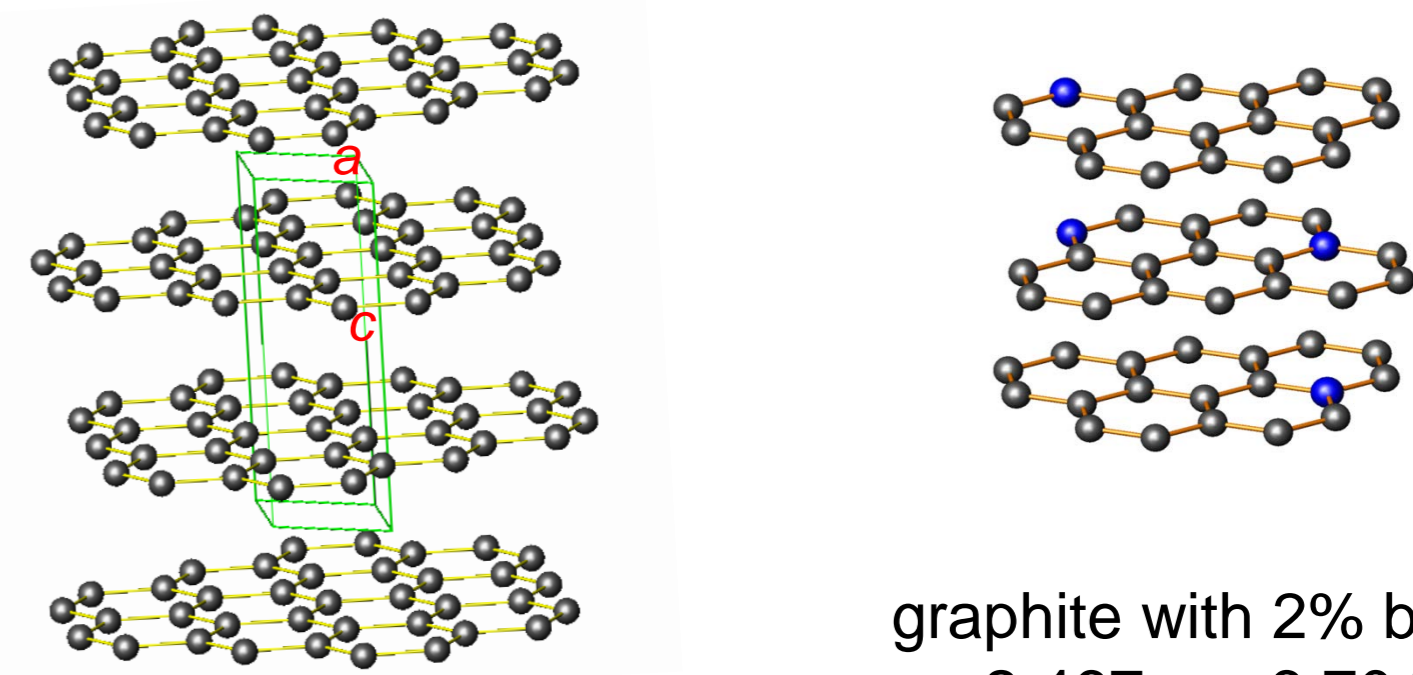
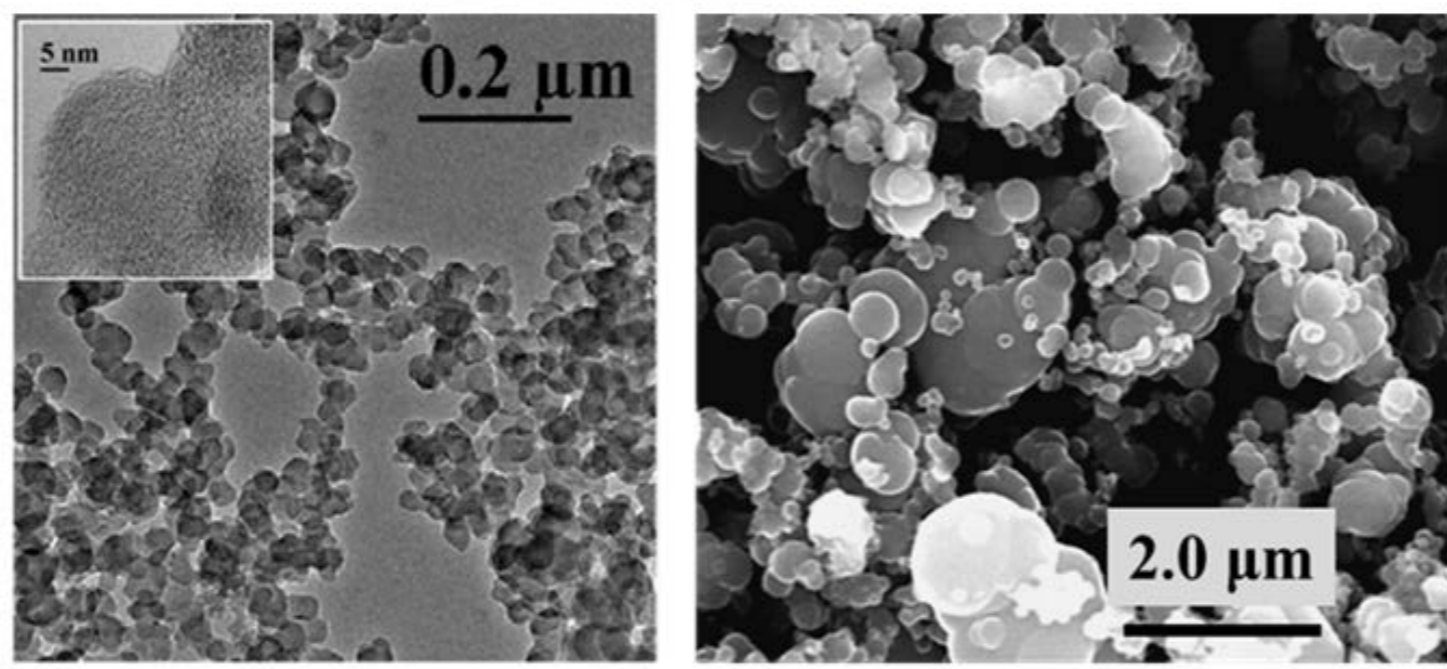




# Effect of boron on the structure of graphite formed at high pressure

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globular nanocarbon + boron

