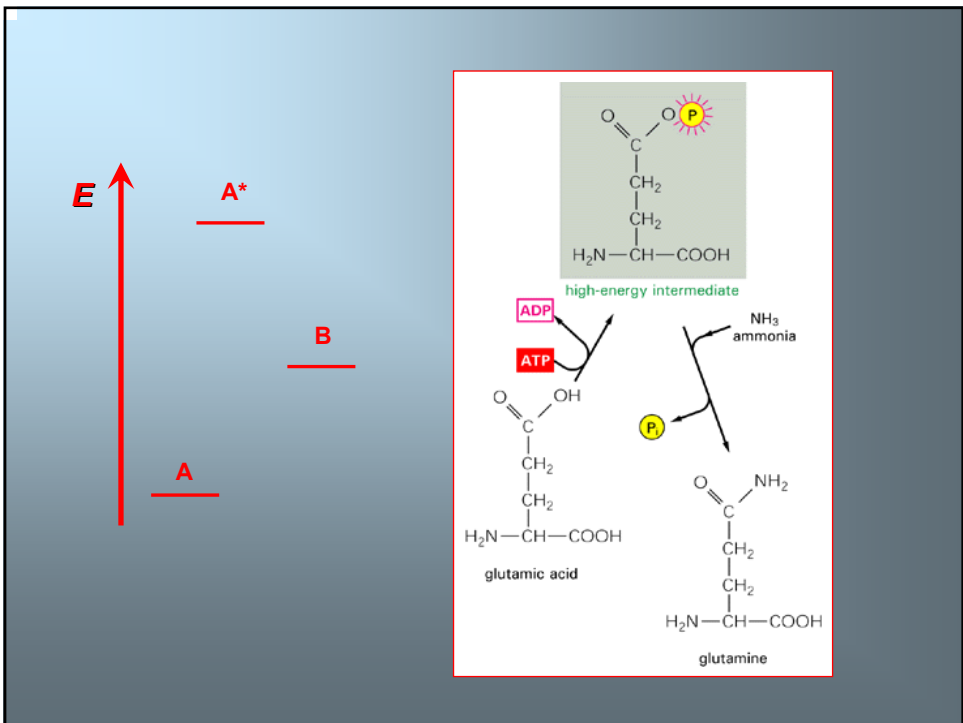
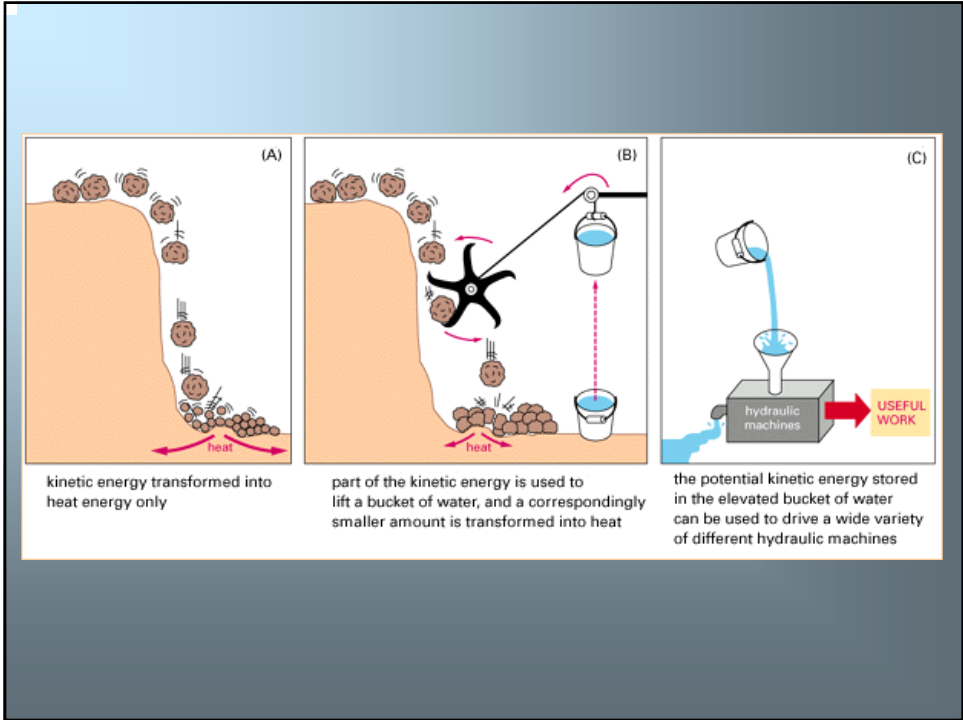


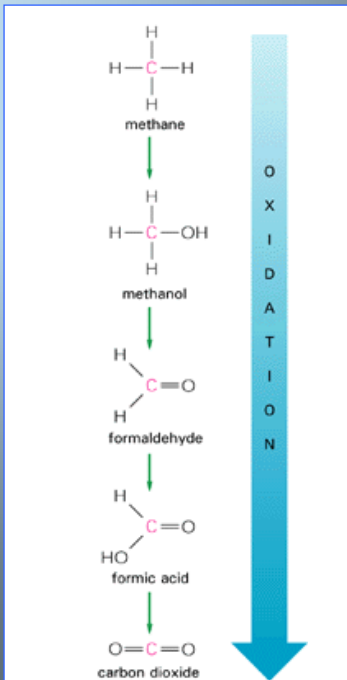
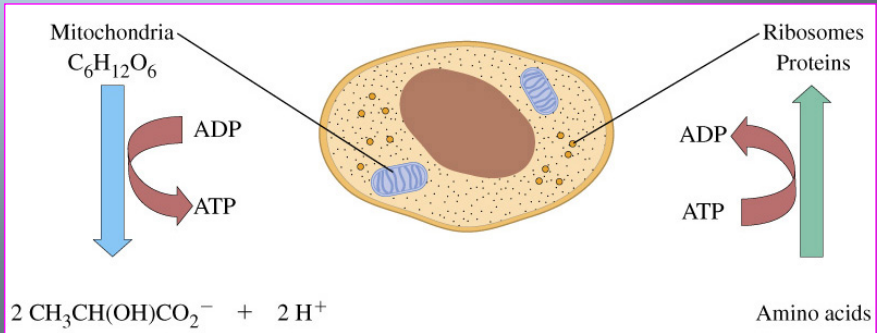
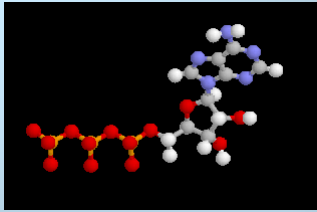
TIME FOR A RECHARGE

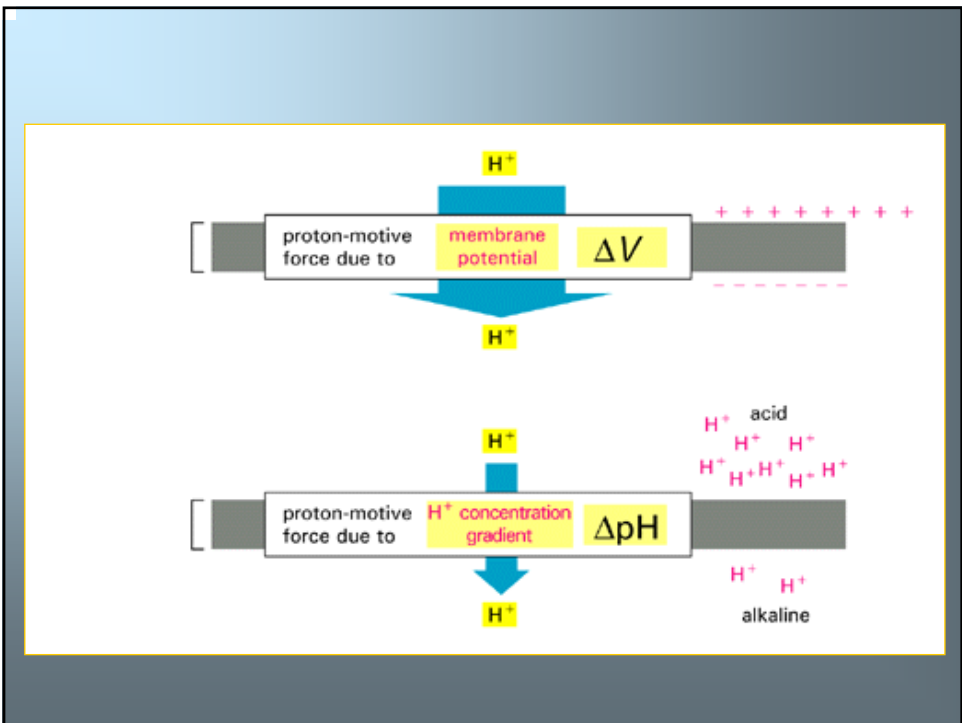
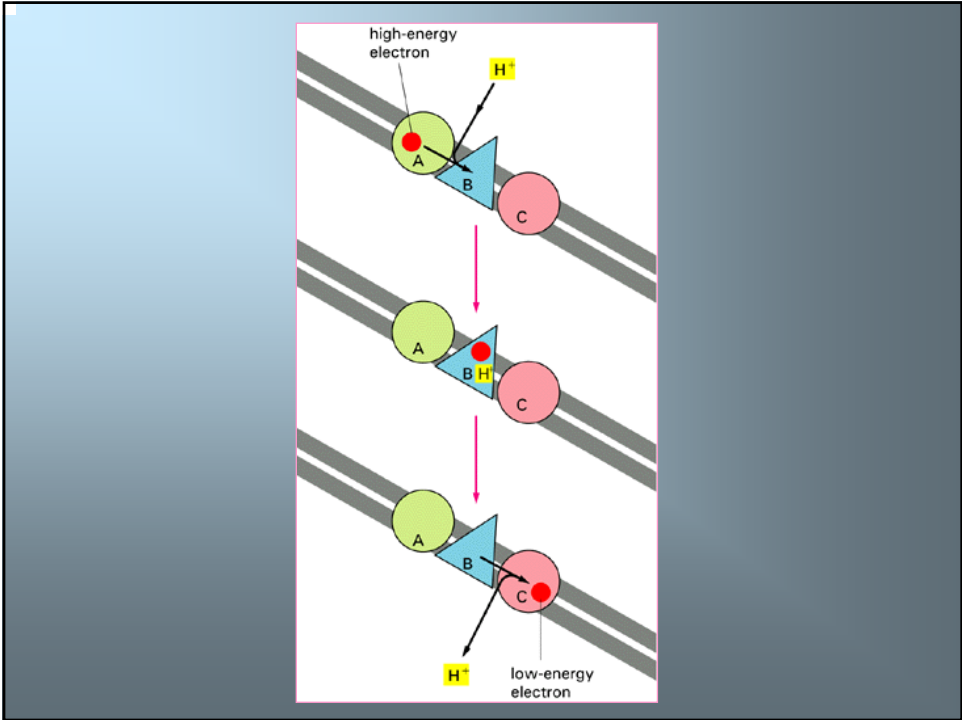


