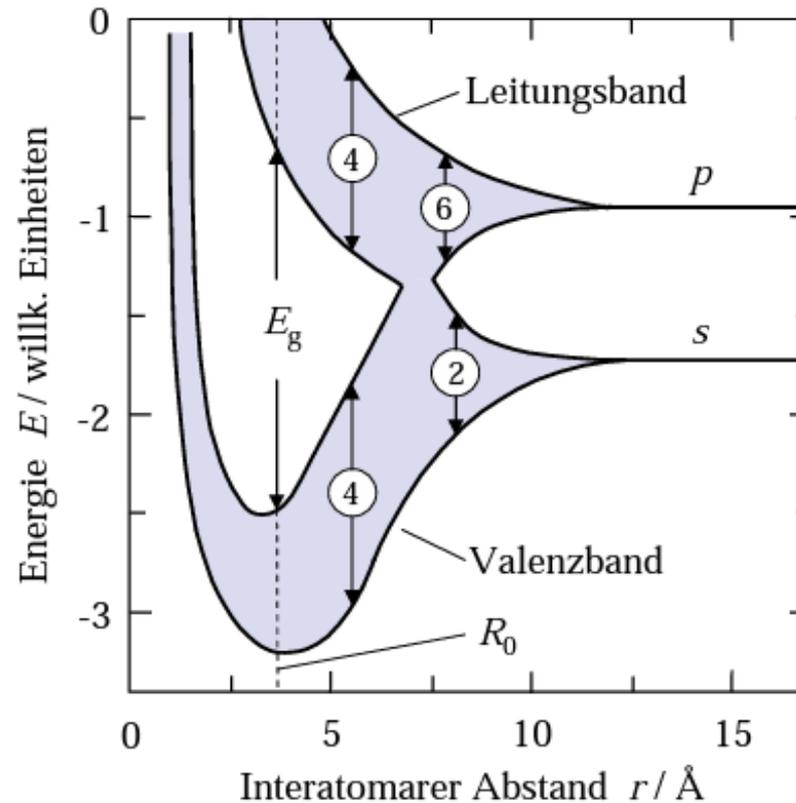


- 1. Elektronen in einem schwachen periodischen Potential**
- 2. Bloch-Theorem; Bloch-Funktionen**
- 3. Tight-binding-Model: Stark-gebundene Elektronen**
- 4. Energiedispersionskurven  $E(k)$**
- 5. Brillouin-Zonen und Fermi-Flächen**

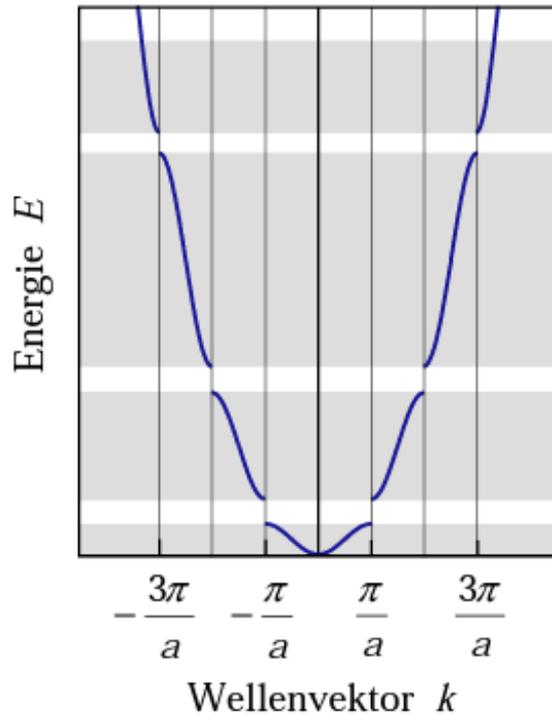
# LCAO Rechnung für Diamant

Abstandsabhängigkeit der Energiebänder in Diamant:  $R_0 = 3,57 \text{ \AA}$  ;  $E_g = 5,5 \text{ eV}$

