

# Kristall = Gitter x Basis

Basis = periodisch  
wiederkehrende Struktureinheit



Fig-FK- 2.1

# Kristall = Gitter x Basis

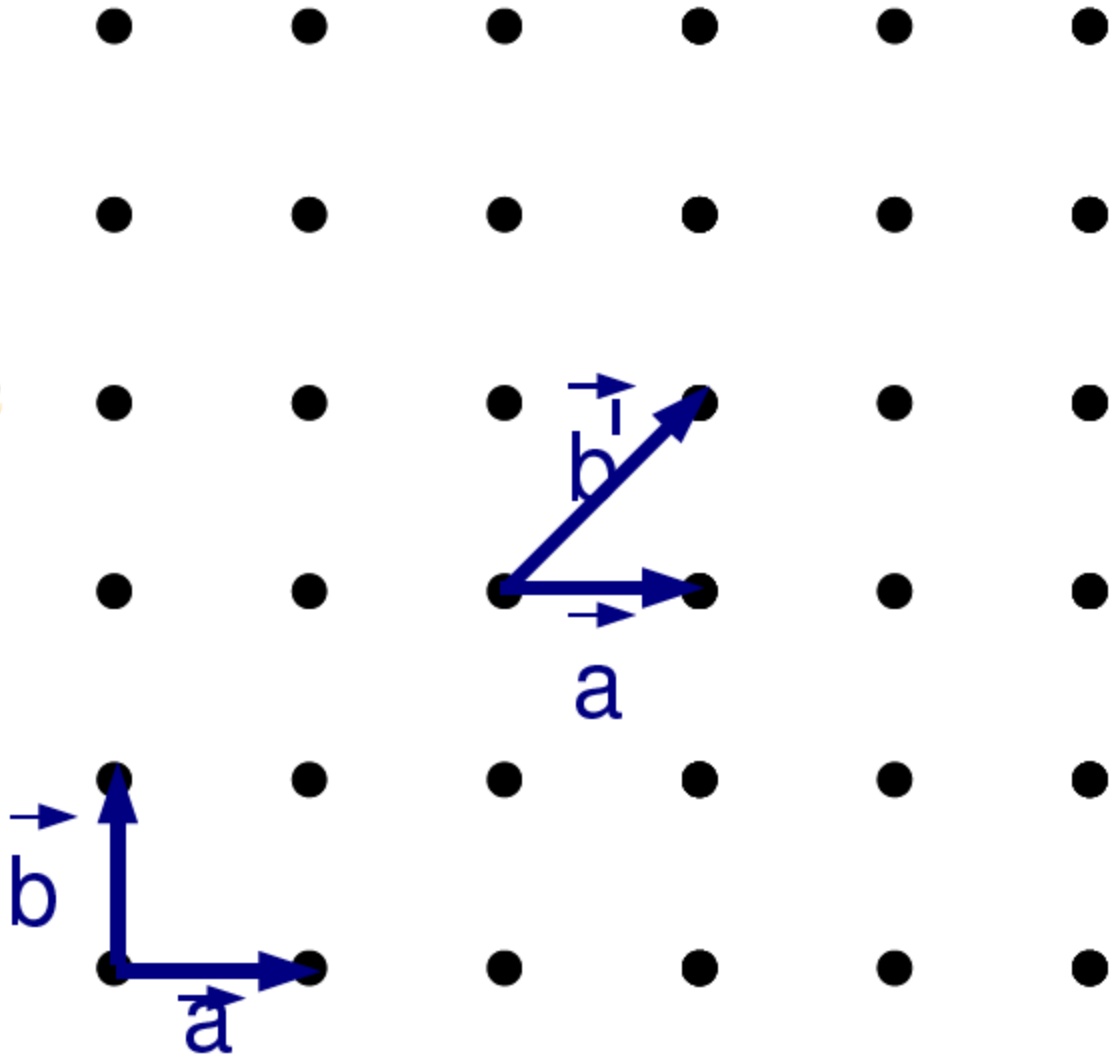
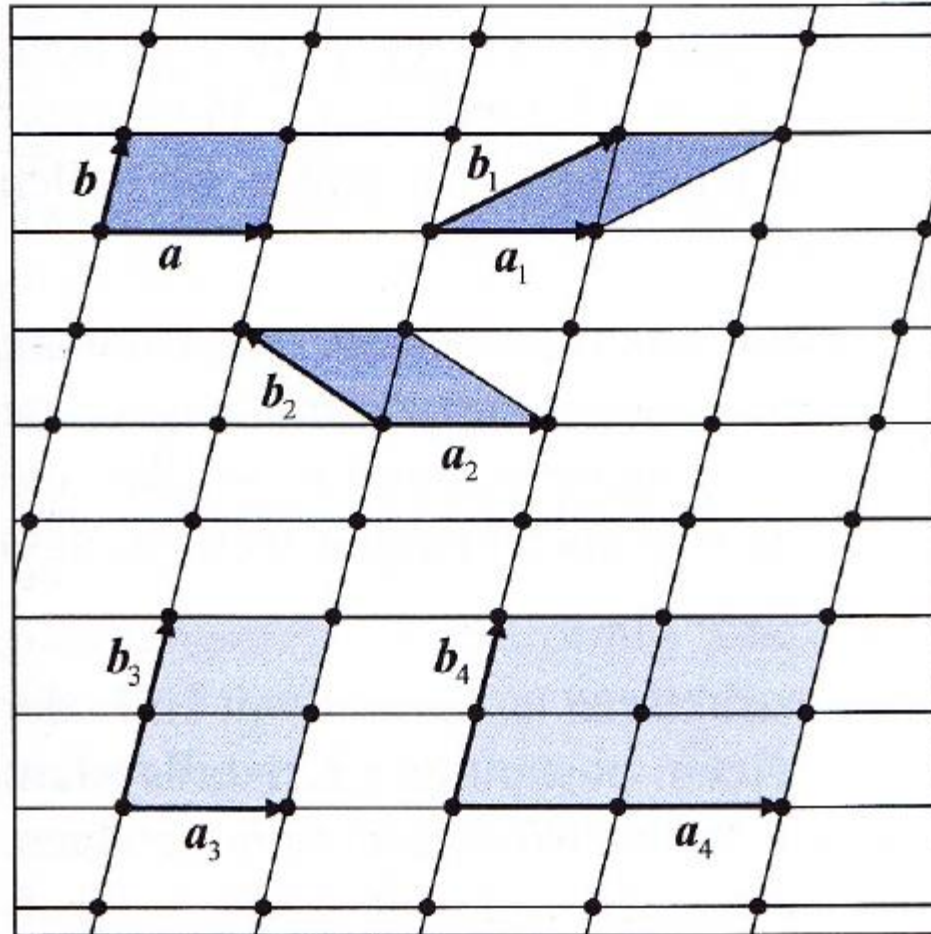