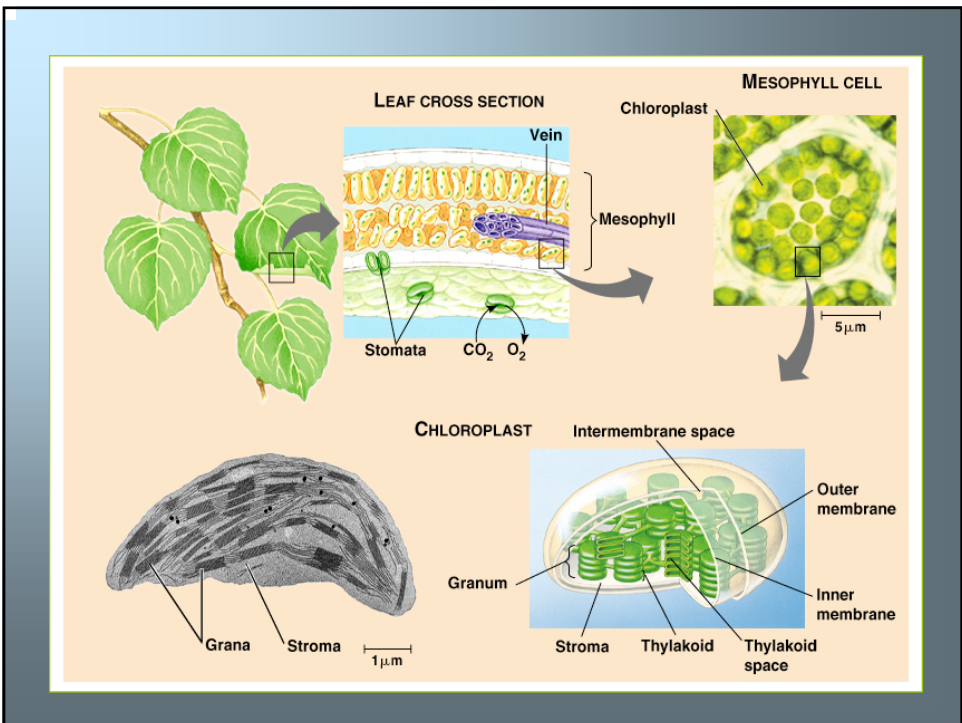
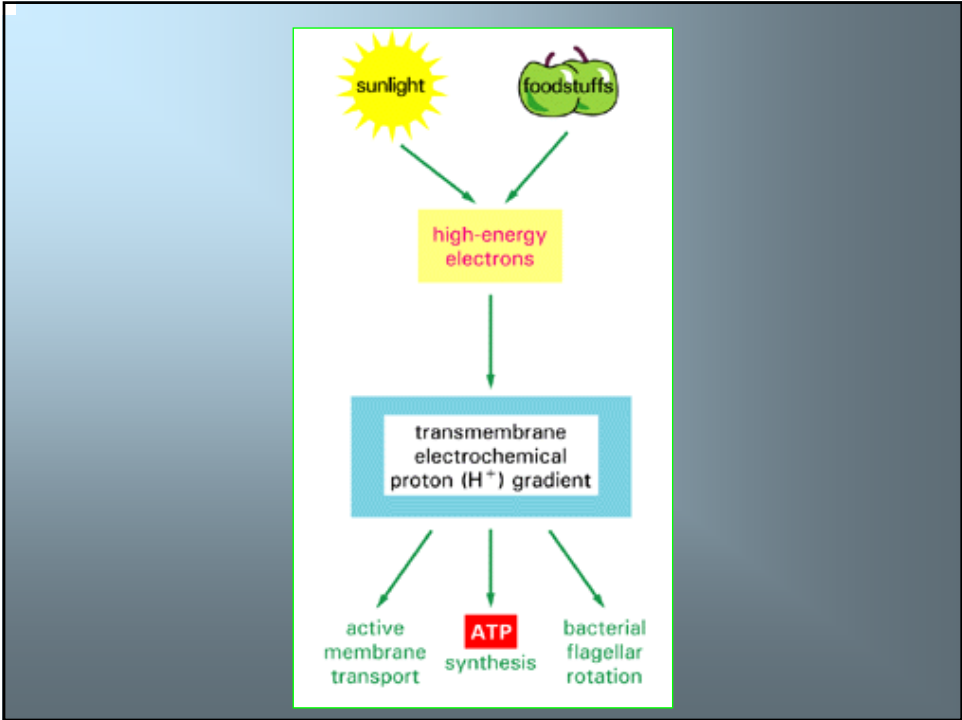
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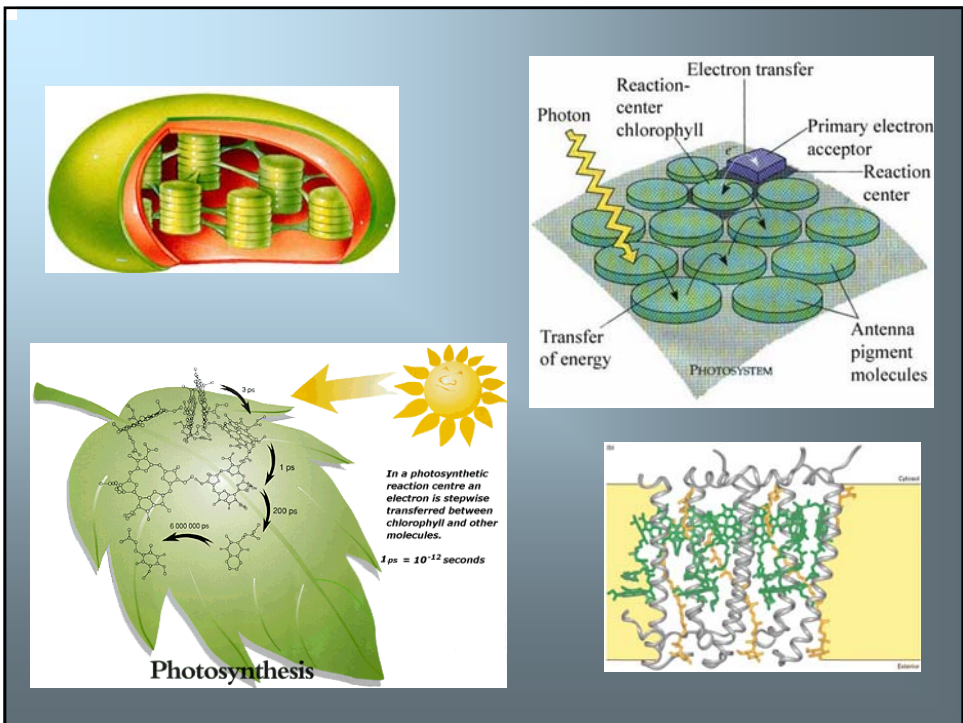
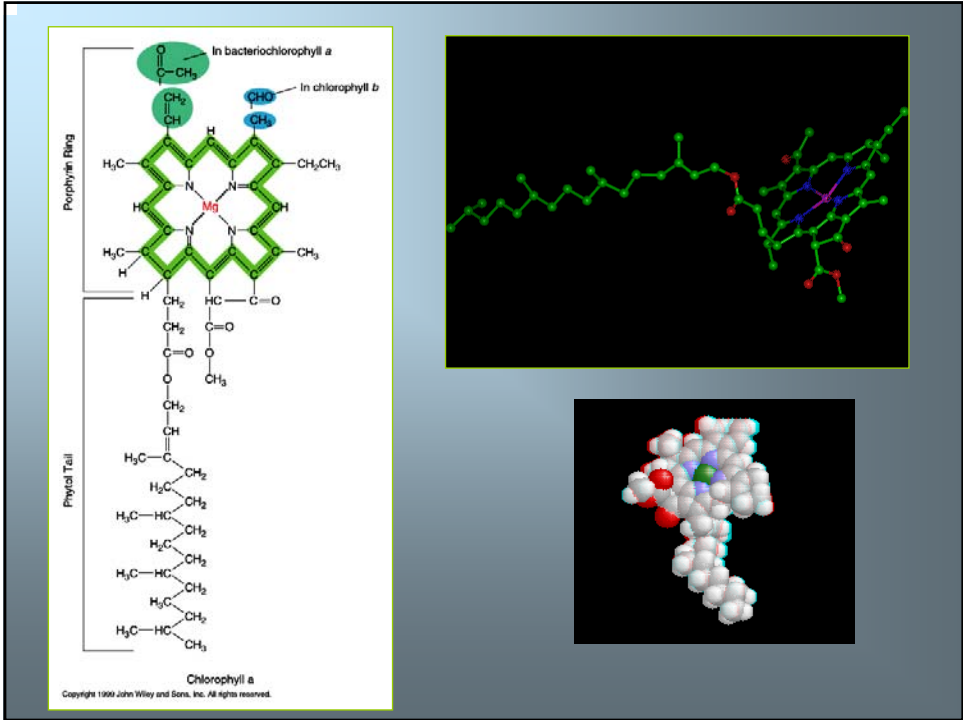


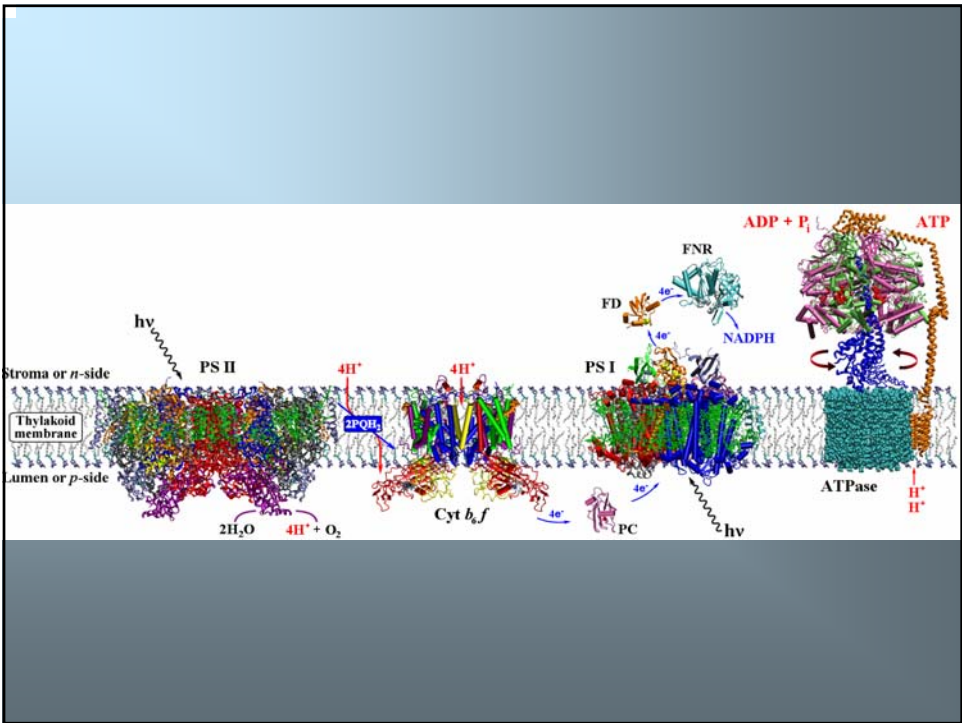
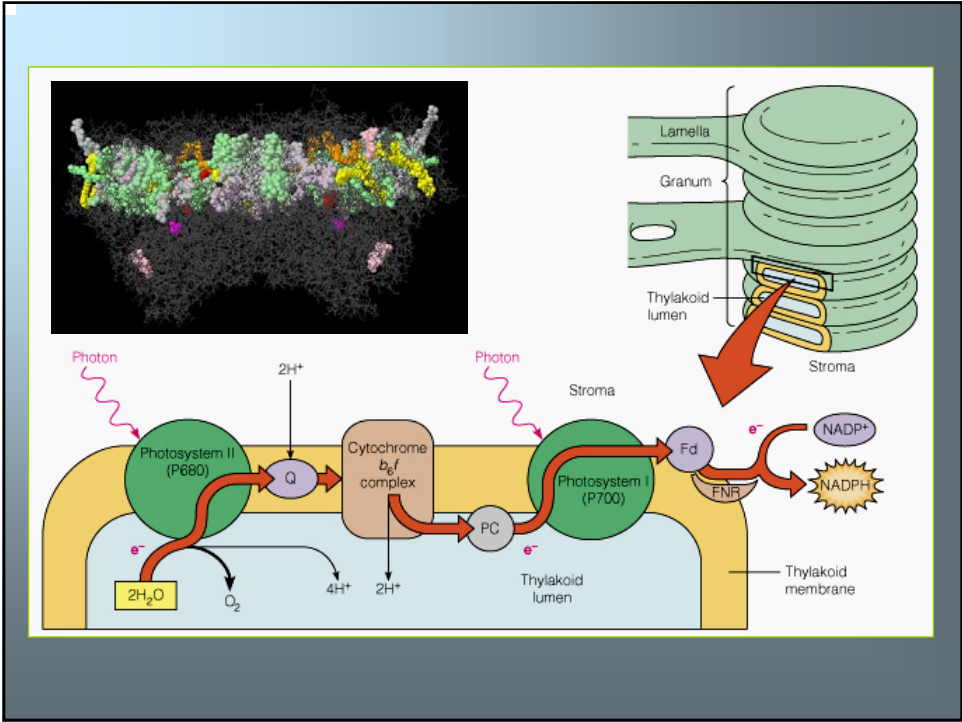