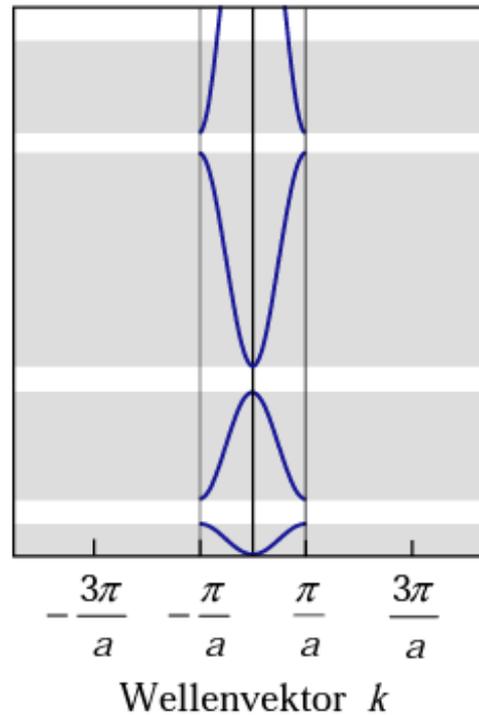


© A.H. Wilson, *The Theory of Metals* (1965)

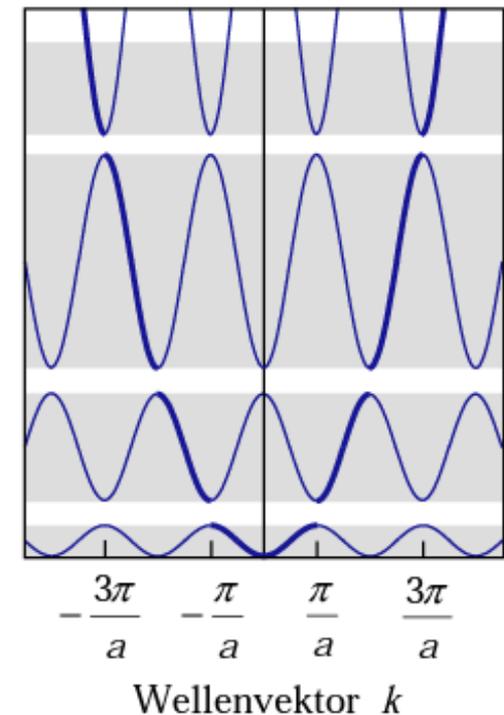
# Energiedispersionskurven



erweitertes  
Zonenschema



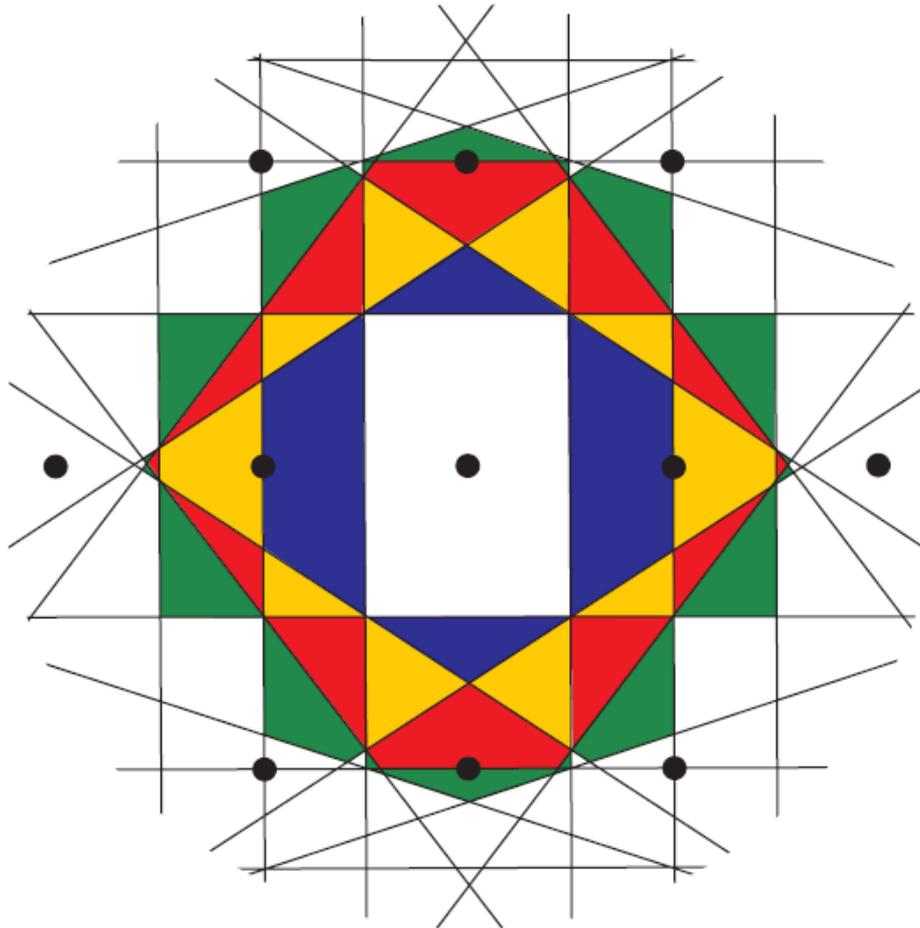
reduziertes  