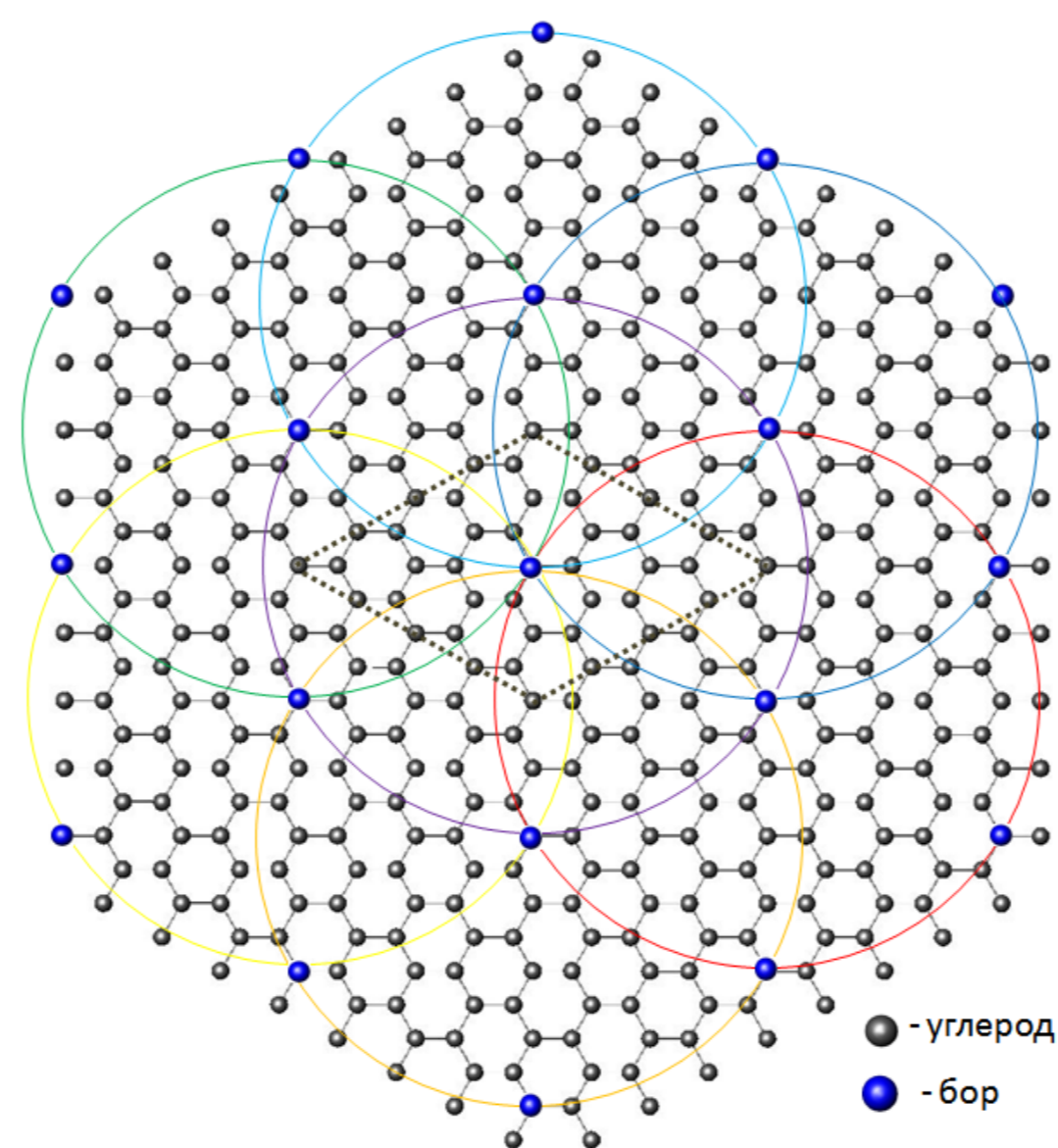
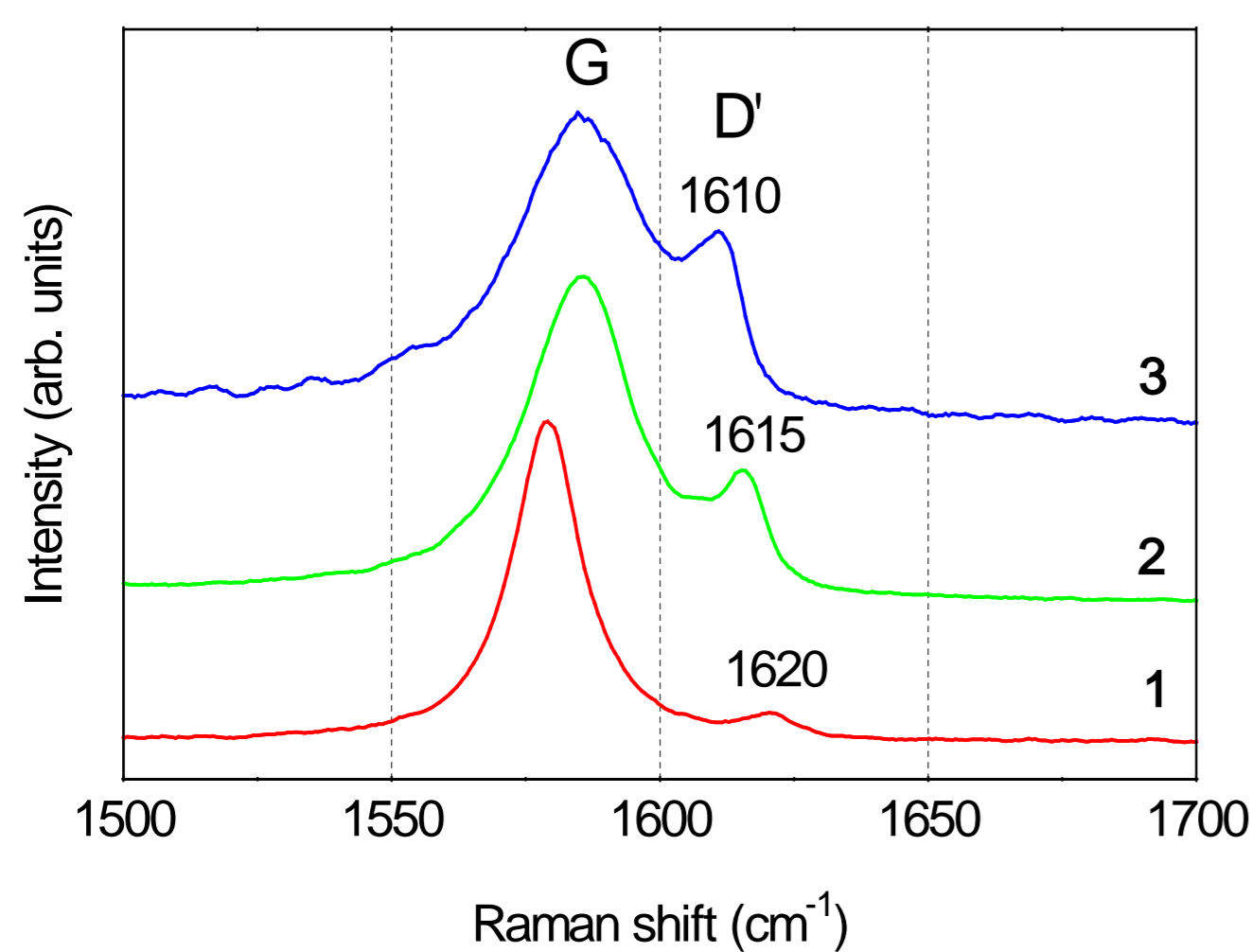
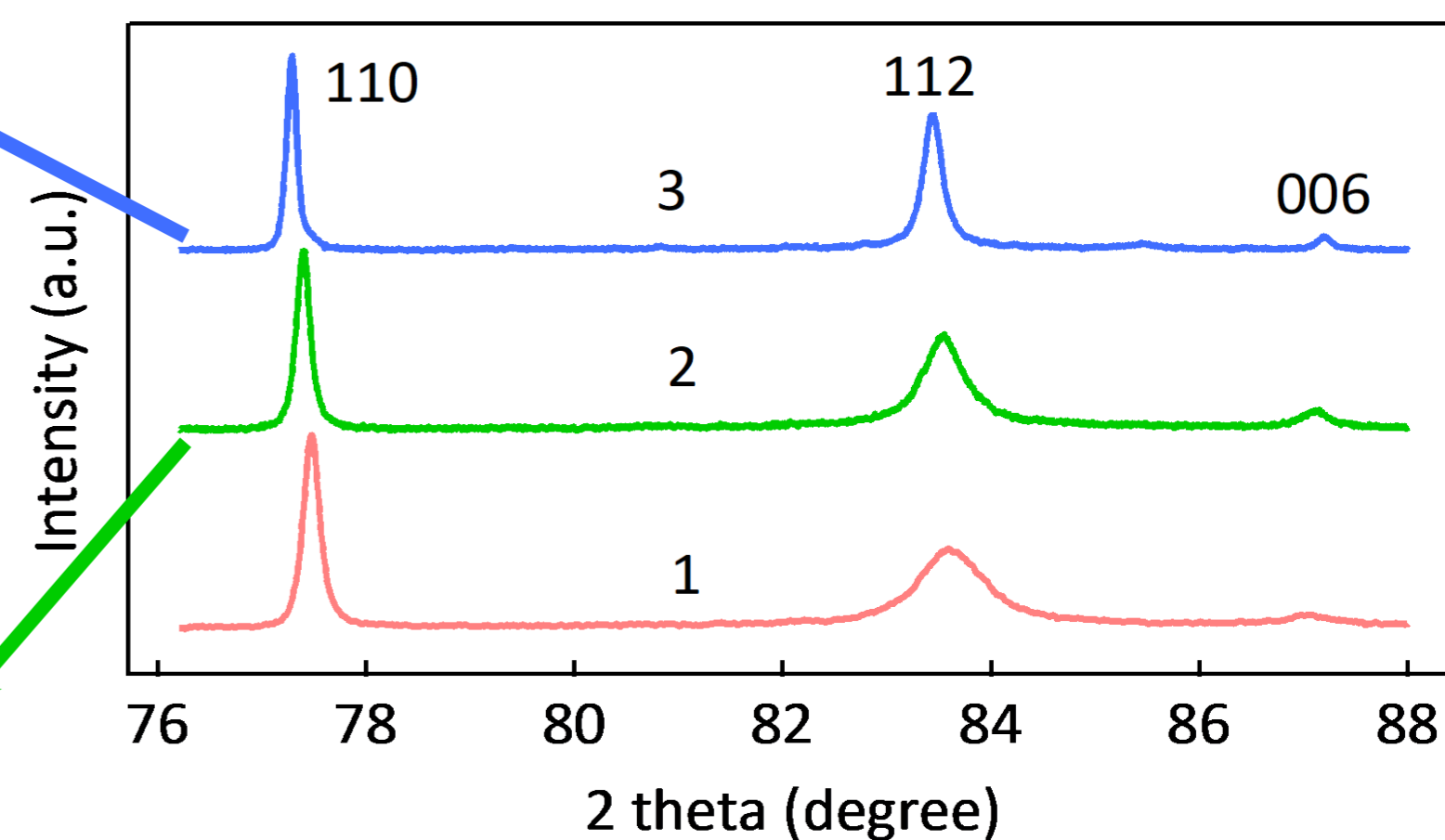
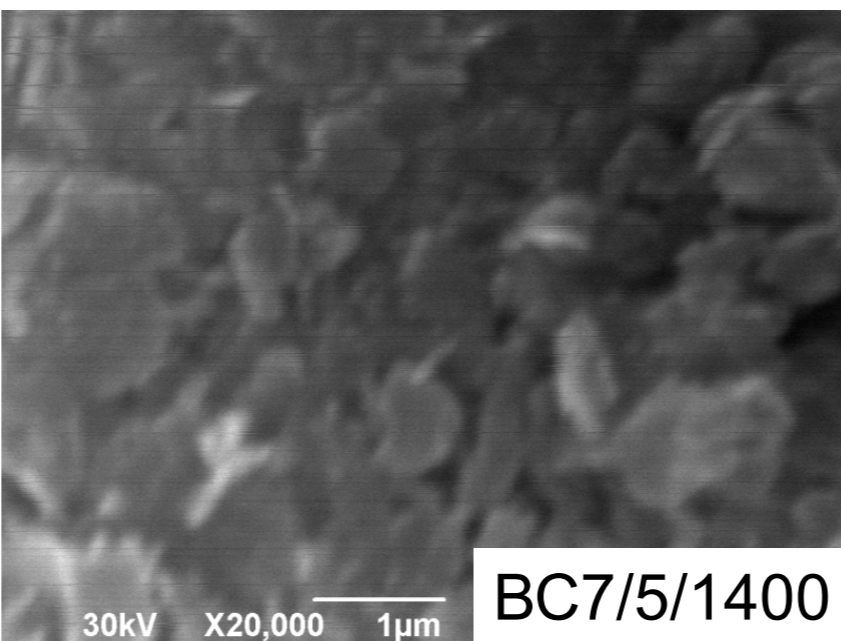
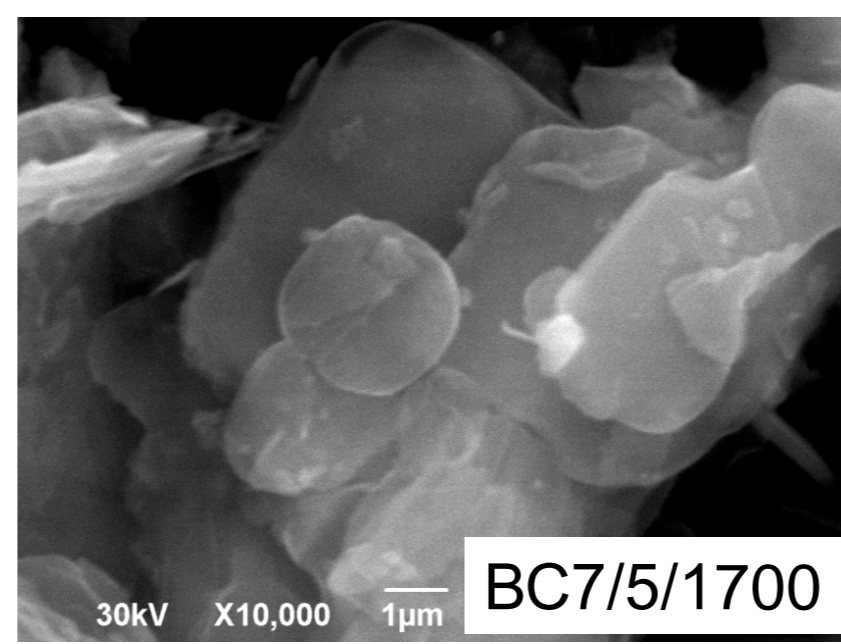
