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Journal of the European Ceramic Society 26 (2006) 2035-2038

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Crystal structure and microwave dielectric properties of $(Ba_{1-\alpha}Sr_{\alpha})_{6-3x}Sm_{8+2x}Ti_{18}O_{54}$ solid solutions

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Available online 8 November 2005

Abstract

The tungstenbronze-type-like $(Ba_{1-\alpha}Sr_{\alpha})_{6-3x}R_{8+2x}Ti_{18}O_{54}$ (*R* = rare earth) solid solutions have been studied. Microwave dielectric properties of these solid solutions link to the substitution of large cations such as Ba ions by small Sr ions. In these solid solutions for *x* = 0, the quality factor (*Q*·*f*) exibits extremely low due to the inner strain resulting by the occupation of large Ba ions at relatively small A1-sites. However, dielectric constant (ε_r) of this composition show a high value. In this paper, improvement of quality factor (*Q*·*f*) by substituting small Sr ions for Ba ions in $(Ba_{1-\alpha}Sr_{\alpha})_{6-3x}Sm_{8+2x}Ti_{18}O_{54}$ solid solutions, where *x* = 0, is reported. In addition, the relationship between crystal structure and microwave dielectric properties is discussed from the viewpoint of structural formula and occupational state of large cations such as Ba, Sr and Sm in A1-sites. \bigcirc 2005 Elsevier Ltd. All rights reserved.

Keywords: Tungstenbronze-type-like structure; Dielectric properties; Powders-solid state reaction; X-ray methods

1. Introduction

Microwave dielectric materials have been used widespread for applications such as oscillators and filters in the microwave communications. The desirable properties in microwave dielectric materials are a high dielectric constant (ε_r), low dielectric loss $(\tan \delta)$ (i.e. high quality factor $(Q \cdot f)$), and temperature coefficient of resonate frequency ($\tau_{\rm f}$) close to zero. Materials with a high dielectric constant are possible to reduce the size of microwave devices because the wavelength (λ) in dielectrics is inversely proportional to $\sqrt{\varepsilon_r}$ of the wavelength (λ_0) . One of dielectric materials with high dielectric constant (ε_r) is tungstenbronze-type-like Ba_{6-3x} R_{8+2x} Ti₁₈O₅₄ (*R* = rare earth) solid solutions. Fig. 1 shows the crystal structure of tungstenbronze-type-like $Ba_{6-3x}R_{8+2x}Ti_{18}O_{54}$ solid solutions. The crystal data of R = Sm system with x = 0.71 is as follows: orthorhombic, *Pbnm* (No. 62), *a*=12.13 (1), *b*=22.271 (5), c = 7.639 (5) Å and Z = 2.1 These solid solutions are constituted of Ti(1) - Ti(5) octahedral sites with Ti, A2(1) - A2(2)pentagonal sites with Ba and A1(1) - A1(5) rhombic sites with

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0955-2219/\$ - see front matter © 2005 Elsevier Ltd. All rights reserved. doi:10.1016/j.jeurceramsoc.2005.09.086 rare earth and Ba. These solid solutions have superlattice with doubled *c*-axis by tilting of TiO_6 octahedron (Fig. 2).¹ The relationship between these crystal structure and microwave dielectric properties has been discussed. Ohsato et al.² reported that the quality factor (Q,f) of R = Sm system increases from 429 to 10548 GHz as increasing x value from 0.3 to 2/3. The structural formula of R = Sm system with $0 \le x \le 2/3$ is as follows: $[Ba_{2-3x}Sm_{8+2x}V_x]_{A1}[Ba_4]_{A2}Ti_{18}O_{54}$ (V: Vacancy). Where $0 \le x < 2/3$, Ba ions occupy A1-sites that is slightly small for Ba ions, and create inner strain. This inner strain makes quality factor (Q:f) decrease. On the other hand, quality factor (Q:f), where x = 2/3, shows maximum value, which resulted from the lowest inner strain and the highest lattice stability. These behaviors are lead by rare earth ions and Ba ions occupying A1 and A2-sites, respectively. Imaeda et al.³ studied $(Ba_{1-\alpha}Sr_{\alpha})_{6-3x}Sm_{8+2x}Ti_{18}O_{54}$ (BSST) solid solutions where x = 0.6 and $0 \le \alpha \le 0.2$. As according to this report, a quality factor $(Q \cdot f)$ of these solid solutions was improved by increasing lattice stability, which resulted from substituting Sr ions with smaller ionic radius for Ba ions.