Fig-FK- 2.2

# 2 dimensionales Punktgitter: Verschiedene Elementarzellen

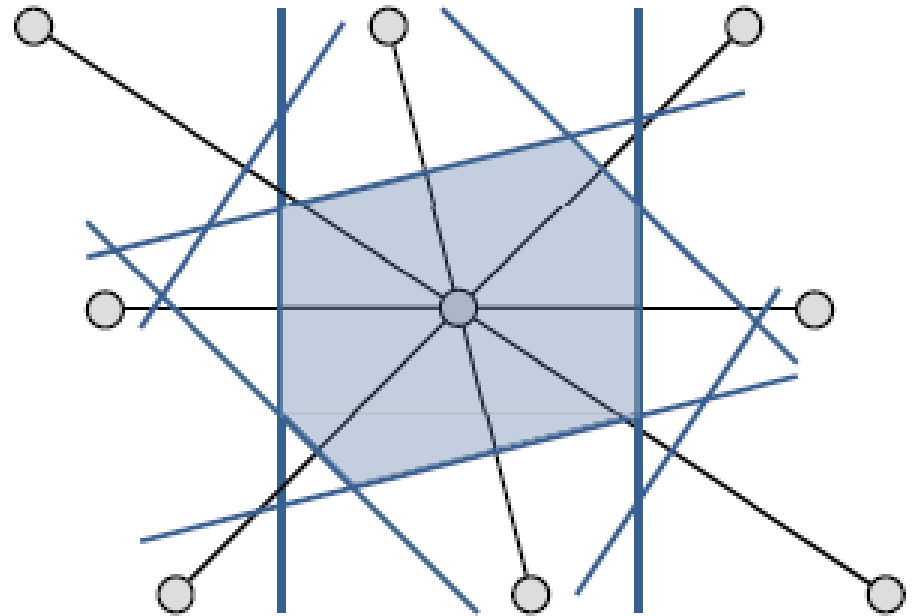
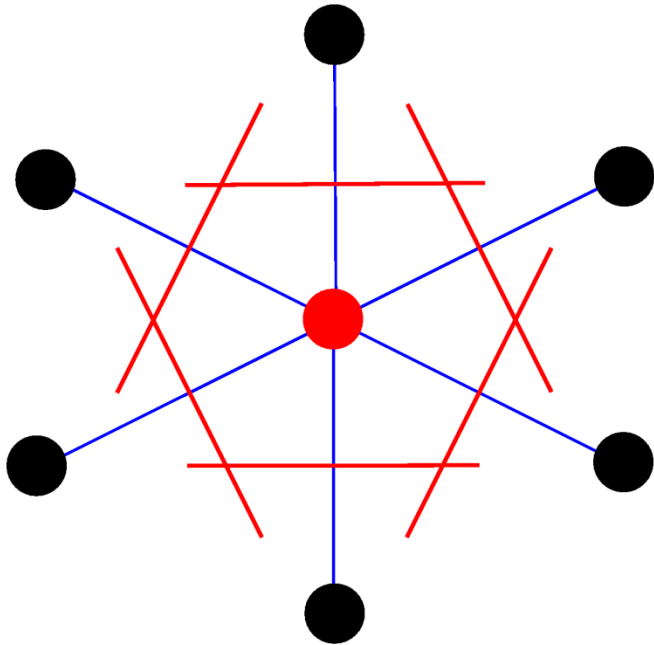