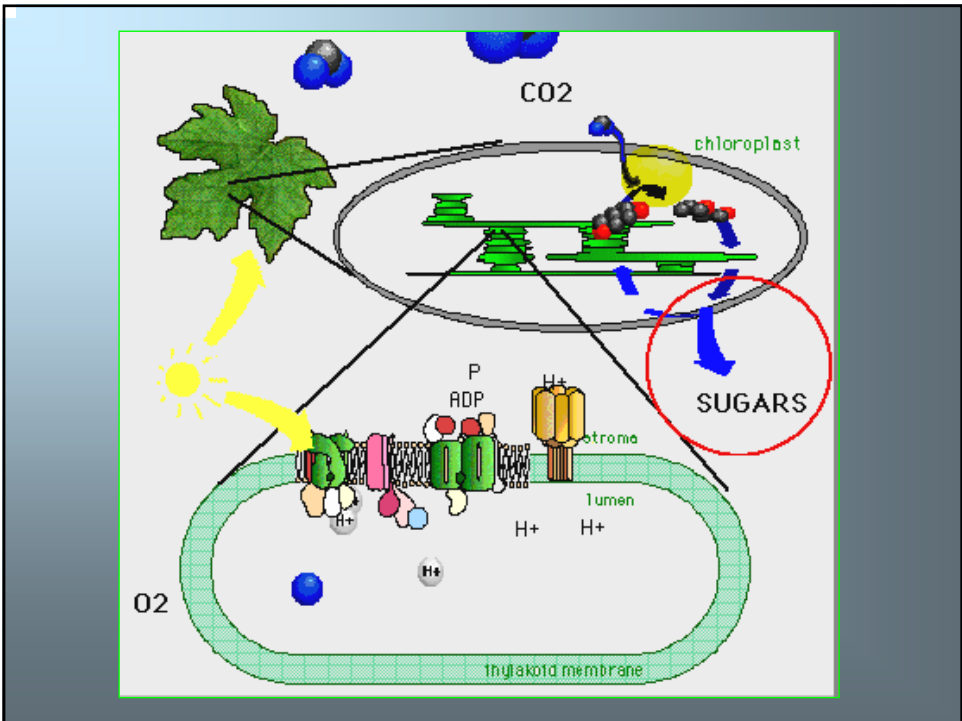
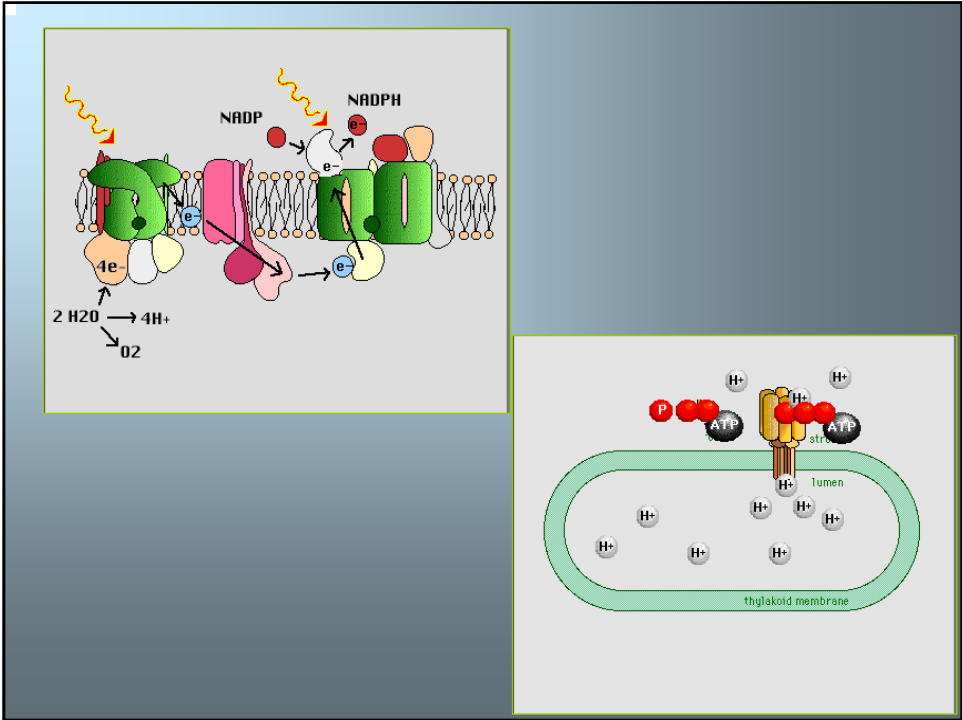


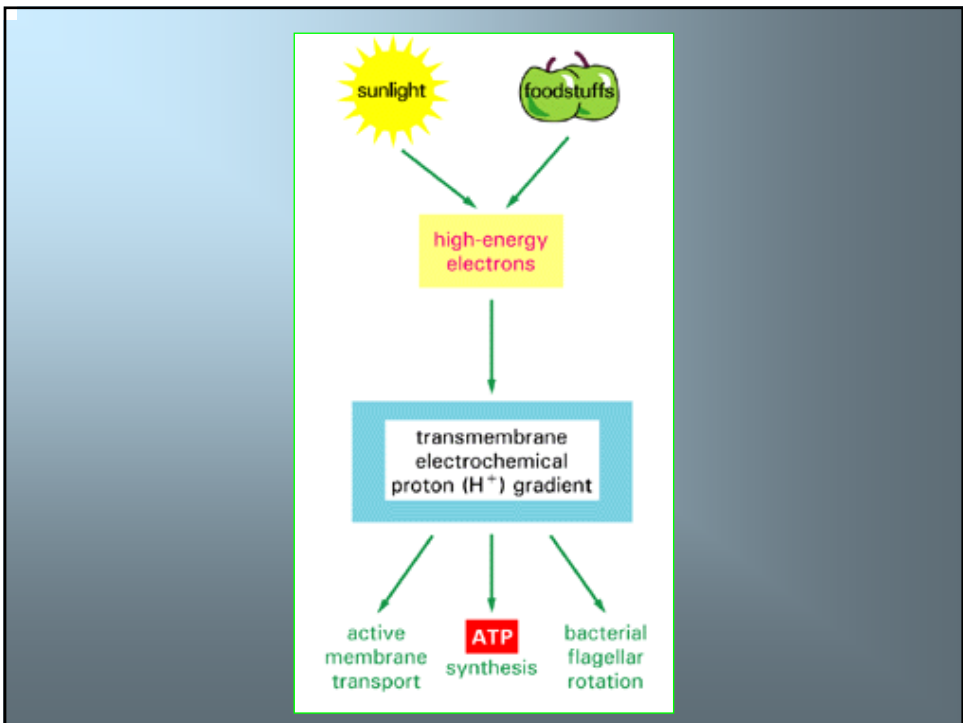
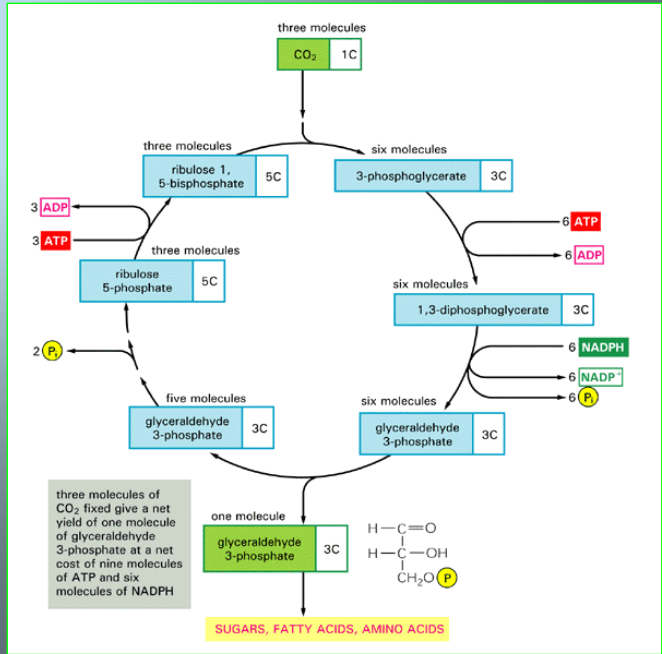
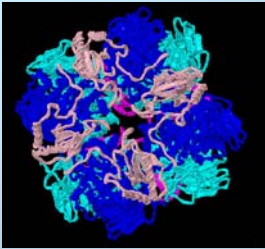


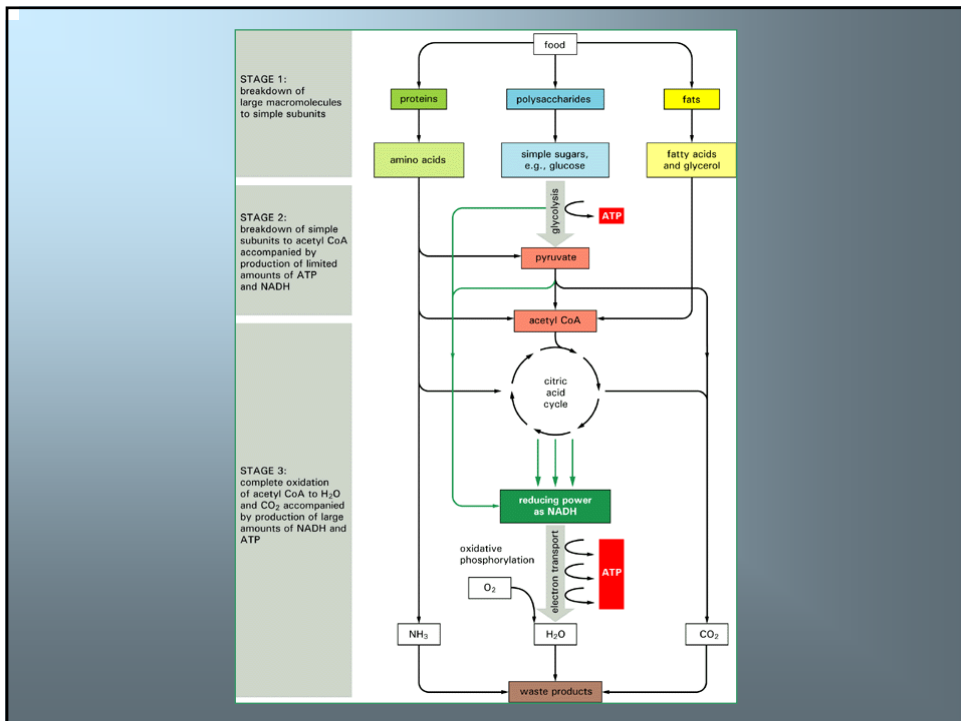
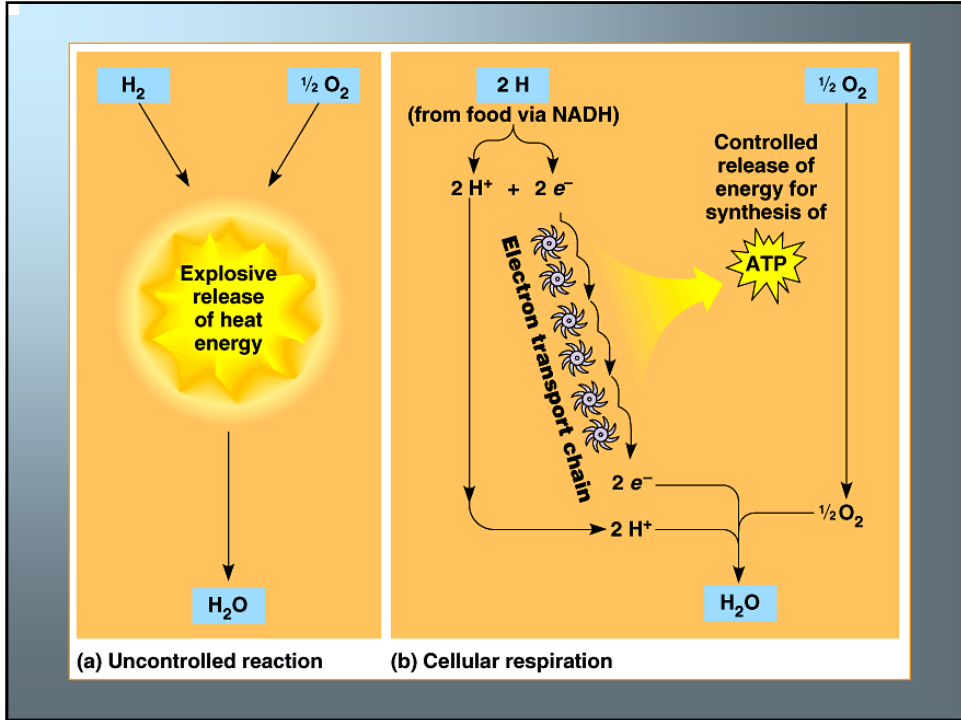