Zonenschema



periodisches  
Zonenschema

© S. Hunklinger

# Brillouin-Zonen des ebenen Rechteck-Gitters



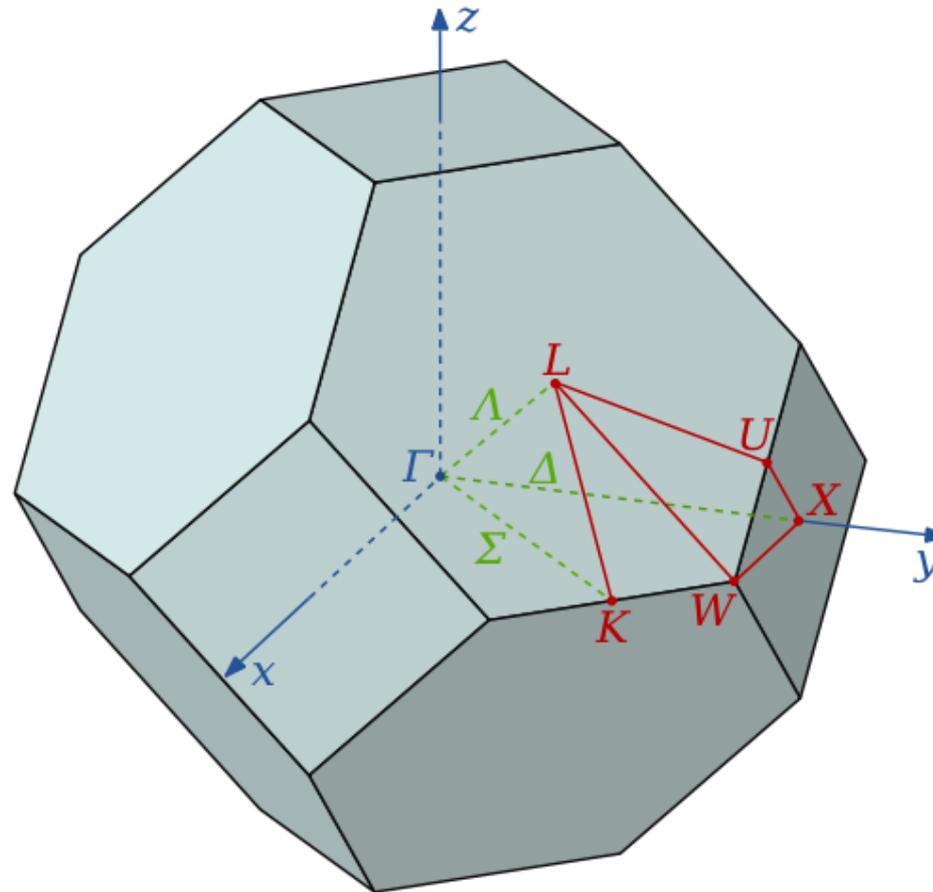
-  1. Brillouin-Zone
-  2. Brillouin-Zone
-  3. Brillouin-Zone
-  4. Brillouin-Zone
-  5. Brillouin-Zone