primitiv

Fig-FK- 2.3

# Wigner-Seitz Zelle

Fig-FK- 2.4

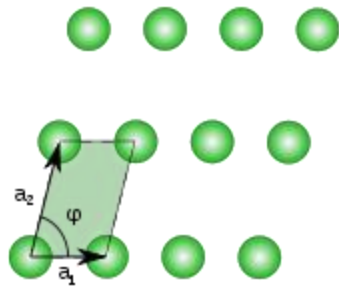


Im reziproken Gitter bezeichnet man die WSZ als 1. Brillouin-Zone.

# 2 dimensionale Bravais-Gitter:

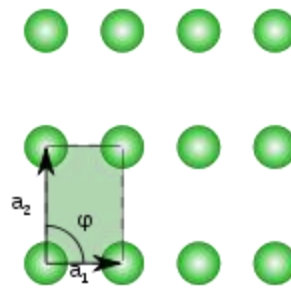
Allg. schiefwinkelige Zelle

+ 4 Spezialfälle



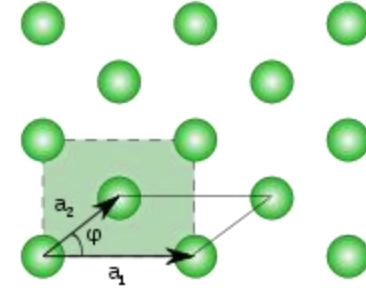
$$|a_1| \neq |a_2|, \varphi \neq 90^\circ$$

1



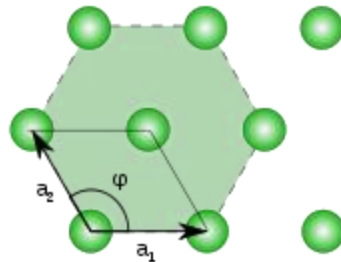
$$|a_1| \neq |a_2|, \varphi = 90^\circ$$

2



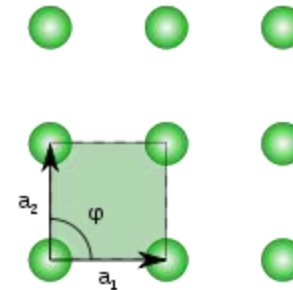
$$|a_1| \neq |a_2|, \varphi \neq 90^\circ$$