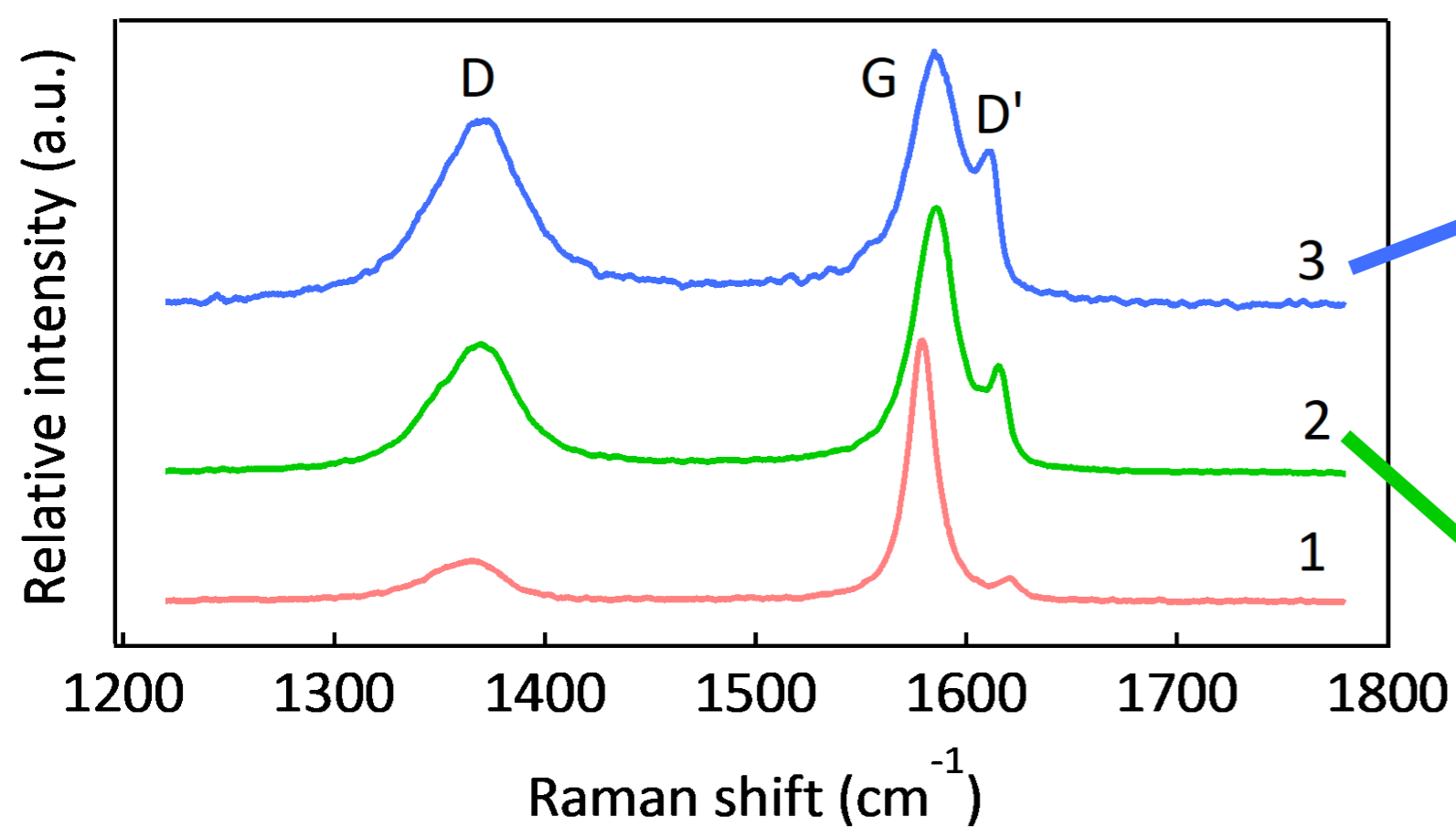


graphite with 2% boron:  
a - 2.467, c - 6.704

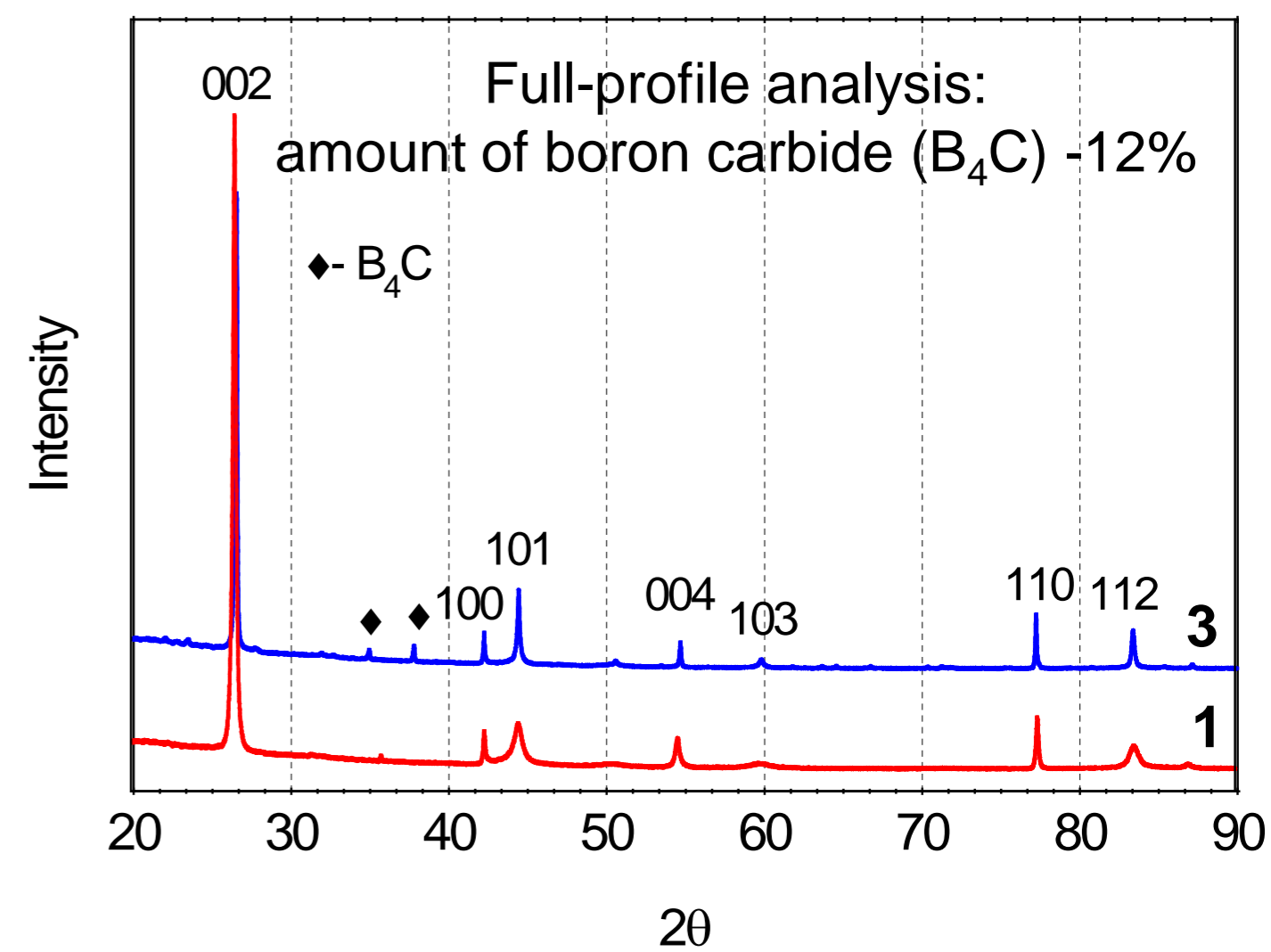
graphite: a - 2.462, c - 6.720

Influence of thermobaric treatment on the parameters of the crystal lattice of the formed graphite