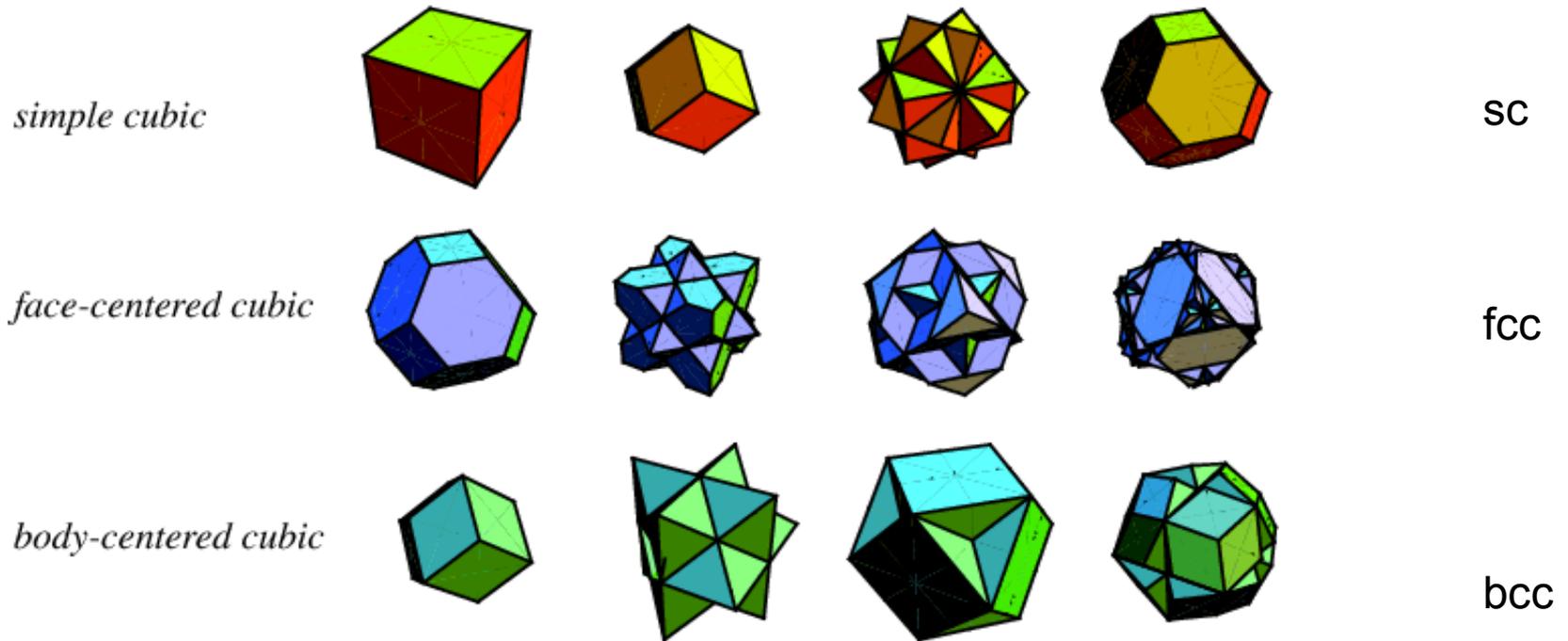
© H.v.Löhneysen

# Erste Brillouin-Zone eines kubisch-flächenzentrierten Gitters