3



$$|a_1| = |a_2|, \varphi = 120^\circ$$

4



$$|a_1| = |a_2|, \varphi = 90^\circ$$

5

Fig-FK- 2.5

# 3-dimensionale Bravais-Gitter

## 7 Kristallsysteme

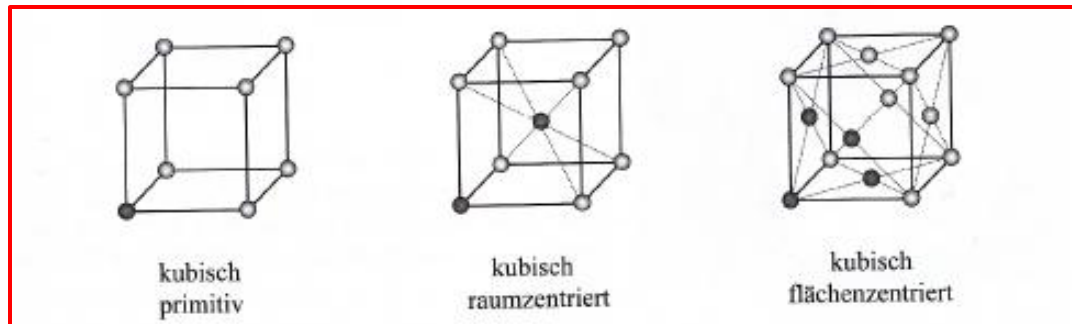
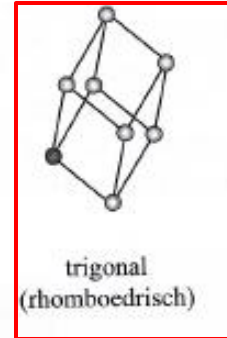
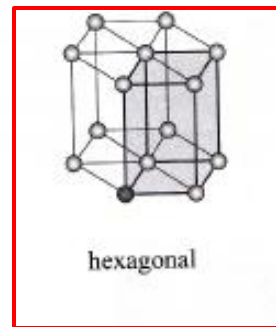
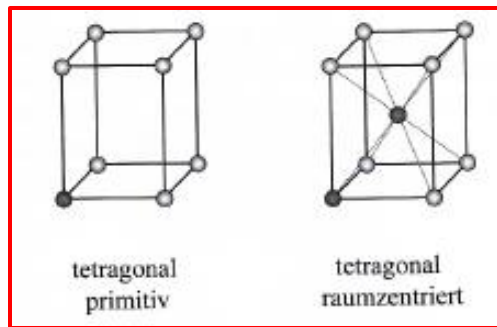
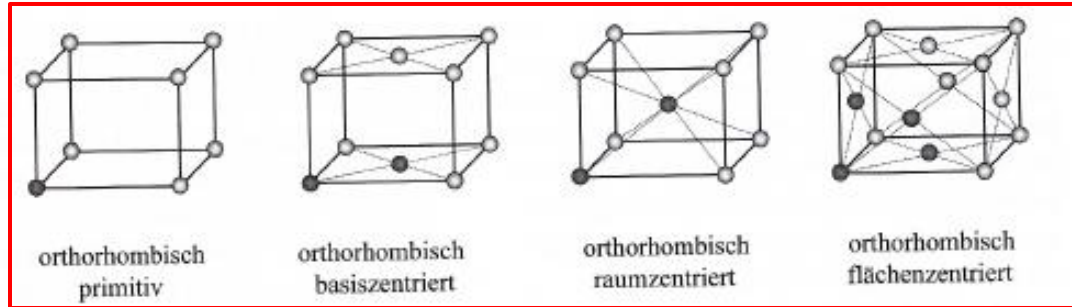
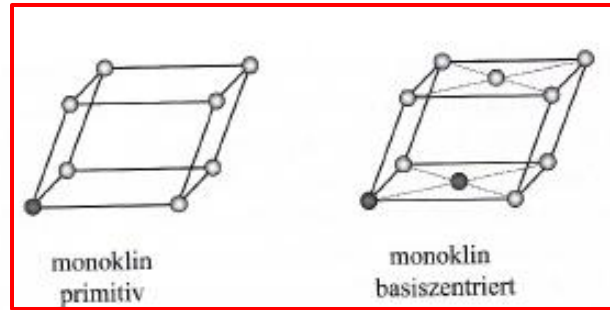
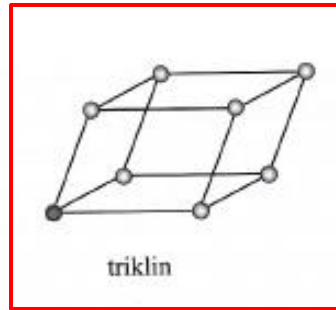
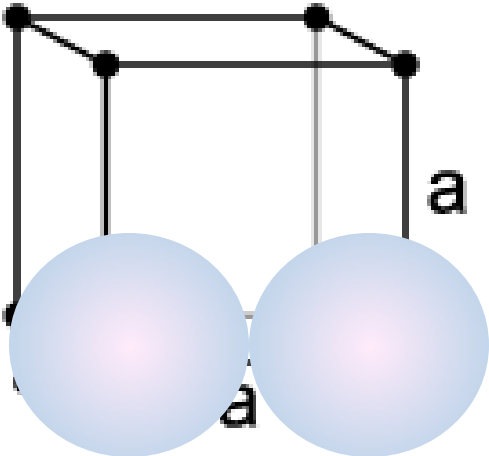


Fig-FK- 2.6

# Kubische Kristallgitter

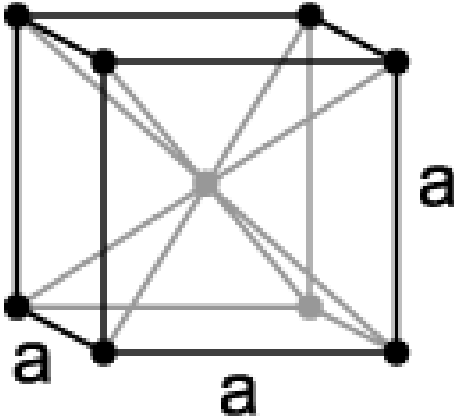
sc



$$a=2R$$

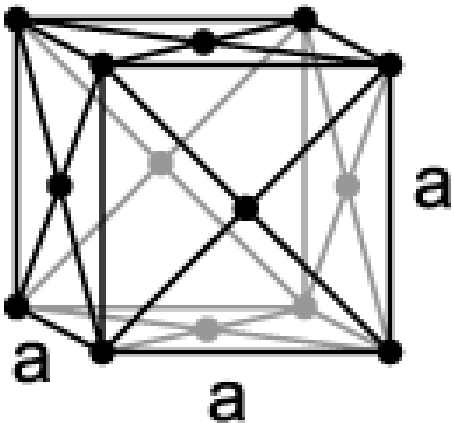
Raumerfüllung:  
52%

bcc



Raumerfüllung:  
68%

fcc



Raumerfüllung:  
74%

Fig-FK- 2.7

## Dichteste Packungen:

fcc: ABCABC (111) Ebene

hcp: ABABAB

Hexagonal Dichte  
Packung (hcp)

