

Dielectric and Transport Behavior of Dual Doping (Sm-Ce) in $\text{Bi}_2\text{Ca}_2\text{CoO}_6$ Perovskite Cobaltite



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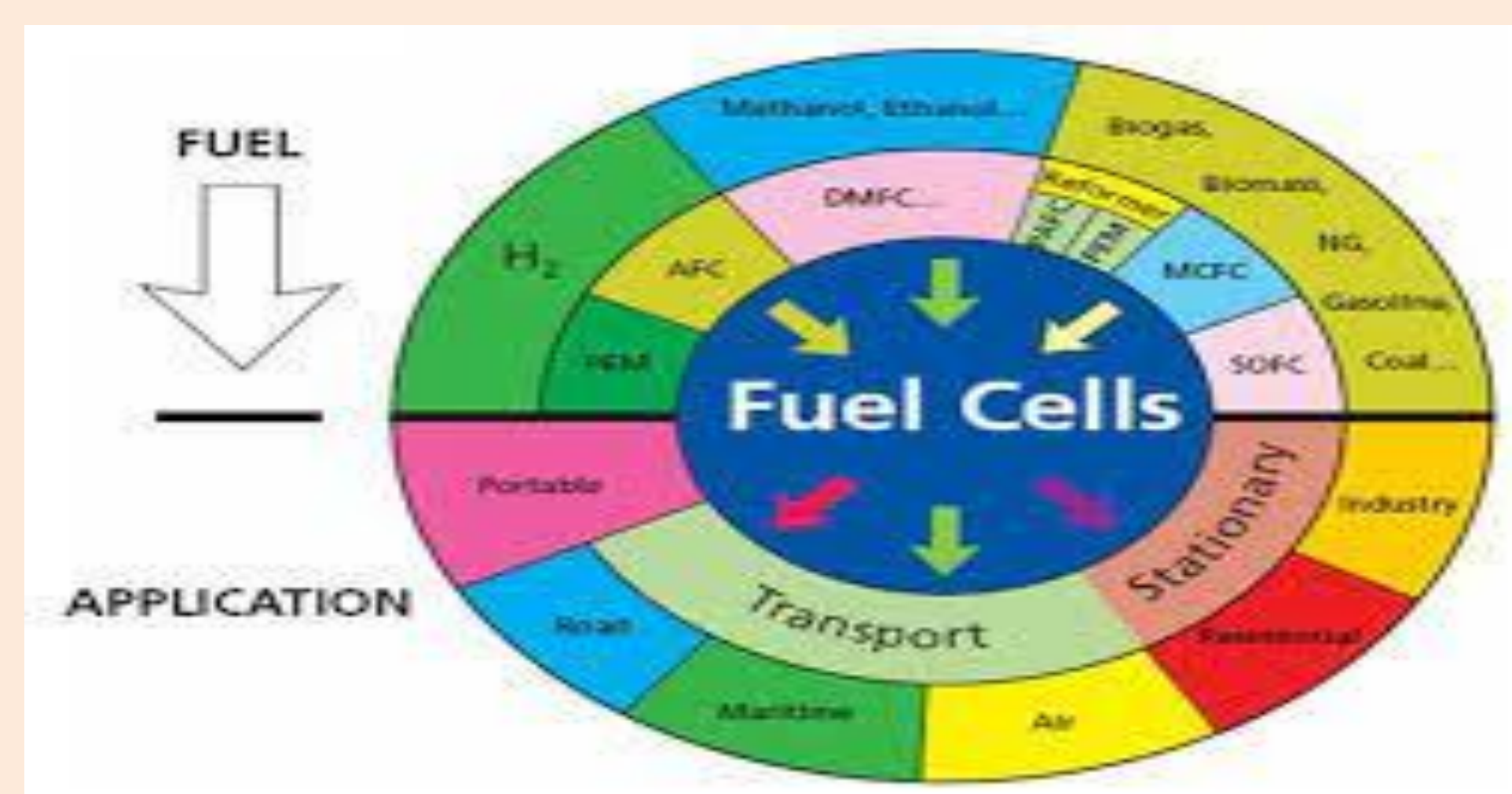


Abstract

- $\text{Bi}_2\text{Ca}_{1.80}\text{Sm}_{0.10}\text{Ce}_{0.10}\text{CoO}_6$ (BCSCCO) perovskite were synthesized the using co-precipitation route
- XRD analysis confirms the monoclinic structure with space group P21/m
- The obtained results are validated using the Jonscher power law.

