

substance: boron compounds with group VIII elements
property: properties of ternary borides with group VIII elements

(superconducting compounds)

New structure variants of the CeCo_3B_2 -type formed by the ternary alkali metal platinum borides LiPt_3B , $\text{NaPt}_3\text{B}_{1+x}$ and $\text{Na}_3\text{Pt}_9\text{B}_5$ [87M].

monoborides

LnRh_3B (Ln= Sm, Gd, Er, Yb)

Perovskite-type structure

Preparation by the flux method and structure [84T].

diborides

Compound	T_c [K]		
LaRh_3B_2	2.82	structure type CeCo_3B_2	80K2, 80K1
LaIr_3B_2	1.65	structure type CeCo_3B_2	80K2 91F
ThIr_3B_2	2.09	structure type CeCo_3B_2	80K2
CeRu_3B_2	1.10	structure type CeCo_3B_2	85A 80K1
$\text{LaRu}_{2.7}\text{B}_2$	4.0	structure type YOs_3B_2	80K1 91F
YOs_3B_2	6.00	structure type YOs_3B_2	80K1 85A
ThOs_3B_2	3.46	structure type YOs_3B_2	80K1 85A
CeOs_3B_2	3.34	structure type YOs_3B_2 (?)	85A
YRuB_2	7.8	structure type LuRuB_2	80K3 91F
LuRuB_2	9.99	structure type LuRuB_2	80K3
ScOsB_2	1.34	structure type LuRuB_2	80K3
YOsB_2	2.22	structure type LuRuB_2	80K3
LuRuB_2	2.66	structure type LuRuB_2	80K3

tetraborides

Compound	T_c [K]		
YRh ₄ B ₄	11.34	structure type CeCo ₄ B ₄	77M
NdRh ₄ B ₄	5.4	structure type CeCo ₄ B ₄	77M 91F
SmRh ₄ B ₄	2.72	structure type CeCo ₄ B ₄	77M
ErRh ₄ B ₄	8.7	structure type CeCo ₄ B ₄	77M 91F
TmRh ₄ B ₄	9.5	structure type CeCo ₄ B ₄	77M 91F
LuRh ₄ B ₄	11.76	structure type CeCo ₄ B ₄	77M 91F
ThRh ₄ B ₄	4.34	structure type CeCo ₄ B ₄	77M 91F
HoIr ₄ B ₄	2.12		80K1 91F
ErIr ₄ B ₄	2.34		80K1
TmIr ₄ B ₄	1.75		80K1

CeT₃B₂ (T=Co, Ru, Rh, Ir)

Intermediate valence behaviour compared with UT₃B₂ [84Y].

CeCo₃B₂

Preparation and structure in [69K].

Superconducting and magnetic properties of borides with the CeCo₃B₂ structure (MRu₃B₂, M = La, Ce, Pr, Nd, Sm, Gd, Tb, Dy, Ho, Er, Tm, Yb, Lu, Y, Th, U); MRh₃B₂ (M = La, Ce, Pr, Nd, Sm, Eu, Gd); MOs₃B₂ (M = Lu, U), MIr₃B₂ (M = La, Th, U) [80K1, 80K2, 80K3].

RRh₃B₂ (R = La to Gd)

Magnetic behaviour of RRh₃B₂ (R = La to Gd) ternary borides [83M, 83V].

CeRh₃B₂

Preparation of single crystals by Czochralski method [87K2, 87H, 86H].

Structure: hexagonal

Space group: P6/mmm

lattice parameters

a	5.057(3) Å	$T = 300$ K	Weissenberg and precession method	87H
c	3.036(2) Å			

Atomic coordinates, temperature factors and interatomic distances in [87H].

Simple cubic anti-perovskite-type structure with a significant number of vacancies at the rhodium position [86H].

occupancies

Ce (1(a) site)	100 %	$T = 300$ K		86H
Rh (3(c) site)	95.4(5) %			
B (1(b) site)	90(6) %			

PrRh_{4.8}B₂

Structure: orthorhombic

Space group: Immm

lattice parameters

(in Å)

<i>a</i>	9.697(4)	$T = 300$ K	X-ray diffraction	88H
<i>b</i>	5.577(2)			
<i>c</i>	25.64(3)			

Atomic positions and interatomic distances in [88H].

occupancy of rhodium sites (B sites, Pr1, 2 sites = 100 % each)

Rh(1) (16(j))	100 %	$T = 300$ K	X-ray diffraction	88H
Rh(2) (16(j))	100 %			
Rh(3) (8(i))	100 %			
Rh(4) (8(i))	100 %			
Rh(5) (16(n))	27(4) %			
Rh(6) (16(l))	33(5) %			

DyB₂

Entropy in [86B].