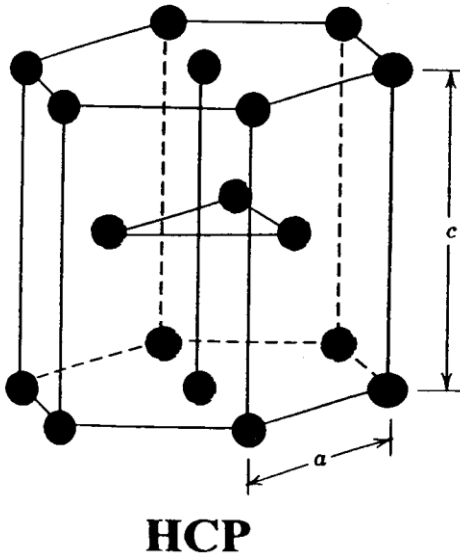
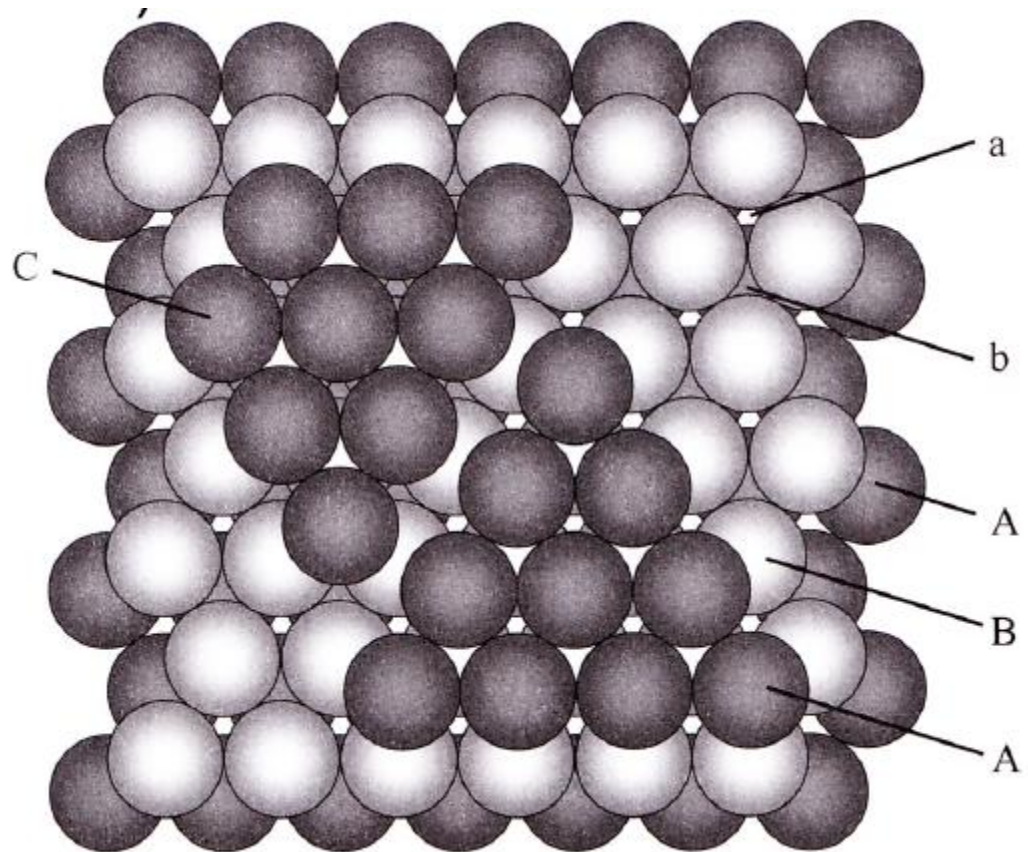