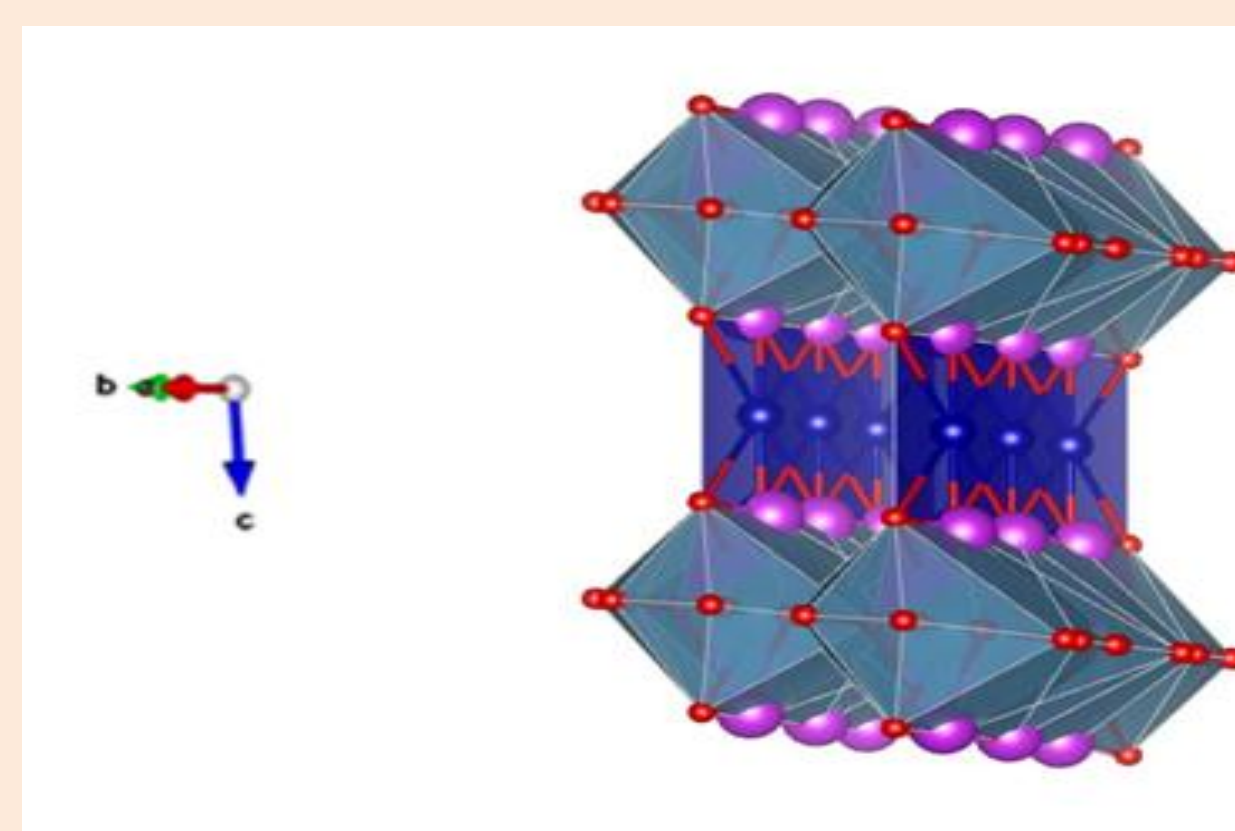
Introduction

- Fossil fuels are the major sources of our energy production in current era, which are limited by nature. Geologists are expecting the current reserves to last up-to 2042 [1].

