Ba_{2-x} Pb_xCa₂Cu₃O_{8+δ} superconductors Up 99K

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ABSTRACT

In the present study the effect of replacement of Pb at the Ba place in the oxygen deficient BaO layer of $Ba_{2-x} Pb_x Ca_2 Cu_3 O_{8+\delta}$ with ($0 \le x \le 0.3$) cuprate superconductor has been investigated. Bulk polycrystalline specimens have been prepared by the solid state reaction process. Four probes technique is used to find the T_c. The highest $T_{c(offset)}$ were 88, <77, 97and 99 K for all specimens with x = 0.0, 0.1, 0.2, and 0.3 respectively. The optimum $T_{c(off)}$ of ~ 99 K and $T_{c(onset)} \sim 106.38$ K has been found for the composition $Ba_{1.7} Pb_{0.3}Ca_2Cu_3O_{8+\delta}$. X-ray diffraction analysis showed an orthorhombic arrangement with an increase of the c-axis for the specimens substitution with Pb as compared with those of no Pb content. The change of the Pb content of all specimens created a change in the volume of the unit cell which has an important effect to obtain the superconductor state.

Keywords: superconducting transition temperature T_c , electrical resistivity, Pb content, Ba_{2-x} Pb_xCa₂Cu₃O_{8+ δ} superconductor

INTRODUCTION

uperconductor compounds are Multilayered cuprates and generally described by the formula of MCa_{n-1}Cu_nO_{2n}, where M refer to Bi₂Sr₂[1], Bi₂Ba₂[2], HgSr₂[3], HgBa₂[4,5], ✓ TlBa[6]. The general formula for the Superconducting series of Ba-based is given with the following equation $Ba_2Ca_{n-1}Cu_nO_{2n+2}[Ba-02(n-1)n]$, where n is the number of CuO₂ layers, this compound consists of a mixture of Ba₂Ca₁Cu₂O₆ (0212), Ba₂Ca₂Cu₃O₈ (0223) phases, $Ba_2Ca_3Cu_4O_{10}(0234)$ phasesandBa₂Ca₄Cu₅O₁₂(0245) phases. These systems exhibit superconductive transition temperatures T_c of 90 - 120-105 K and 90 K, respectively [7]. It found that the transition temperature (T_c) to the Ba₂Ca_{n-1}Cu_nO_{2n+2} compound do not depend on the number n of Cu-O planes, with increasing n from5to9,But on the high pressure where temperature T_c reached to about 80 K [8]. Also the superconductivity was found in apical fluorine system of Ba₂CaCu₂O_{6-v}F_v. The highest T_c of about 108 K has been achieved for the synthesized specimen from a nominal composition of $Ba_2CaCu_2O_{4.4}F_{1.6}$ under high pressure with different (F) [9] "contents. SKOBA, et al"[10],"deposited of superconducting films Ba₂Ca_n- $_1Cu_nO_{2n}$ (O,F)₂ on a SrTiO₃(100) substrate by using Nd-YAG pulsed laser ablation under an O₂ pressure of 16-19 Pa. The films exhibited a superconducting onset temperature ($T_{c(onset)}$) of 76.5 K and a zero resistance temperature (T_{c(offset)}) of 22.5 K" [10]. In this research fabrication of Ba based [02(n-1)n], n=3 superconductors of nominal composition Ba_{2-x} Pb_xCa₂Cu₃O_y has been investigated and study the structural and resistance measurements in the temperature range 77K.

Experimental

Purity oxide powders of (BaO, PbO, CaO and CuO) were used as starting materials. The calculated weight depends on molecular weights of these Oxides. The synthesis of specimens have been performed by solid state reaction method, the powders were mixed together by using vortex mixer with grinding time of about 12hrs. Then the powders were pressed into disc shape

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of (25 mm in diameter and 5 mm thickness) using hydraulic press under a pressure of (100MPa) to insure a uniform transmission of pressure during compact and high density. The specimens were heated using carbolite furnace at 860 °C for 24 hours with heating rate 5 °C/min, and then cooled to the room temperature by the same rate. The X-ray diffraction technique was used to determine the crystal structure analysis of the prepared specimens of Ba_{2-x} Pb_xCa₂Cu₃O₈₊₆ with ($0 \le x \le 0.3$) by X-ray diffraction and a computer program (Full Prof Suite toolbar) was used to calculate the lattice parameters [11] and the density of the specimens (d_m) were calculated [12]. "The volume fraction of any phase (Vphase) in the specimen was determined by using the relation"[13]:

"
$$V_{ph} = \frac{\sum I_0}{\sum I_0 + \sum I_1 + \sum I_2 + \dots + \sum I_n} \times 100\%$$
"

"Where I_0 is the XRD peak intensity of the phase which were determined, $I_1, I_2, ... I_n$ are the peaks intensity of all XRD"[13]. A four-point probe is the most common method to determine the T_c of a superconductor.