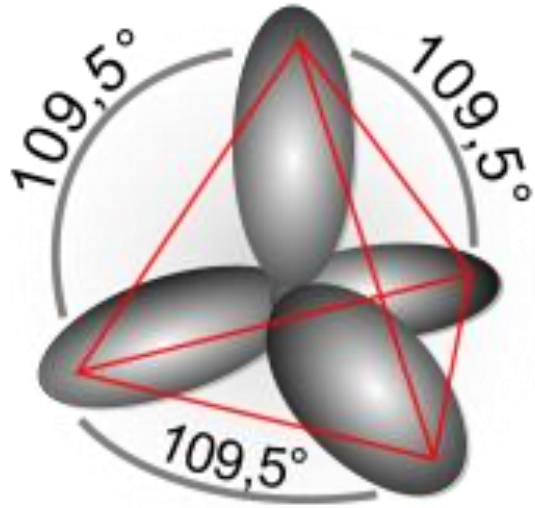
Fig-FK- 2.8



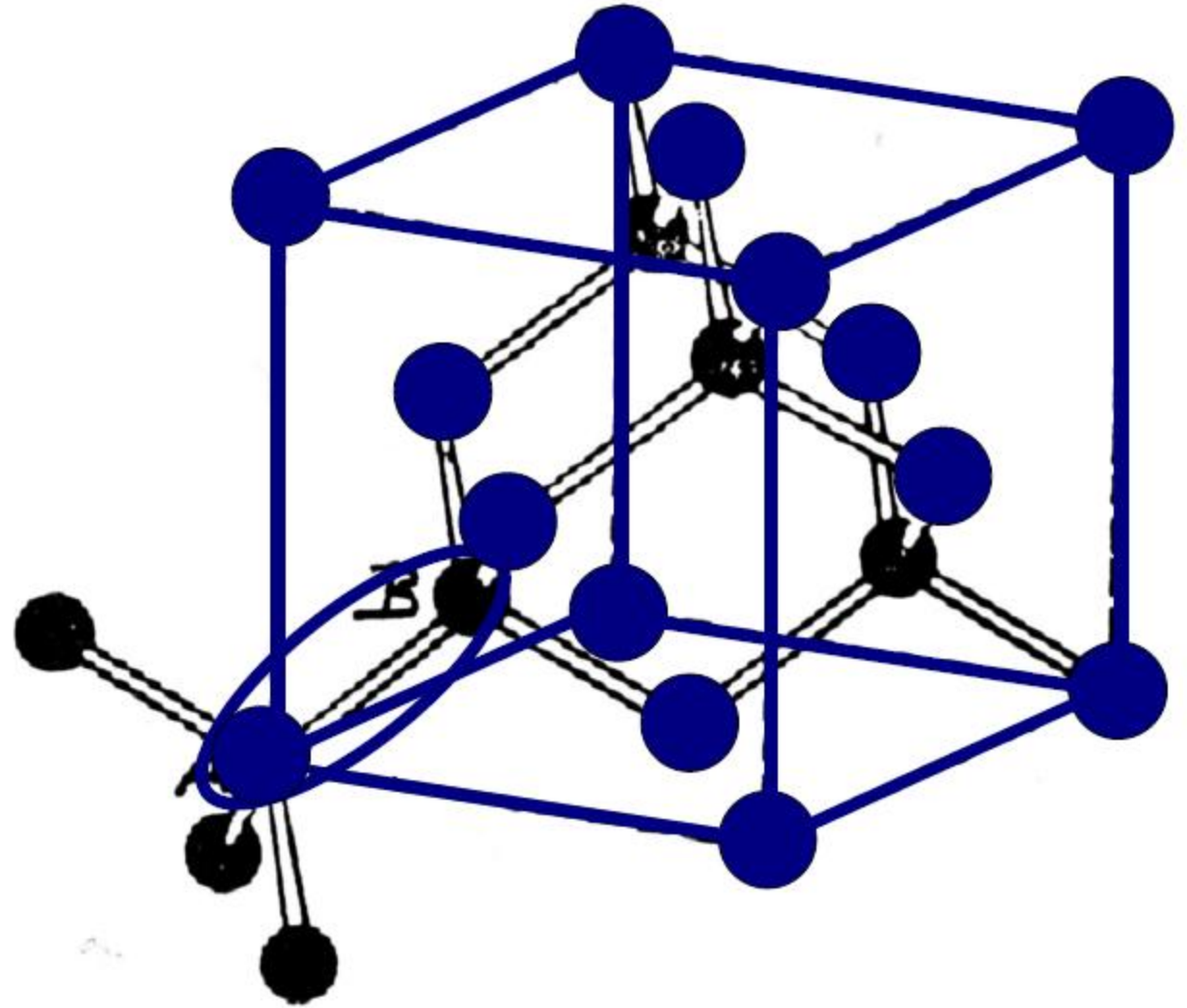
Raumerfüllung: 74%



# Diamantstruktur



$sp^3$   
orbitale



fcc Struktur mit 2-atomiger Basis:  
 $(000) + (\frac{1}{4}, \frac{1}{4}, \frac{1}{4})$