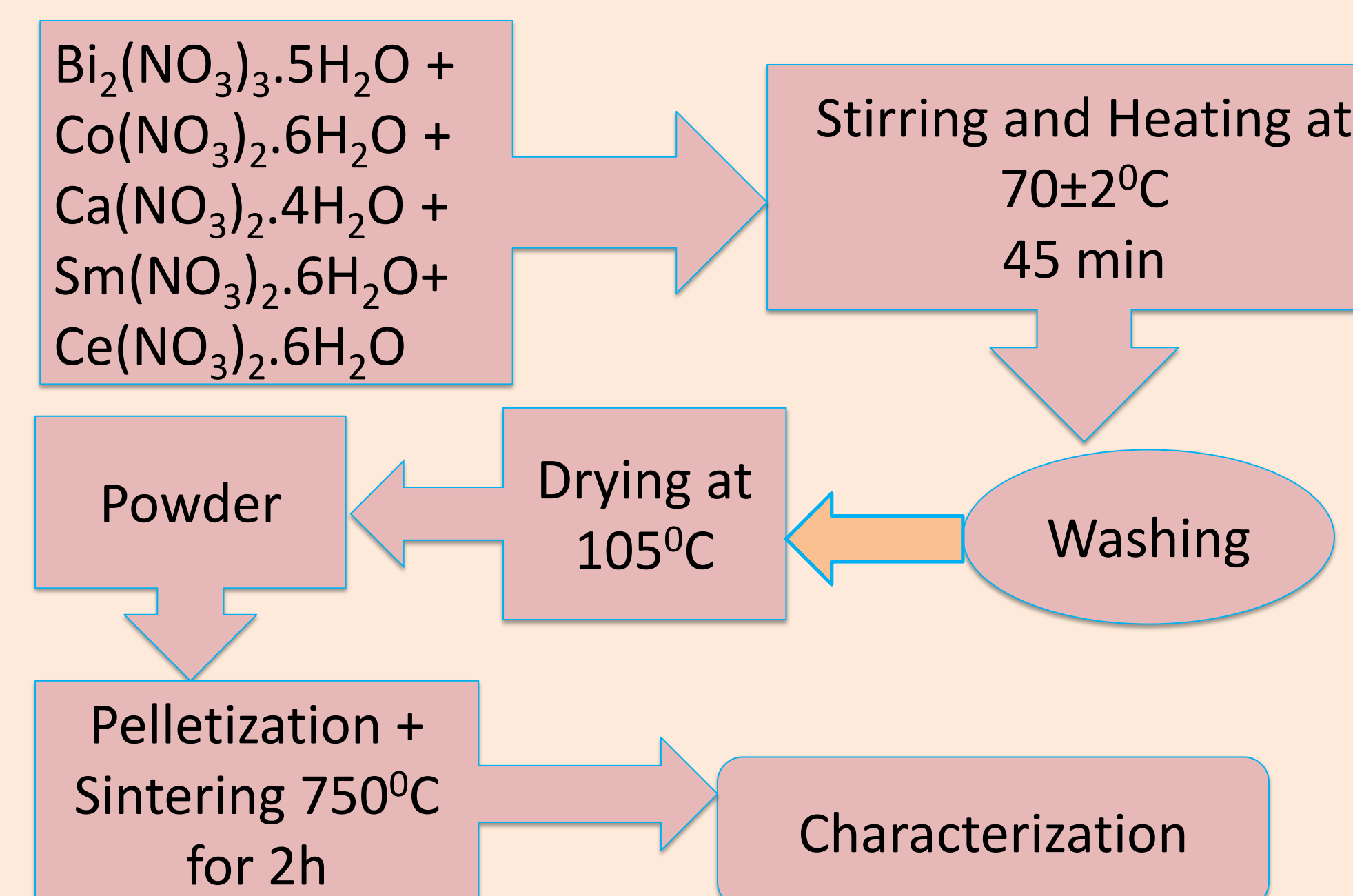


Crystal Structure

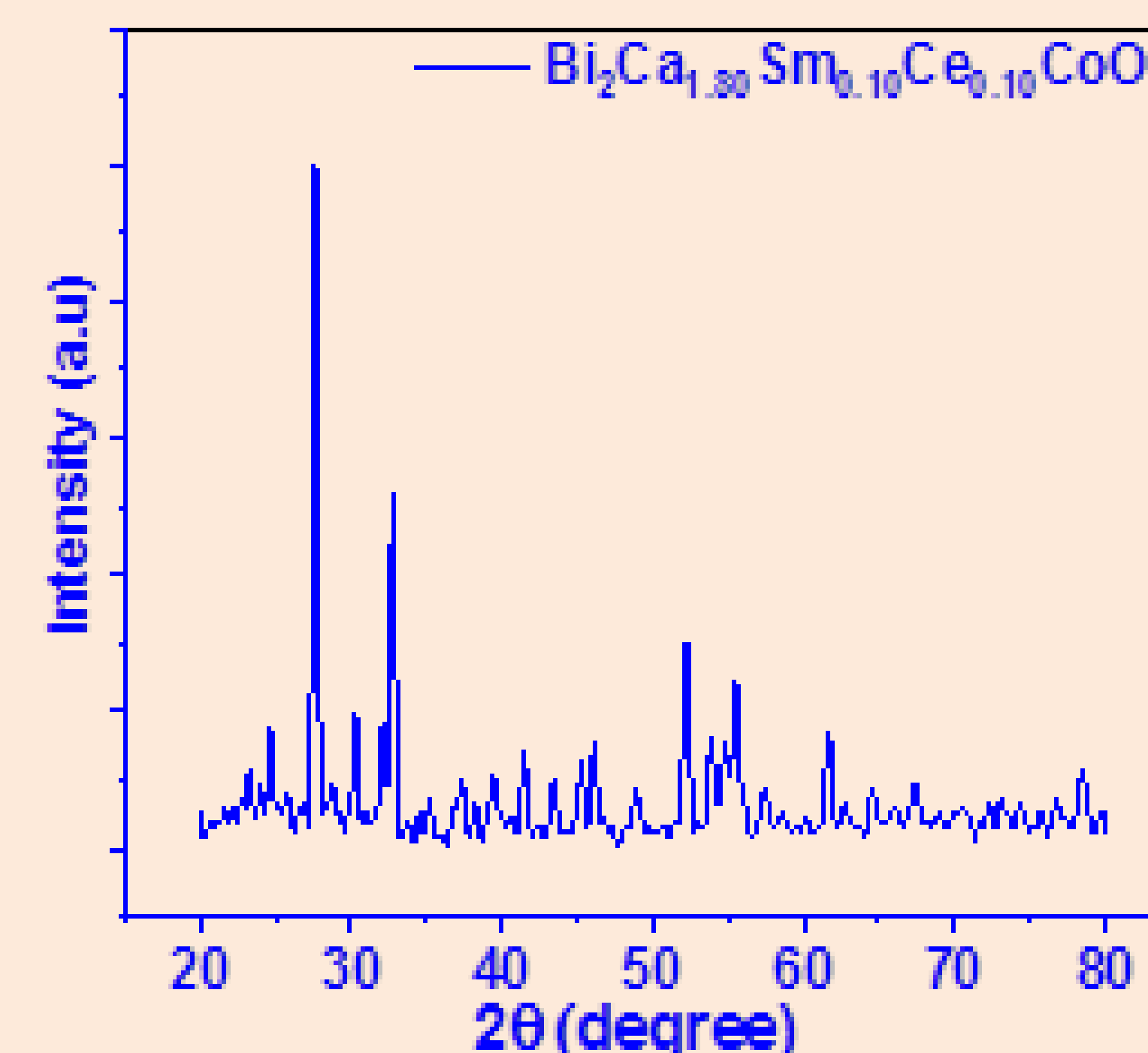
$\text{Bi}_2\text{Ca}_2\text{CoO}_6$
Crystal System