In this paper, the crystal structure of BSST solid solutions is demonstrated. Furthermore, microwave dielectric properties of BSST solid solutions, where x = 0 and $\alpha = 0-1/3$, are reported. Relationship between crystal structure and microwave dielectric



Fig. 1. The tungstenbronze-type-like structure of $Ba_{6-3x}R_{8+2x}Ti_{18}O_{54}$ (*R* = rare earth) solid solutions.

properties is discussed from the viewpoint of structural formula and occupational state of large cations such as Ba, Sr and Sm in A1-sites.

2. Experimental

Single crystal of BSST solid solutions was synthesized by cooling method from the liquid phase. High purity BaCO₃, SrCO₃, Sm₂O₃ and TiO₂ were weighed out composition x = 0 and $\alpha = 1/3$. After mixing for 24 h in ethanol, the dried mixture was calcined in air at 1000 °C for 2 h. The calcined powder was melted at 1650 °C for 5 h and cooled 2 °C/h for 1650–1400 °C, 4 °C/h for 1400–1300 °C, 1 °C/min for 1300–1050 °C, and 5 °C/min for 1050 °C to room temperature. The X-ray diffraction data were obtained by a single crystal diffractometer with imaging plate (Rigaku; R-AXIS RAPID). The structural parameters were refined by a full-matrix least-squares method that is termed RADY (S. Sasaki, 1982).⁴

Ceramics resonators of BSST solid solutions, where x=0 and $\alpha = 0-1/3$, were synthesized conventional solid-state reac-



Fig. 2. Superlattice and fundamental lattice of $Ba_{6-3x}R_{8+2x}Ti_{18}O_{54}$ (*R* = rare earth) solid solutions.¹

tion method. High purity BaCO₃, SrCO₃, Sm₂O₃ and TiO₂ were weighed out various α values from 0 to 1/3 and ball-milled for 24 h in ethanol. The mixture was dried and calcined in air at 1250 °C for 2 h. The calcined powders with organic binder were mixed using alumina mortar and pressed into discs 12 mm in diameter. These discs were sintered in air at 1300 °C for 2 h. Sintered discs were polished and annealed in air at 1000 °C for 2 h. Microwave dielectric properties were measured by the Hakki & Coleman method.⁵ The crystalline phase of samples were identified by X-ray diffraction using Cu K α with Ni filter (Phillips; X'Pert MPD).

3. Results and discussion

Crystal data of BSST solid solutions with superlattice were as follows: orthorhombic, *Pbnm* (No. 62), a = 12.174 (4), b = 22.378 (9), c = 7.670 (2) Å and Z = 2. The structural parameters were refined with compositions which were fixed at x = 0 and $\alpha = 1/3$. According to the refined results, *R*-factor and weighted *R*-factor were 4.01 and 5.14%, respectively. Structural parameters of BSST (x = 0, $\alpha = 1/3$) solid solutions are listed in Table 1. All Sr ions occupied A1-sites. Especially, A1(1) and A1(3)-sites had more amount of Sr ions as compared with another A1-sites.

The XRD patterns of sintered ceramics, where $0 \le \alpha \le 1/3$, are shown in Fig. 3. Ohsato et al.⁶ reported that the domain of solid solutions is limited by the composition, where $0.3 \le x \le 0.7$, in Ba_{6-3x}Sm_{8+2x}Ti₁₈O₅₄ solid solutions. Therefore, BSST solid solutions, where x = 0.3, were obtained in the case of the starting composition that was weighed out as x = 0. As a result, Ba and Sr were surplus and formed secondly phases such as (Ba, Sr)TiO₃ solid solutions with all composition α .

Microwave dielectric properties of BSST solid solutions, where x=0 and $\alpha=0-1/3$, are shown in Fig. 4. Quality factor (*Q*:*f*) of Ba_{6-3x}Sm_{8+2x}Ti₁₈O₅₄ solid solutions, where x=0, shows extremely low values.¹ However, quality factor (*Q*:*f*) of BSST solid solution, where x=0, was improved with increasing α . The aforesaid remarked that BSST solid solutions, where x=0.3, are obtained in the case of the starting composition that is weighed out as x=0. Structural formula of Ba_{6-3x}Sm_{8+2x}Ti₁₈O₅₄ solid solutions, where x=0.3, is as follows: [Ba_{1.1}Sm_{8.6}V_{0.3}]_{A1}[Ba₄]_{A2}Ti₁₈O₅₄ (V: Vacancy). When Ba ions are substituted by Sr ions, at first Ba ions in



Fig. 3. XRD patterns of BSST solid solutions where x = 0 and $\alpha = 0$ to 1/3.

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