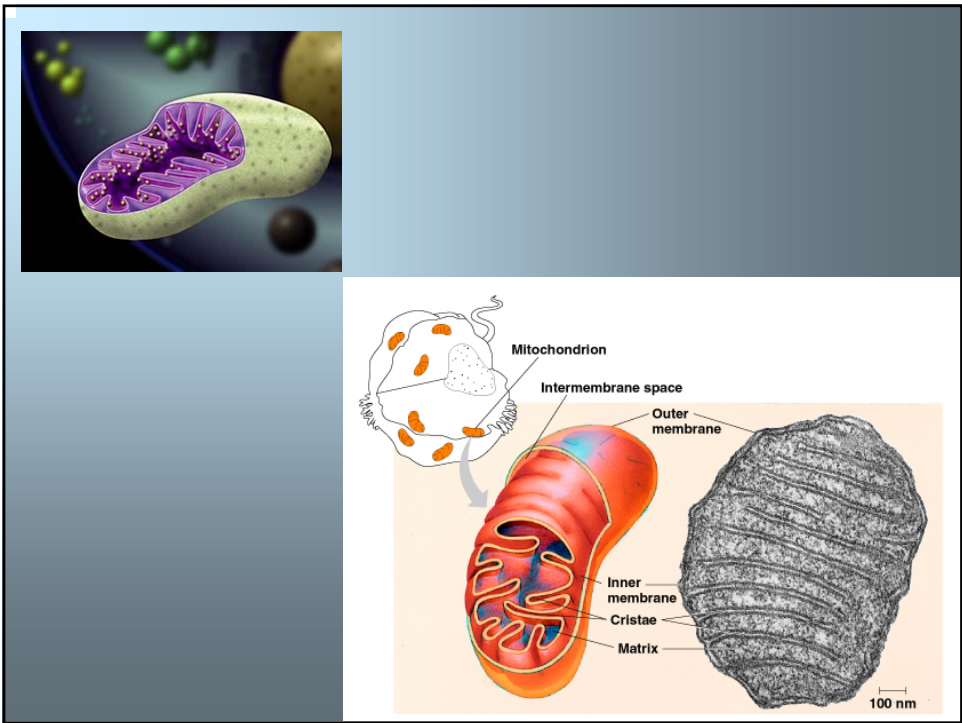
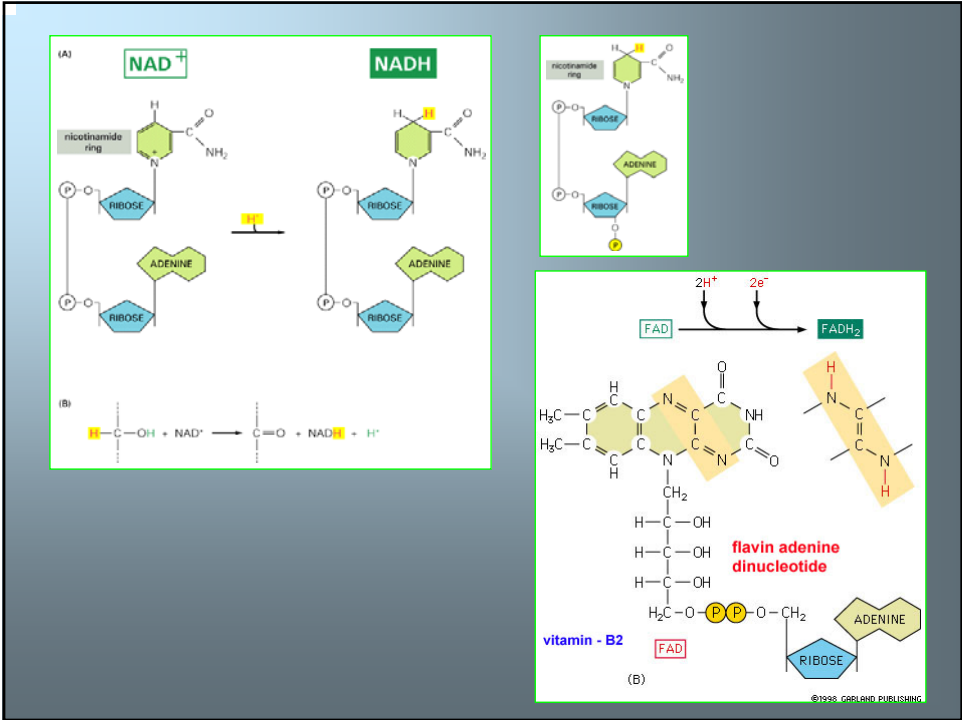


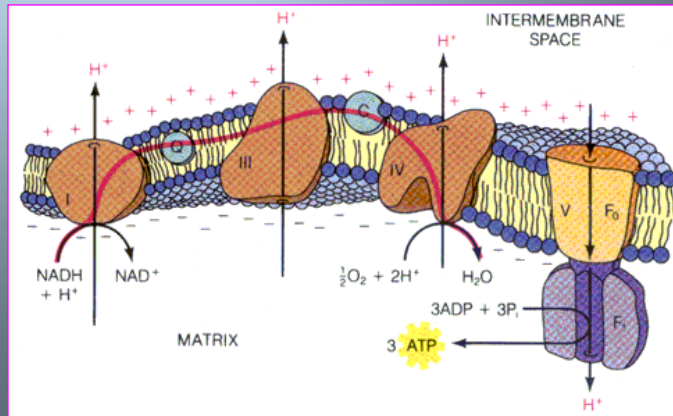
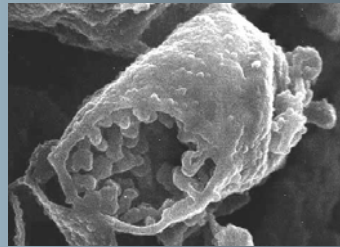
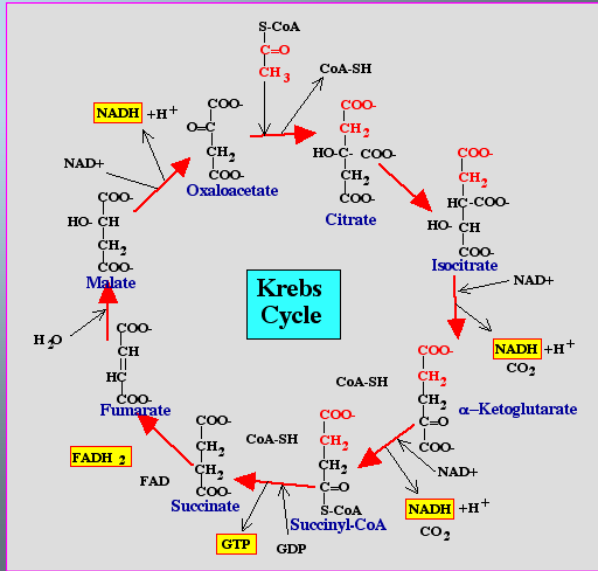
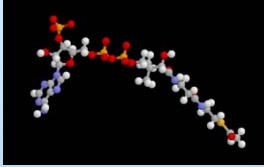


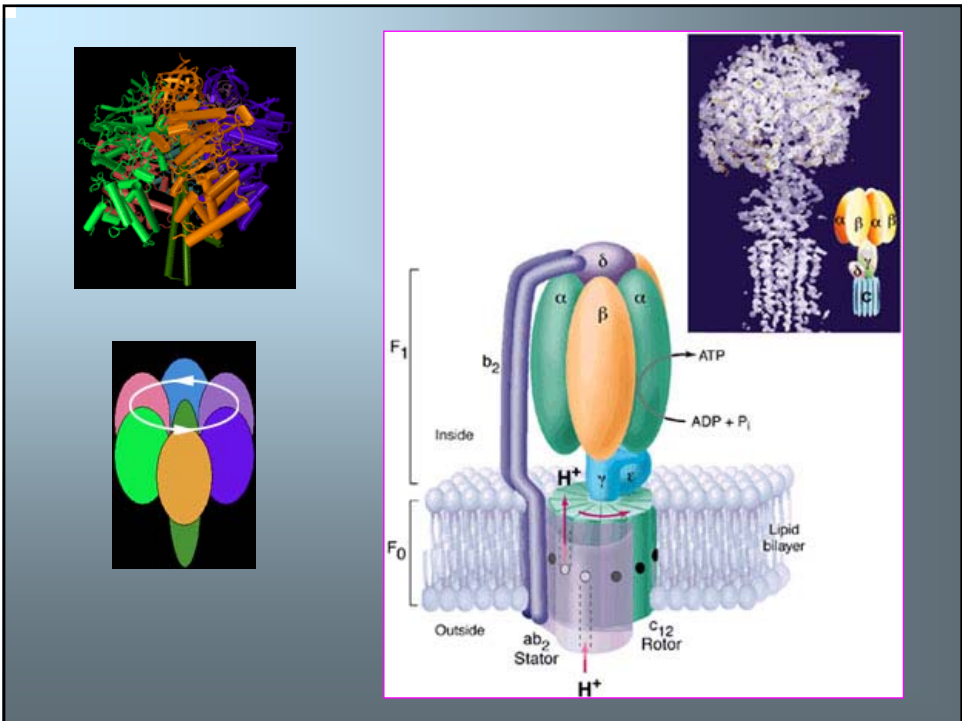