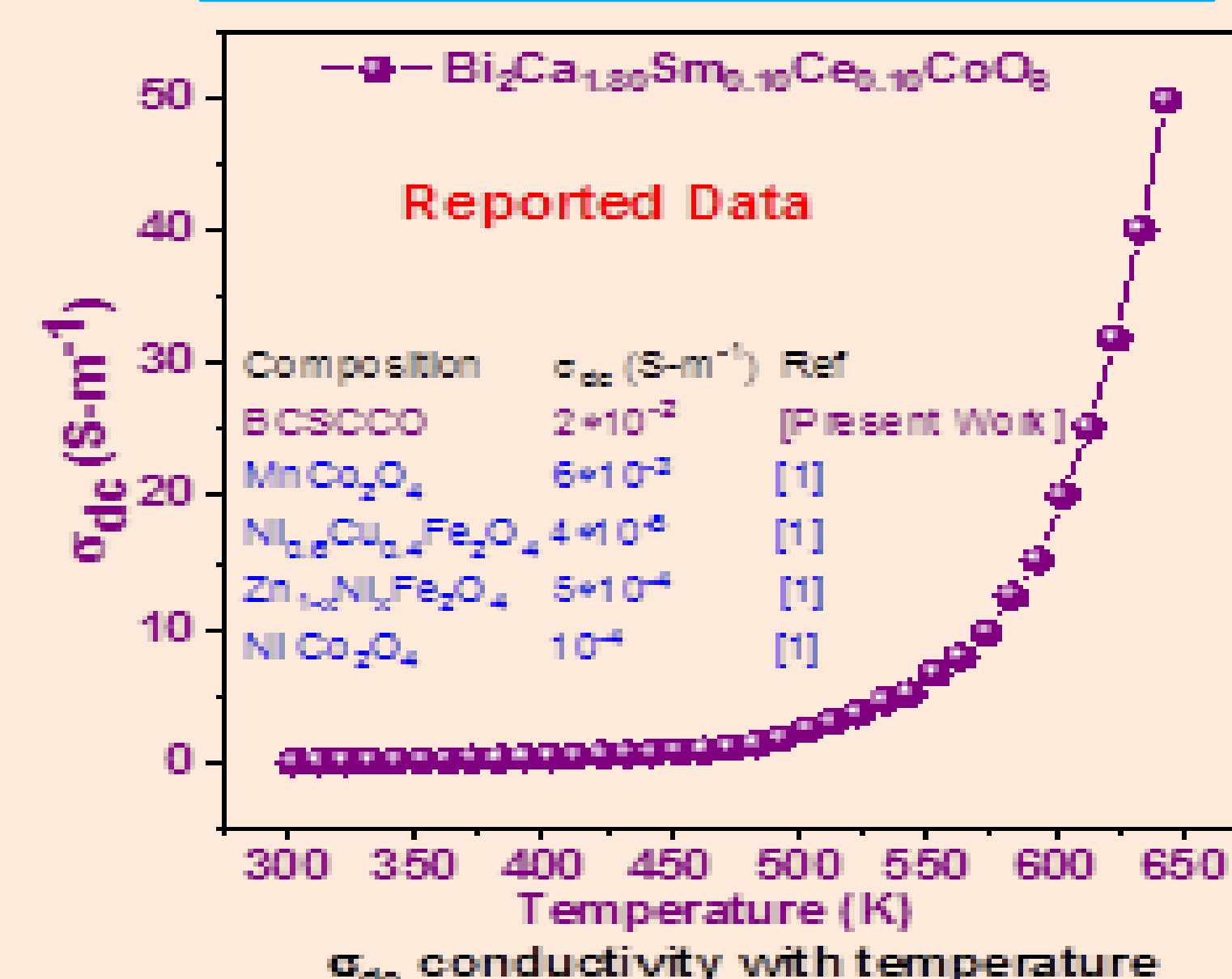
Synthesis Route



Results



Structural analysis