Fig-FK- 2.9

# Kristallgitter

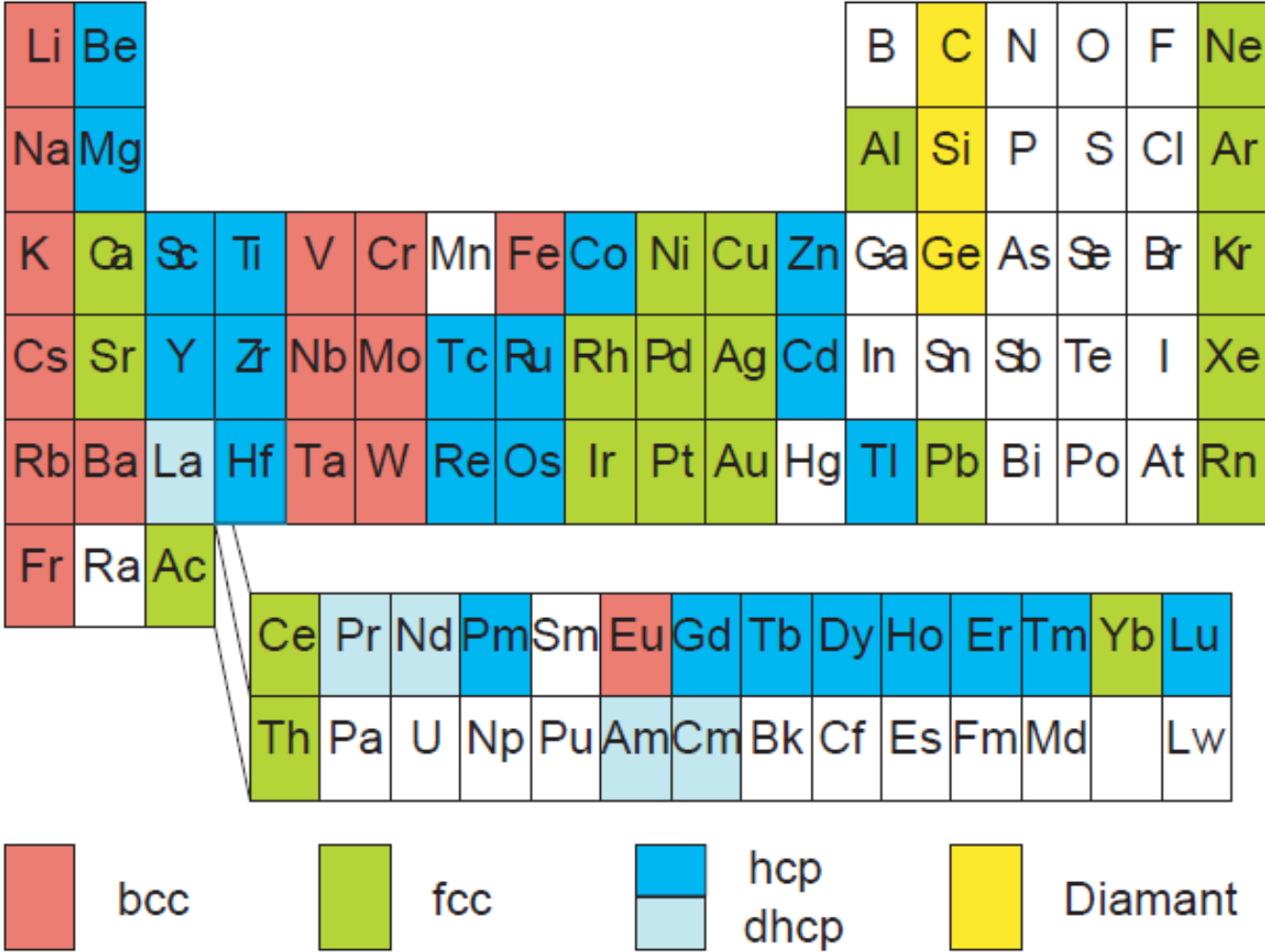
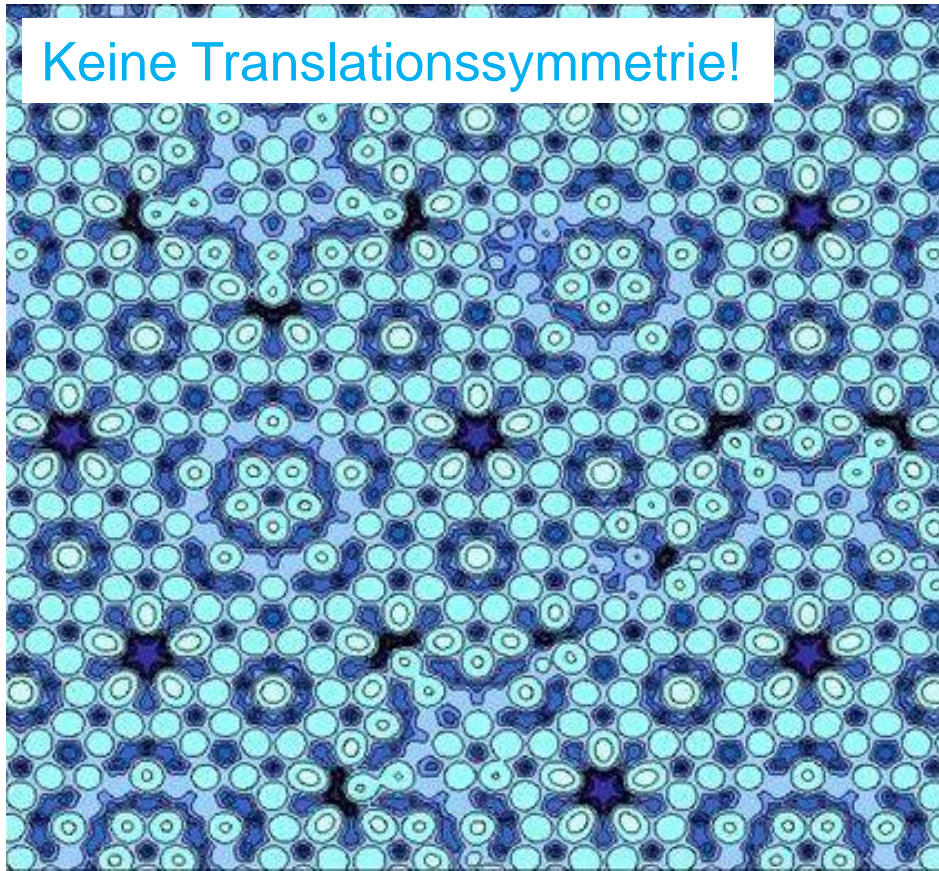