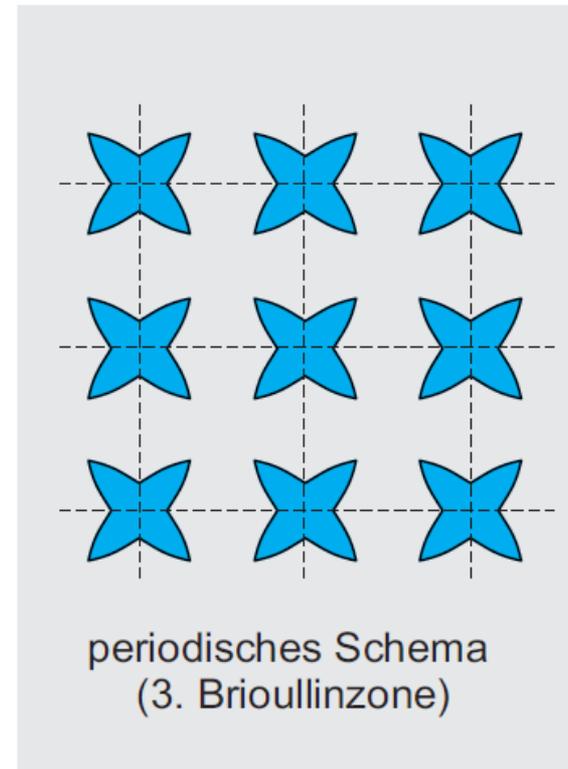
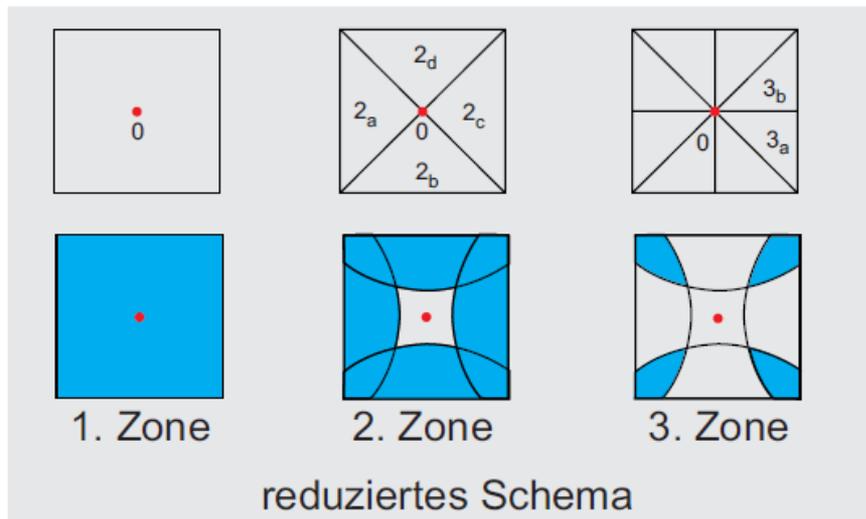
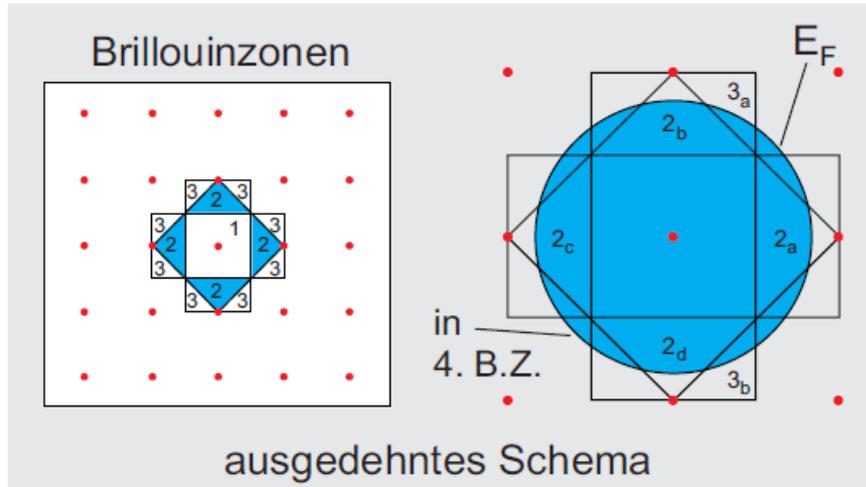
1.BZ