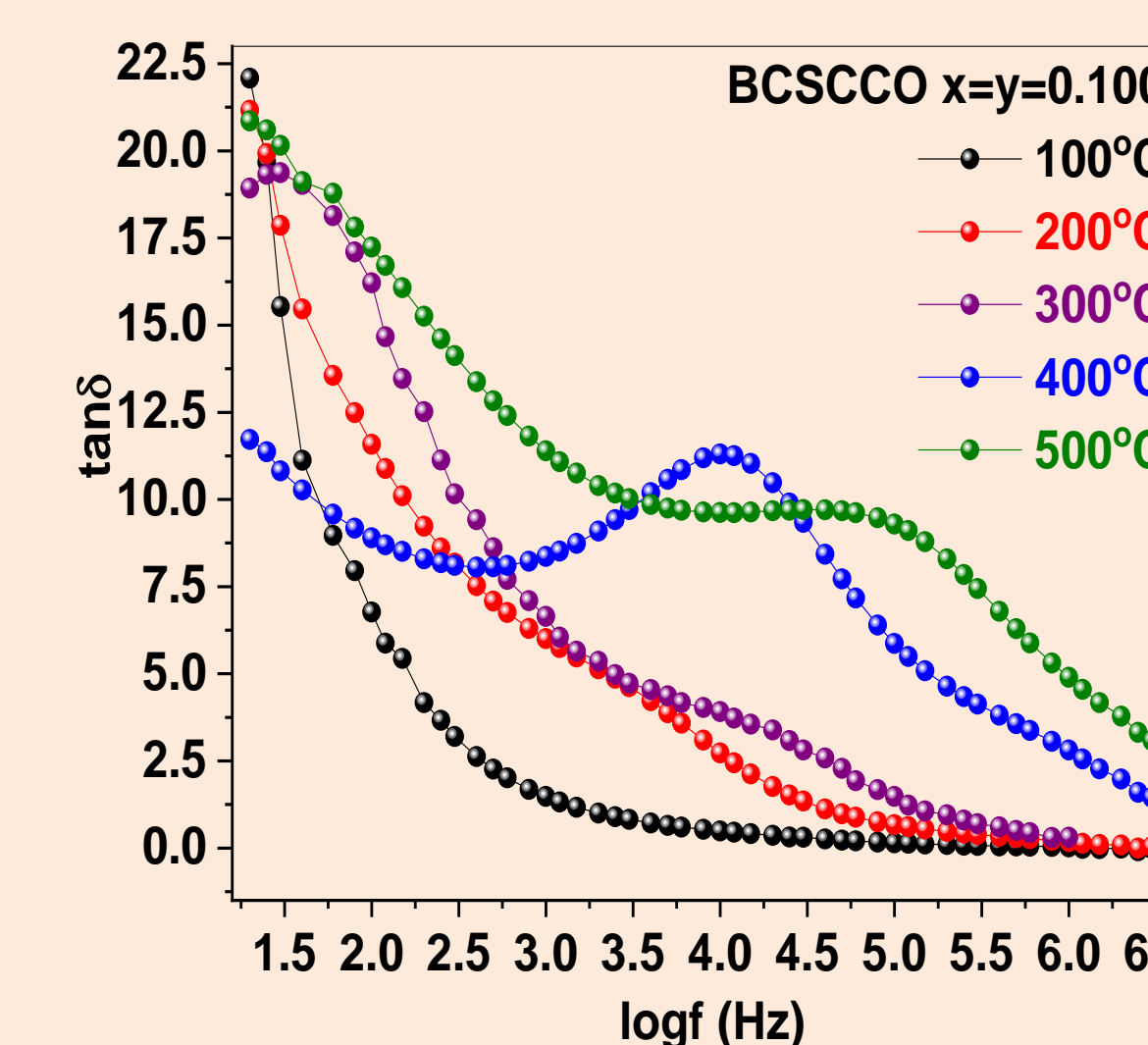


DC electrical analysis

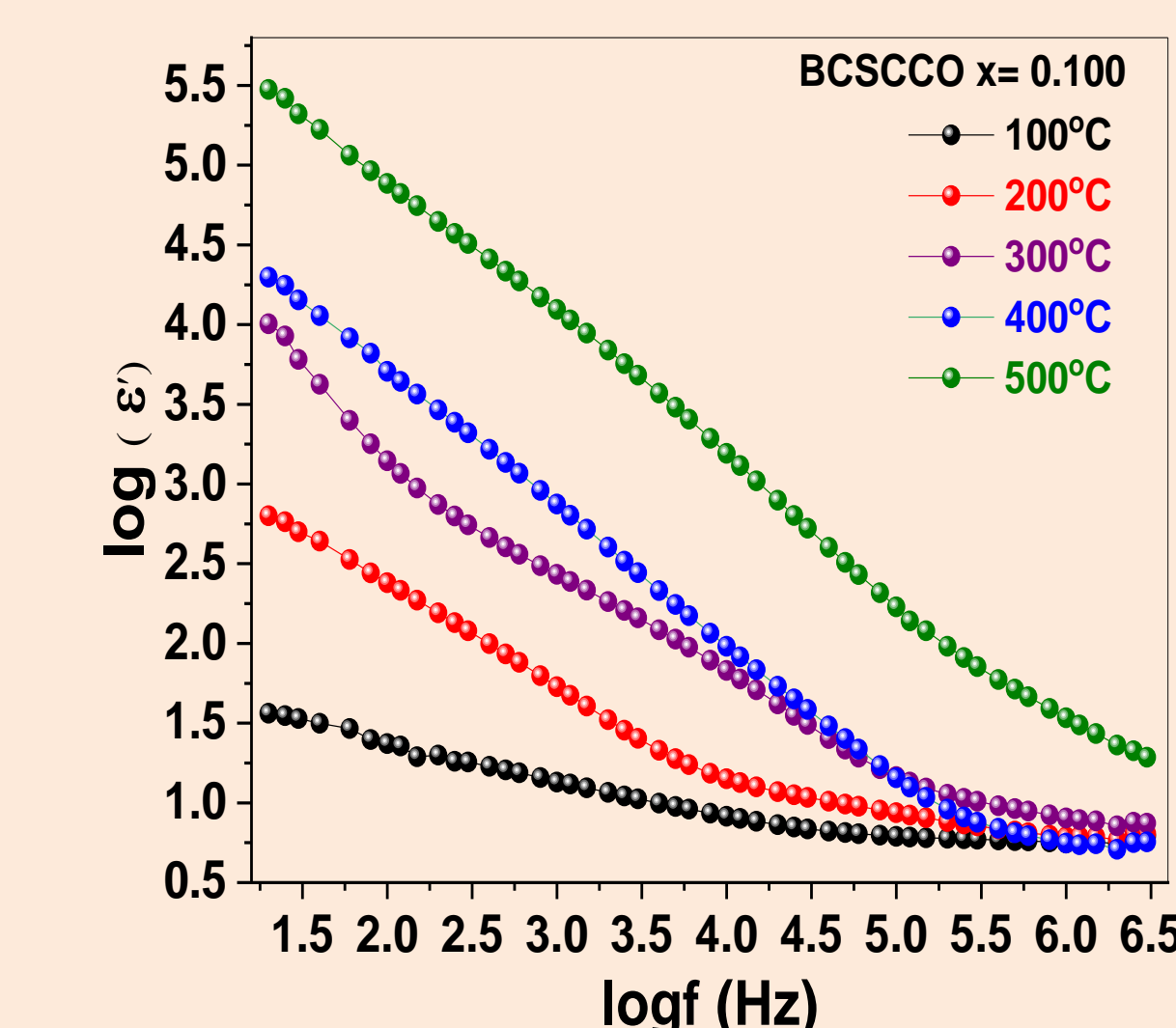
Acknowledgement