LuB₂

Entropy in [86B].

ErRh₃B₂

Crystal structure: monoclinic

Space group: C2/m

Crystal structure of ErRh₃B₂ in Fig. 1 [91B].

lattice parameters

(*a*, *b*, *c* in Å)

<i>a</i>	5.362	<i>T</i> = 300 K	powder, X-ray diffraction	81K2
<i>b</i>	9.288			
<i>c</i>	3.099			
α	90.9°			
<i>a</i>	5.32(2)		single crystal	91B
<i>c</i>	9.21(8)			
<i>c</i>	3.077(6)			
α	90.6°			
<i>a</i>	5.3561(4)	<i>T</i> = 300 K	ErRh ₃ ¹¹ B ₂ , X-ray powder diffraction	93B
<i>b</i>	9.2819(7)			
<i>c</i>	3.1013(3)			
α	90.898(8)°			
<i>a</i>	5.385(5)	<i>T</i> = 300 K	neutron diffraction	93H
<i>b</i>	9.25(1)			
<i>c</i>	6.194(5)			
α	90.35(7)°			

Magnetic susceptibility in Fig. 2 [91B], for susceptibility results on ¹⁰B and ¹¹B enriched samples see [93B].

On crystal and magnetic structure see [94H].

Review on magnetic and superconducting properties of ErRh₃B₂ and the related rare earth transition metal borides (SmCo₃B₂, GdCo₃B₂, ErCo₃B₂, LaRh₃B₂, La_{0.4}Ce_{0.6}Rh₃B₂, CeRh₃B₂, Ce(Rh_{0.8}Ir_{0.2})₃B₂, CeRh₃(B_{0.5}Si_{0.5})₂, CeRu₃B₂, Ce(Ru_{0.5}Rh_{0.5})₃B₂, PrRh₃B₂, NdRh₃B₂, SmRh₃B₂, EuRh₃B₂, GdRh₃B₂, LaIr₃B₂, ThIr₃B₂, TbRh₃B₂, DyRh₃B₂, HoRh₃B₂, ErRh₃B₂, TmRh₃B₂, NdIr₃B₂, SmIr₃B₂, GdIr₃B₂, TbIr₃B₂, DyIr₃B₂, HoIr₃B₂, ErIr₃B₂, TmIr₃B₂) in [94B].

LuOs₃B₂

Space group: P6/mmm

Comparative critical field study of superconducting ternary borides in [87L].

LuRuB₂

Space group: Pnma

Comparative critical field study of superconducting ternary borides in [87L].

Superconductivity of ternary borides with the LuRuB₂ structure in [80K3].

further data on tetraborides

RM₄B₄ (R = rare earth element; M = Ru, Rh, Os, Ir)

Preparation, structure, magnetic properties in [77M; 22J, 79K, 79R1, 80S, 80O, 80W, 81H].

R(Os, Ir)₄B₄ (R = rare earth element)

NdCo₄B₄ type structure

Preparation, structure, magnetic properties in [79R2].

R(Rh_{1-x}Ir_x)₄B₄ (R = Gd, Tb, Dy, Ho, Er, Tm, Lu)

Low temperature phase diagram of ternary magnetic semiconductors in [84K].

Antiferromagnetism in R(Rh_{1-x}Ir_x)₄B₄ compounds [81K1].

RRh₄B₄ (R = rare earth element)

LnRh₄B₄ (Ln = lanthanides such as Nd, Sm, Er, Tm, Lu)

Chemical bonding topology of the superconducting compounds in [87K4].

Superconductivity and magnetism in [81M].

The superconductivity of YRh₄B₄ – an enigma in [81E].

Ferromagnetism of ReRh₄B₄ compounds [80M1].

Roles of crystal fields in magnetic superconducting rare-earth rhodium borides [80M2].

Superconducting T_C , ferromagnetic T_C and antiferromagnetic T_N transition temperatures vs. R for RRh₄B₄ compounds in [91M].

Magnetic ordering in superconducting rare-earth compounds [80I].

Low temperature ordered states due to dipole-dipole and exchange interactions [84M].

DyRh₄B₄

Temperature dependence of the magnetic susceptibility in [91M]

ErRh₄B₄

Structure: primitive tetragonal

Space group: P4₂/nmc

lattice parameters

a	5.292(4) Å	$T = 300$ K	X-ray diffraction	94S
c	7.379(3) Å			

superconductivity data

T_c	8.7 K			91F, 77M 94S
T_{c1}	8.55 K			
ΔT_{c1}	30 mK		width of transition range	
T_{c2}	0.84 K	cooling run		
T_{c2}	0.90 K	heating run		

density

d	10.21 g cm ⁻³ 10.00 g cm ⁻³	$T = 293$ K	pycnometric X-ray	94S
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Summary of the properties, prospects and perspectives of borides of the type ErRh₄B₄ [82F]

Magnetic properties in [94S].