fcc

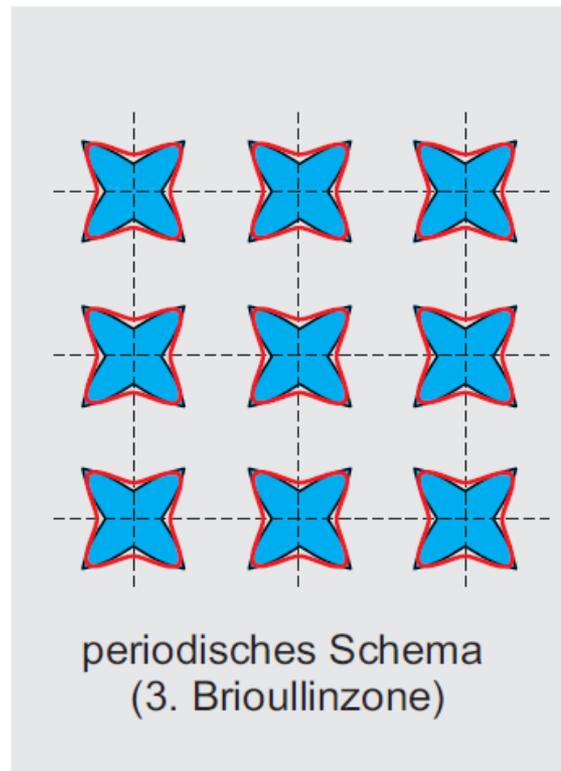
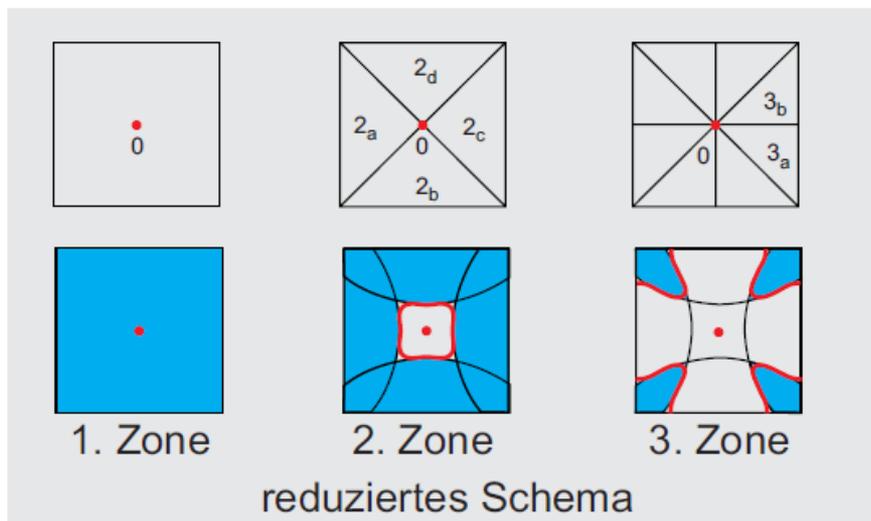
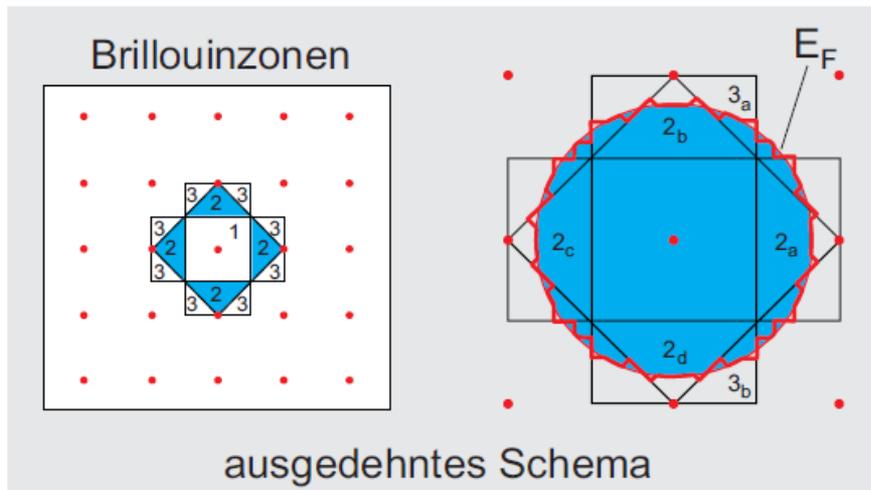