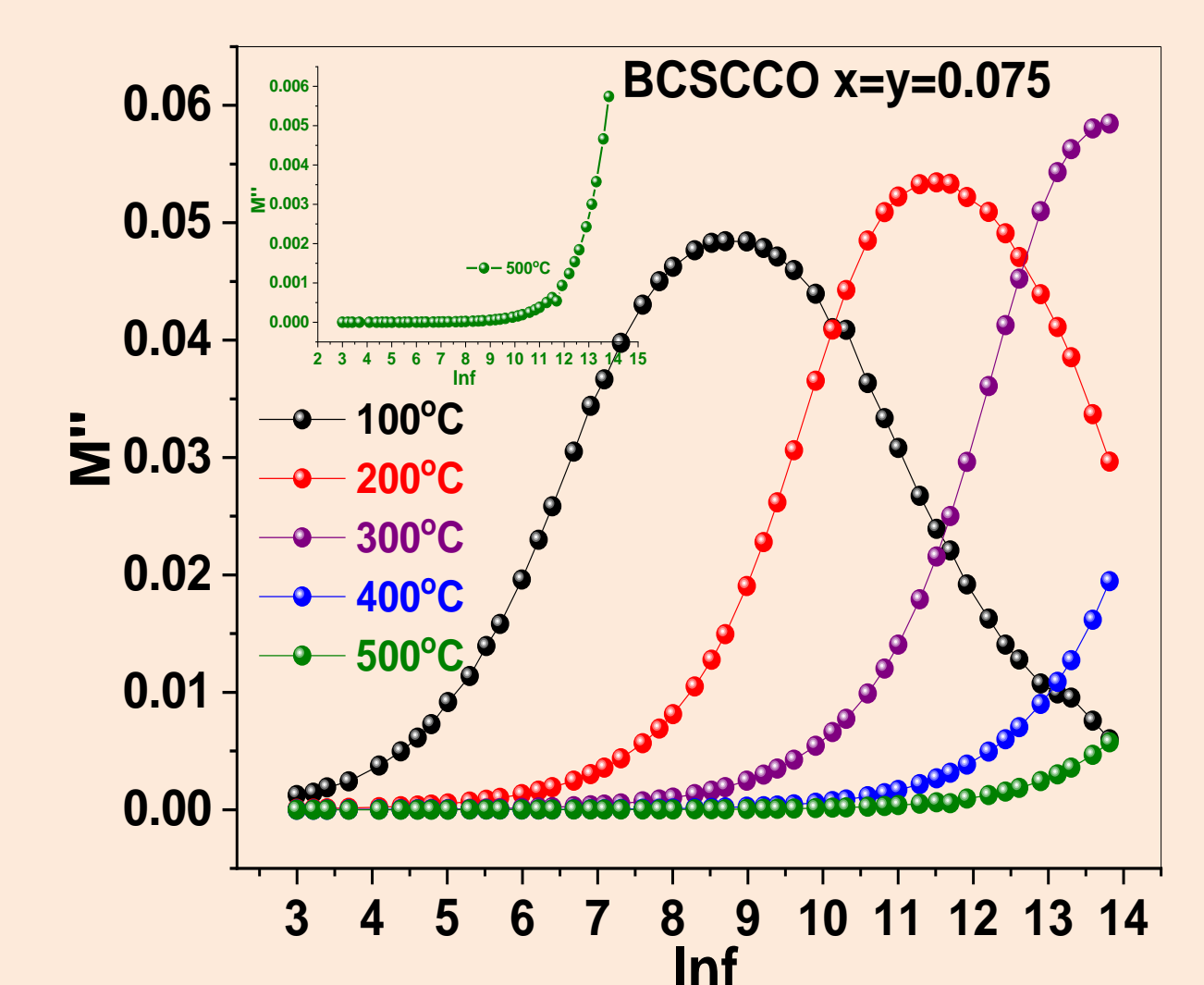
Results



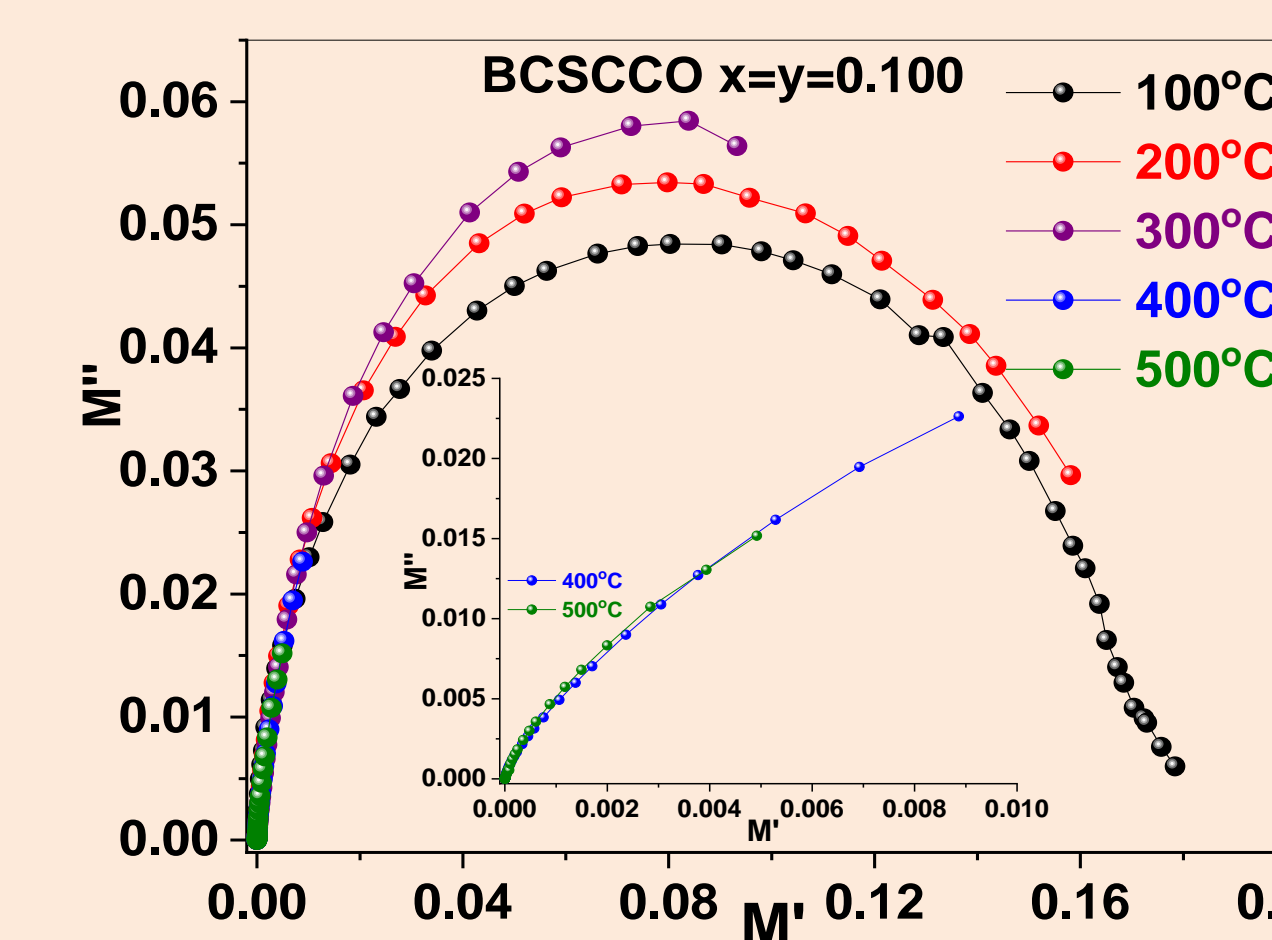
