

substance: boron compounds with group III elements
property: properties of boron-scandium compounds

Sc-Al-B

Crystal growth of scandium aluminium borides and scandium borides by Sc-Al-B ternary system solutions [99O].

Sc₃B_{0.75}C₃

Sc₃B_{0.75}C₃ a novel compound of rare-earth boron carbide [99S].

lattice parameters

<i>a</i>	3.3308(3) Å	<i>T</i> = 300 K	O content 0.13 wt. %	99S
<i>c</i>	7.680(2) Å			

density

<i>d</i>	3.53 g cm ⁻³	<i>T</i> = 300 K		99S
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Physical properties of a layered scandium boron carbide Sc₃B_{0.75}C₃ [99M].

Sc₂B_{1.1}C_{3.2}

Sc₂B_{1.1}C_{3.2} a novel compound of rare-earth boron carbide with graphite-like layer structure [99S].

Structure: trigonal

Space group: P3m1

lattice parameters

<i>a=b</i>	23.710(9) Å	<i>T</i> = 300 K	X-ray diffraction	99S
<i>c</i>	6.703(2) Å			

Sc₂BC₂

Electronic structure of Sc₂BC₂ [90H].

On boron-carbon chains [97B].

ScB₂

metallic conduction (?); preparation [70P, 75S], crystalline structure [70P, 75S], electronic structure [79A], entropy [86B].

ScB₄

preparation [79S]

ScB₁₂

preparation [70P, 75S], crystalline structure [70P, 75S], electronic structure [77P]

LaB₆ – "ScB₆"-system

preparation [79S]

ScRu₄B₄

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

critical temperature of superconductivity

T_c	0.39 K	91F, 68M
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entropy

S	96(13) J mol ⁻¹ K ⁻¹	86B
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On the preparation of ScB₆ in cubic and tetragonal structures [94P].

ScB₁₇C_{0.25}

A single crystal XRD and TEM study of ScB₁₇C_{0.25}: a novel structure with boron-based "nanotubular" molecular units [99L].

Structure: hexagonal

Space group: P6/mmm

lattice parameters

a	1.4476 nm	$T = 300$ K	whisker-like crystals	99L
c	0.89375 nm			

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