Ion damage, critical current and tunneling studies of ErRh_4B_4 films in [80R].

LuRh_4B_4

Structure: orthorhombic

lattice parameters

(in nm)

a	0.5898(1)	$T = 300 \text{ K}$	X-ray diffraction	96O
b	1.1420(1)			
c	0.3485(1)			

Comparative critical field study of superconducting ternary borides [87L].

resistivity

ρ	86(6) $\mu\Omega \text{ cm}$	$T = 300 \text{ K}$		96O
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microhardness

H_V	14.0(6) GPa	$T = 300 \text{ K}$	load 100 g	96O
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$\text{Gd}_3\text{Co}_{11}\text{B}_4$

Structure ($\text{R}_{n+1}\text{Co}_{3n+5}\text{B}_{2n}$ type with $\text{R} = \text{rare earth or Y atom}$, $n = 2$) and electrical resistivity in [96K].

Volume of the unit cell in [91P].

C_p of lanthanide tetraborides in [91P].

RRuB_4 ($\text{R} = \text{Gd, Tb, Dy, Ho, Er, Tm, Lu, Y}$)

(YCrB_4 structure)

Low temperature magnetic order [87K16].

HoRu_4B_4

Nature of magnetic order below 4.2 K (two magnetic phase transitions at 2.6 and 2.3K) in [85M].

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Fig. 1.

ErRh_3B_2 . Crystal structure. Top: three-dimensional structure; bottom: (001) plane with the (002) plane projected on it. The monoclinic unit cell and the pseudo-hexagonal unit cell are shown. Note that β is exaggerated and that with $\beta = 90^\circ$ the monoclinic ErIr_3B_2 structure transforms into the hexagonal CeCo_3B_2 structure [91B].

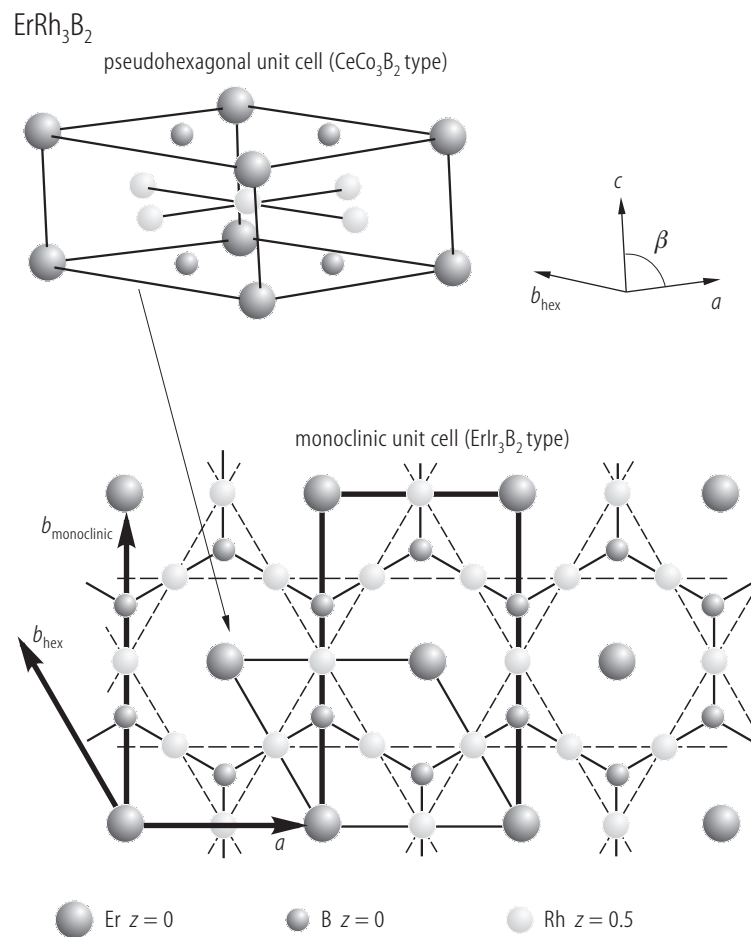


Fig. 2.

ErRh_3B_2 . (a) and (b) Susceptibilities ($\chi = \mu_0 M/B$; in SI units) in low magnetic fields; (c) and (d) inverse magnetic susceptibilities. The b axis in (b) and (d) refers to the monoclinic b axis [91B].