Loss Tangent with temperature



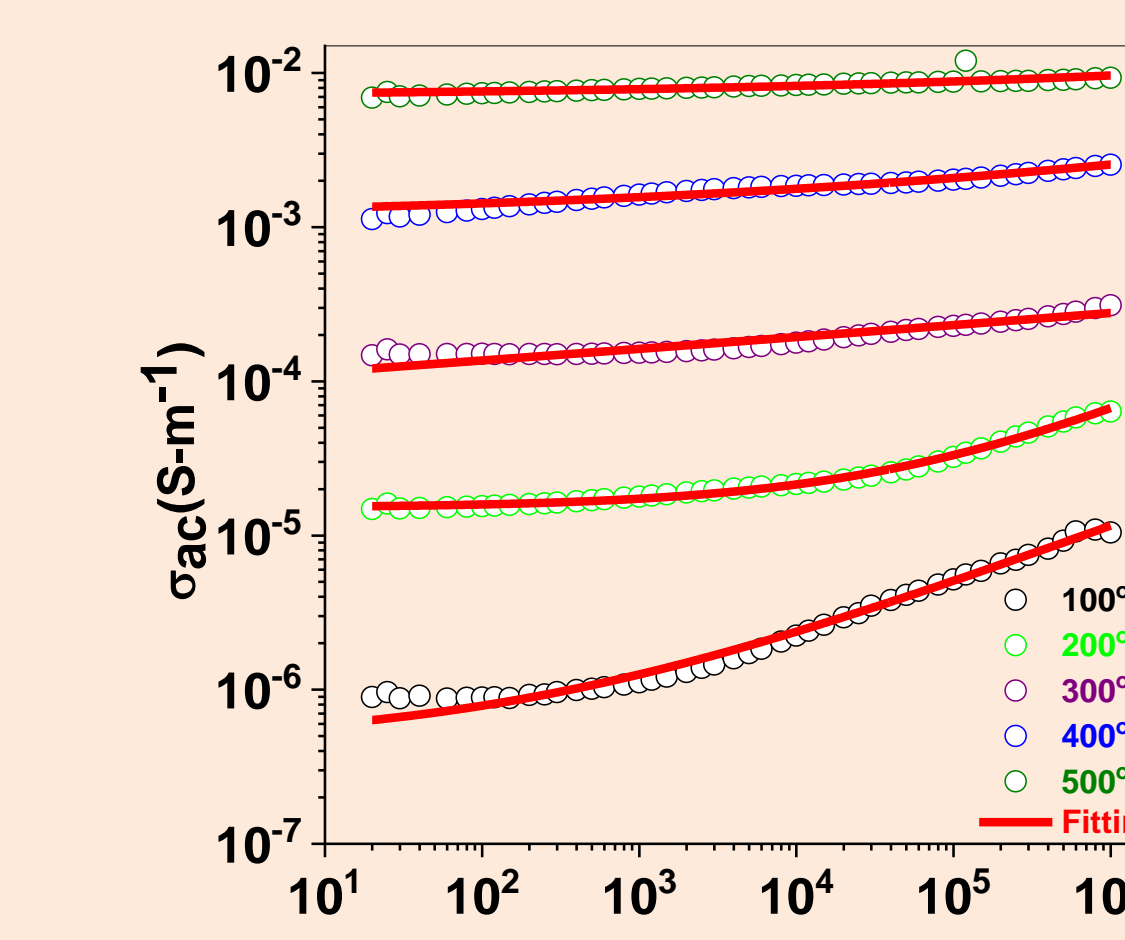
Dielectric Constant



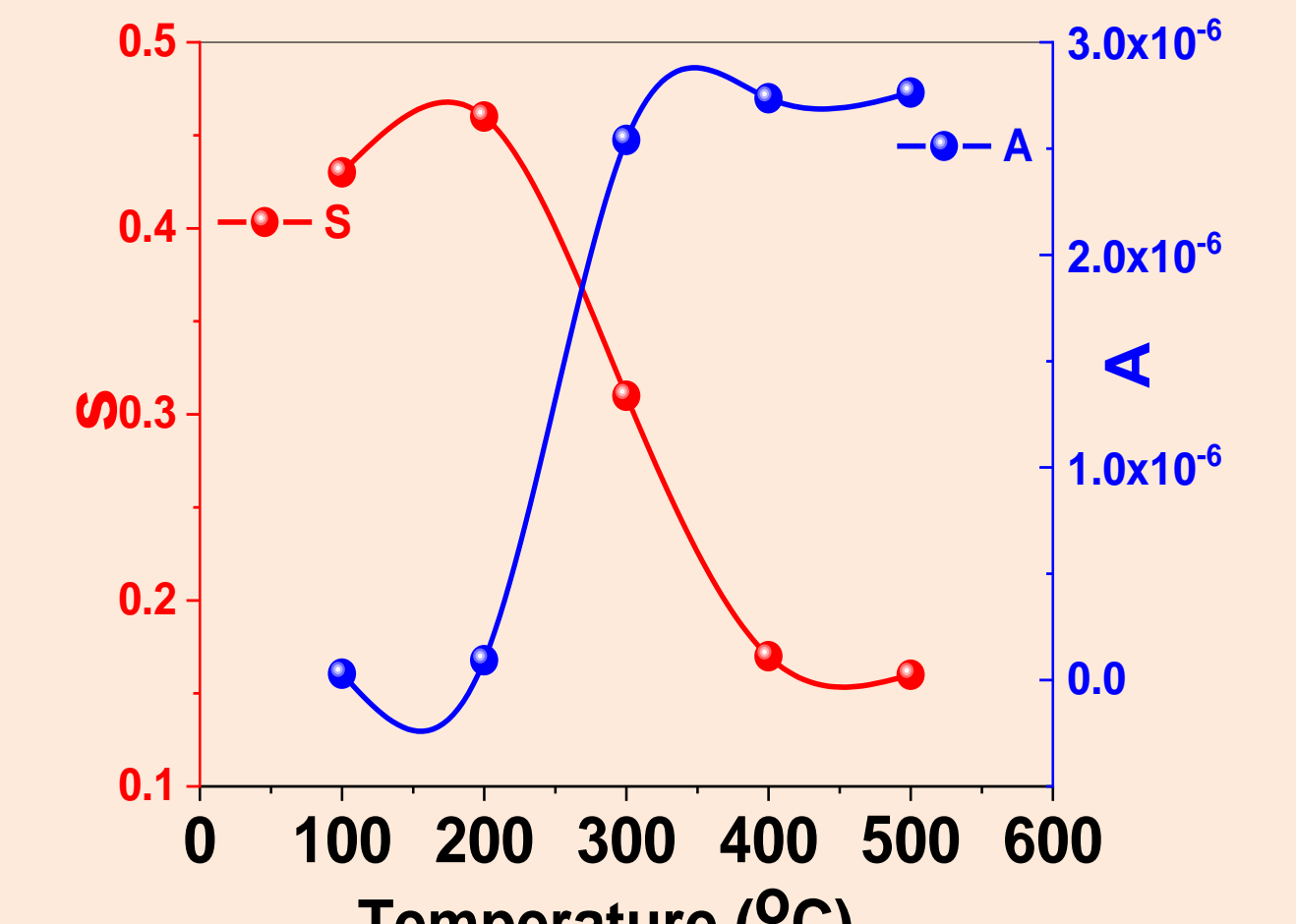