# Fermifläche des freien Elektronengases

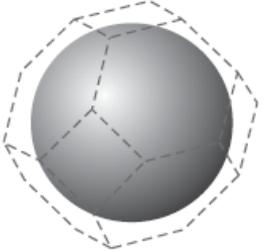
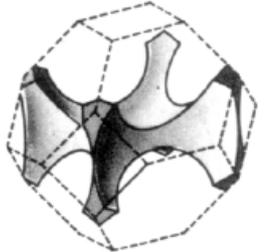
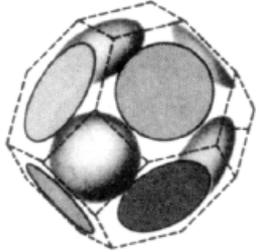
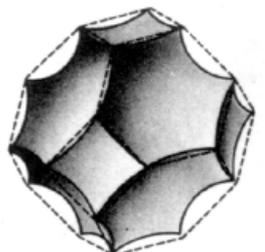
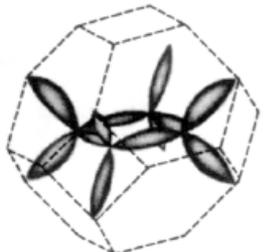
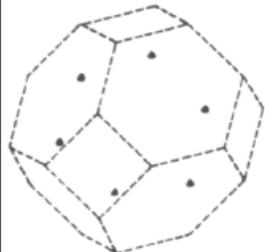


# Fermifläche für Elektronen im periodischen Potenzial



# Fermiflächen von fcc-Metallen für freie Elektronen

Anteile in verschiedenen Brillouin-Zonen

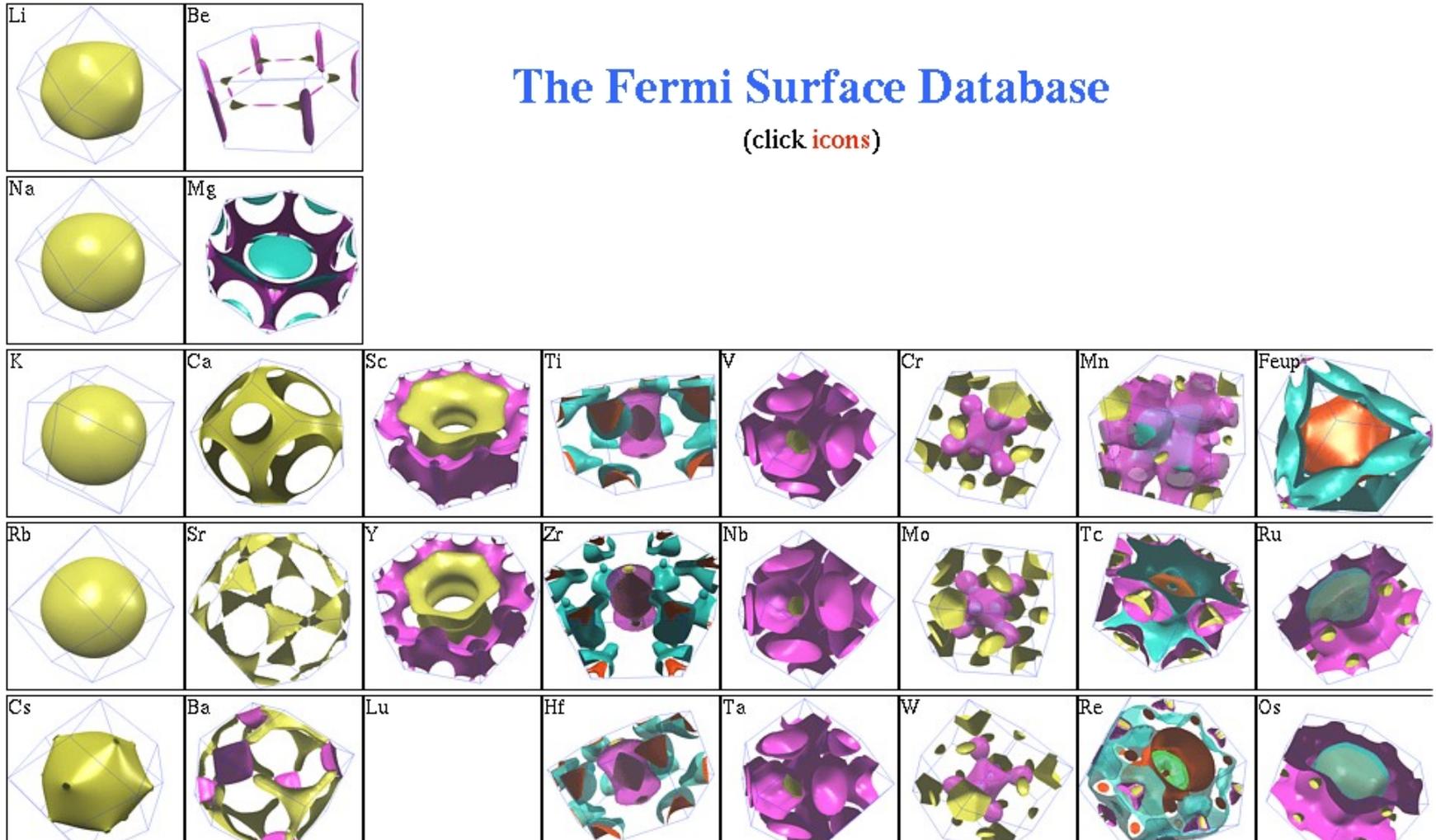
	erste Zone	zweite Zone	dritte Zone	vierte Zone
einwertig		—	—	—
zweiwertig			—	—
dreiwertig	—			