Sample number	Initial state	Processing parameters P, T, τ	a, Å	c, Å	B, at.%
1	Natural graphite	7.0 ГПа, 1700°C, 60sec	2.4622(3)	6.7181(30)	-
2	BC7 mixture	5.0 ГПа, 1400°C, 60sec	2.4644(1)	6.7055(6)	≈ 1
3	BC7 mixture	5.0 ГПа, 1700°C, 60sec	2.4670(2)	6.7035(8)	≈ 2
4	BC7 mixture	7.0 ГПа, 1700°C, 60sec	2.4663(2)	6.7043(13)	≈ 2
5	Globular nanocarbon	7.0 ГПа, 1700°C, 60sec	2.4626(3)	6.7197(21)	-



Perfect graphene layer with 3% boron atoms



The peaks on the X-ray pattern of boron-doped graphite are not broadened, which reflects of a good periodicity of the structure. The apparent contradiction between X-ray and Raman data can be solved by assuming that boron atoms are located in layers periodically and do not disturb the three-dimensional ordering.

Raman spectra of boron-doped graphites show a significant structural imperfection. The width of G peak, intensity of D and D' peaks increase with the graphitization temperature increase.

Solid-state conversion of heavily boron doped graphite into heavily boron doped diamond occurs at a record low pressures of 7.5-8.0 GPa and temperatures of 1600-1650°C.