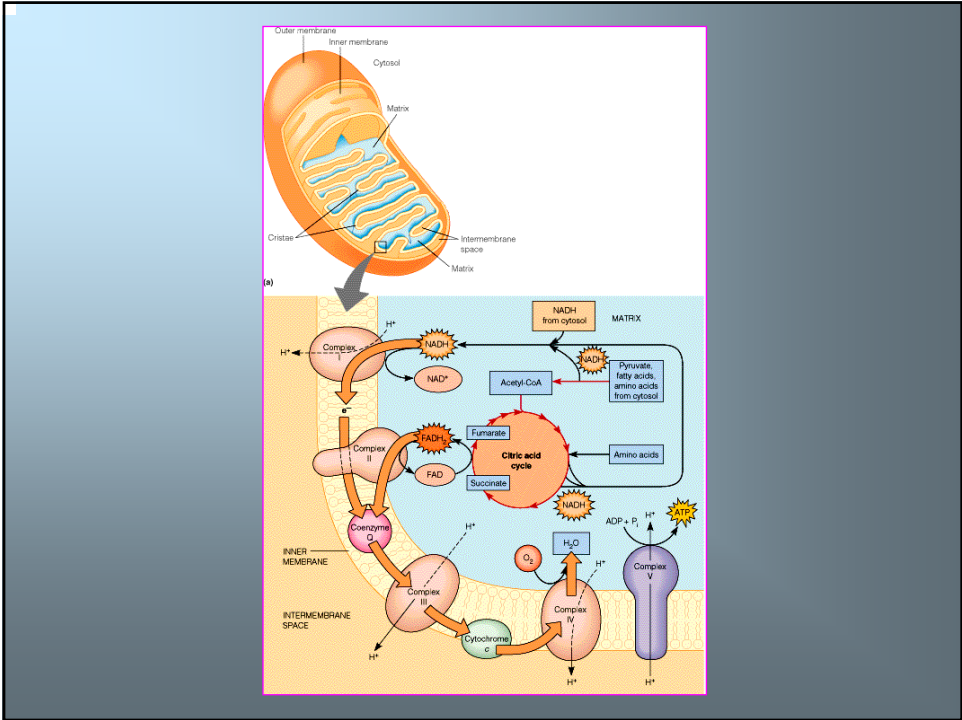


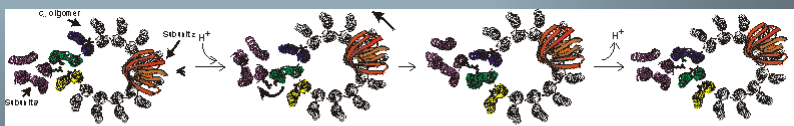
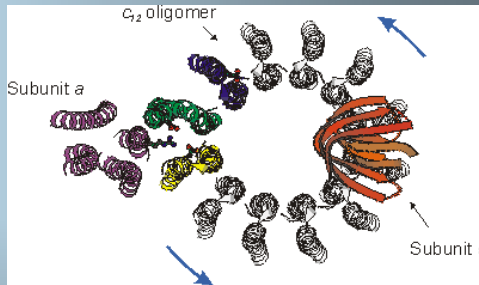
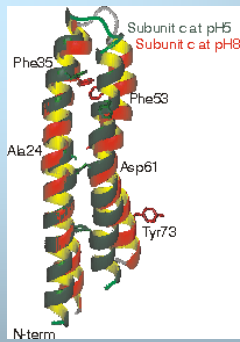












Model for the functioning F_0 complex. (a) Starting state. Arg210 of subunit a lies between protonated and deprotonated Asp61 side chains in the c_2 oligomer. The location of subunit a is shown, positioned according to cross linking data, with its initial orientation indicated by an arrow at right. (b) After protonation of Asp61, the C-terminal helix of the newly protonated monomer (shown in green) rotates towards its stable protonated orientation. (c) Fully protonated intermediate. Subunit a is now at the interface to the next (blue) subunit c . The c_2 ring has rotated by 30° with respect to subunit a . (d) The Asp61 of the next c subunit loses its proton to the F side of the membrane, via a pathway involving both subunit a and subunit c . Its C-terminal helix rotates to adopt the stable conformation of the deprotonated state, regenerating the resting state of the enzyme. The initial position of subunit e is indicated by an arrow at right, to highlight the rotation within the complex.