# The Fermi Surface Database

<http://www.phys.ufl.edu/fermisurface/>

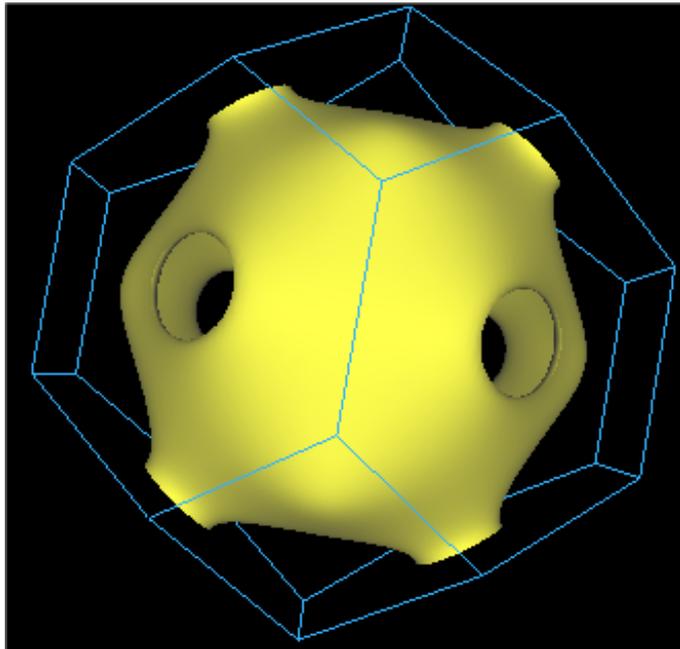
## The Fermi Surface Database

(click icons)

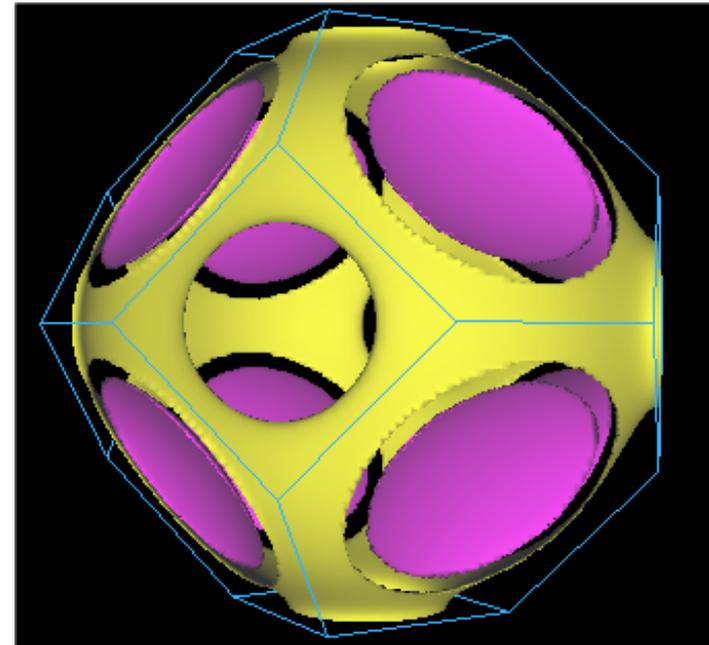


z.B.

<http://www.phys.ufl.edu/fermisurface/>



Cu



Zn