Rustles and Discussion

According to result of the (XRD) data the structures of all specimens are orthorhombic. From the results in the Table (1), It was found that the lattice parameters (a, b and c) are controlled by Pb content with the change of electrons into anti bonding orbital. "The lattice parameters have been estimated by using d-values and (hkl) reflections of the observed X-ray diffraction pattern through the software program based on Full prof Suite toolbar"[11]. The comparison between of XRD patterns of the (Pb) doped specimens with the relative intensities of the same reflection of the specimen with x = 0.0 (Pb)." The change of the lattice parameters effects on the unit cell volume and causes variation of the density d_m, which calculated by equation"[14]:

 $d_m = W_m / N_A V''$ [14]

"Where NA is the Avogadro's number in unit (particles/g.mol), V is the volume of unit cell and W_m is the molecular weight in unit (amu)"[14].Figure (2) shown the Variation density with Pb content for the Ba_{2-x} Pb_xCa₂Cu₃O_{8+ δ} specimens .The increasing of the density d_m could be attributed to the reduction of the porosity and vacancies in the system. Also from the same table it can be observed that there is an increase in the lattice parameters ratios (c/a of unit cell) for specimens by increasing the lead concentration as shown in Figure (3), while the volume unit cell (V) decreased for specimens with the increasing lead content except (0.3 wt % Pb) may be due to (pb) substitution instead of (Ba) sites. From the ($\rho - T$) curves which shown in Figure (4). The transition temperature of the specimens was found to be sensitive to the (Pb) the concentrations, it has been observed that the optimum $T_{c(off)} \sim 99$ K and $T_{c(onset)} \sim 106.38$ K for the composition $Ba_{1,7}Pb_{0,3}Ca_2Cu_3O_{8+\delta}$ where the (Pb) concentration equal to 0.3 as shown in table 1. "The observed behavior may be attributed to the lead content (Pb) which controls the hole concentrations in conducting CuO planes" [14, 15]. "Thus, induced additions may be change an oxygen contents as expected, which might bring changes in carrier concentration causes the observed values of T_c"[14], or the increase of substitution(Pb) which leads to increase the holes in the structures thus enhanced the transition temperature from 88K to 99K and reduced the transition width ΔT , they mean that the increasing multiphase Ba-0212(low phase) and Ba-0223(high phase) as clear in Table (2), but at the addition x=0.1, the specimen shows similar behavior and may be having low T_c, because it behaves like a superconductor, but does not become zero even at the boiling point of liquid nitrogen 77K.

CONCLUSIONS

In this research the effect of replace of Pb by the Ba place in the oxygen BaO_{δ} layer of Bulk polycrystalline specimens $Ba_{2-x} Pb_xCa_2Cu_3O_{8+\delta}$ with ($0 \le x \le 0.3$) cuprate superconductor have been prepared under optimum conditions. X-ray diffraction analysis showed an orthorhombic

4.05369

7.25054

4.05467

4,09540

21.60619

22.07298

21.59905

23.02507

4.25461

8.57678

4.05288

4.09046

0.0

0.1

0.2

0.3

75.86

83.76

84.70

86.16

24.14

16.24

15.29

22.29

construction with an increase of the c-axis for the specimens doped with Pb as compare with those have no Pb concentration. It was found that the change of the lead (Pb) content of all specimens give a change in the volume of unite cell. The transition temperature of as grown specimens is found to be sensitive to the Pb concentrations, the optimum $T_{c(off)}$ of ~ 99 K and $T_{c(onset)} \sim 106.38$ K for $Ba_{1.7}$ Pb_{0.3}Ca₂Cu₃O_{8+ δ}.

volume fraction (vpn) for the $Ba_{2-x}PD_xCa_2Cu_3O_{8+\delta}$ specimens.								
x	a (Å)	b (Å)	c (Å)	c/a	V(Å) ³	d _m (g/cm ³)	V _{ph} (High) %	V _{ph} (Low)%

372.639

1372.63

354.940

385.717

1.76433

0.48406

1.8916

1.7588

5.078297

2.573573

5.329299

5.628966

Table(1): Lattice parameters (a,b, c) , ratio c/a, density (dm), transition width ΔT and
volume fraction (Vph) for the $Ba_{2-x}Pb_xCa_2Cu_3O_{8+\delta}$ specimens.

Table(2):	$T_{c(off)}(k)$,	$T_{c(on)}(k)$	transition width	$\Delta \mathbf{T}$ and	volume fraction	1(Vph)	for theBa ₂₋
			Ph CarCurOa	snecin	iens		

$x I D_x C a_2 C u_3 O_{8+\delta}$ specimens.						
specimens	Tc(off) (k)	Tc(on) (k)	ΔΤ	Eg(ev)		
Ba2Ca2Cu3O8+δ	88	97.38	9.38	0.026793		
Ba1.9 Pb0.1Ca2Cu3O8+δ	<77	<77	-	-		
Ba1.8 Pb0.2Ca2Cu3O8+δ	97	104.9	7.9	0.029533		
Ba1.7 Pb0.3Ca2Cu3O8+δ	99	106.38	7.38	0.030142		







Figures (1) (XRD) pattern for $Ba2_{-x} Pb_x Ca_2 Cu_3 O_x$ specimens for x=0.0, 0.1, 0.2 and 0.3.



Figure (2) Variation density with Pb content for the $Ba_{2-x} Pb_x Ca_2 Cu_3 O_{8+\delta}$ specimens.



Figure (3) Variation ratio of lattice parameters c/a with Pb content for the Ba_{2-x} $Pb_xCa_2Cu_3O_{8+\delta}$ specimens.



Figure (4) The resistivity as function of temperature for $Ba_{2-x} Pb_x Ca_2 Cu_3 O_{8+\delta}$ specimens for x=0.0, 0.10, 0.20 and 0.30.

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