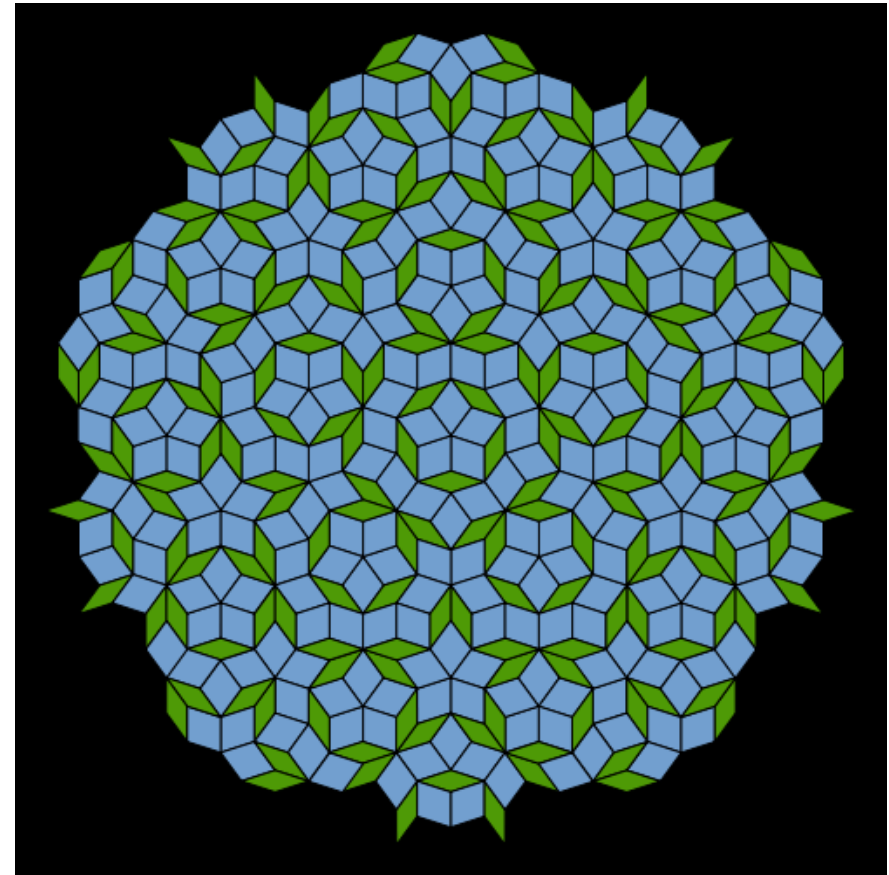
Fig-FK- 2.10

# Quasi-Kristalle = quasi periodische Strukturen



Meist ternäre Legierungssysteme:  
Bsp: Aluminium, Zink, Cadmium

Entdeckt in Al+Mn(14%) Legierung (1984)



Penrose Muster

*Nobelpreis 2011 Chemie  
für Dan Shechtman*