Imaginary part of modulus



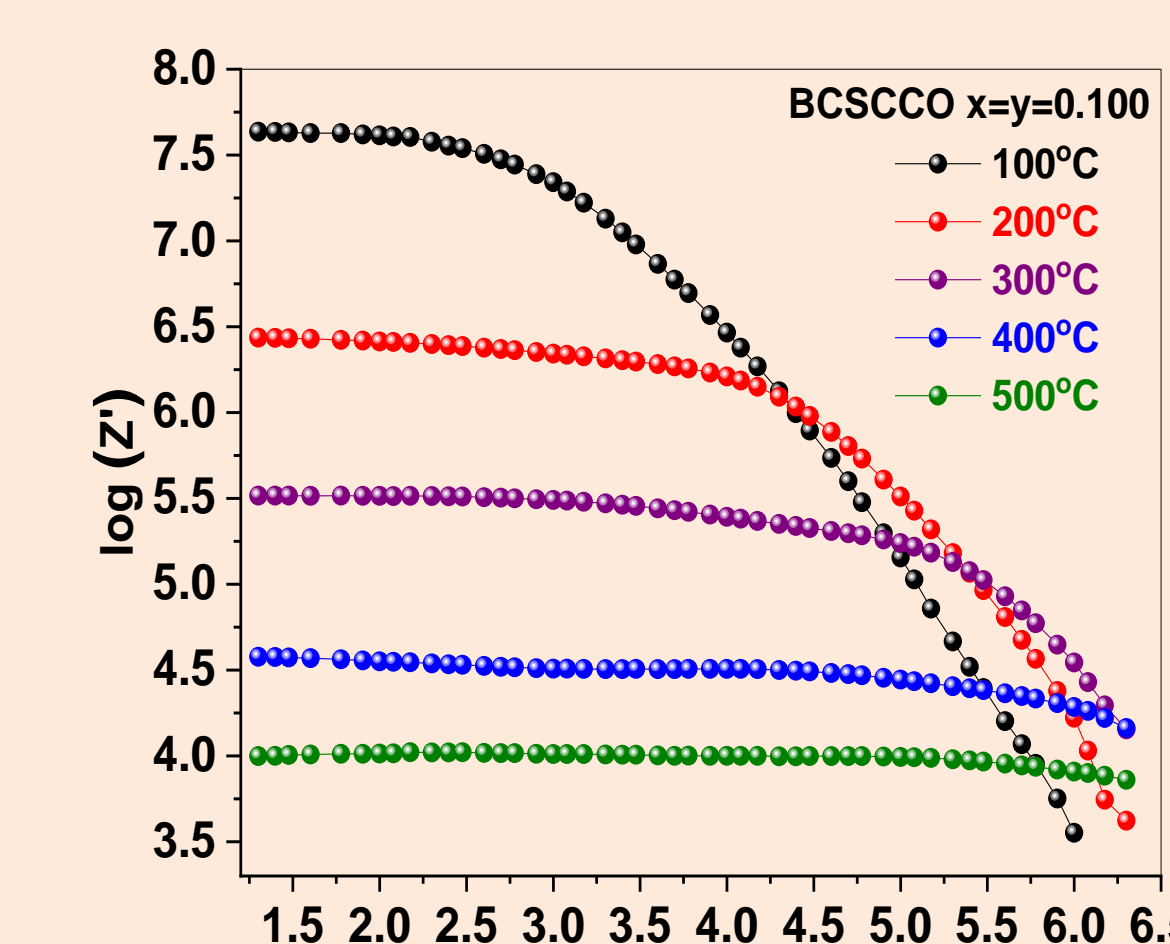
Modulus Spectroscopy



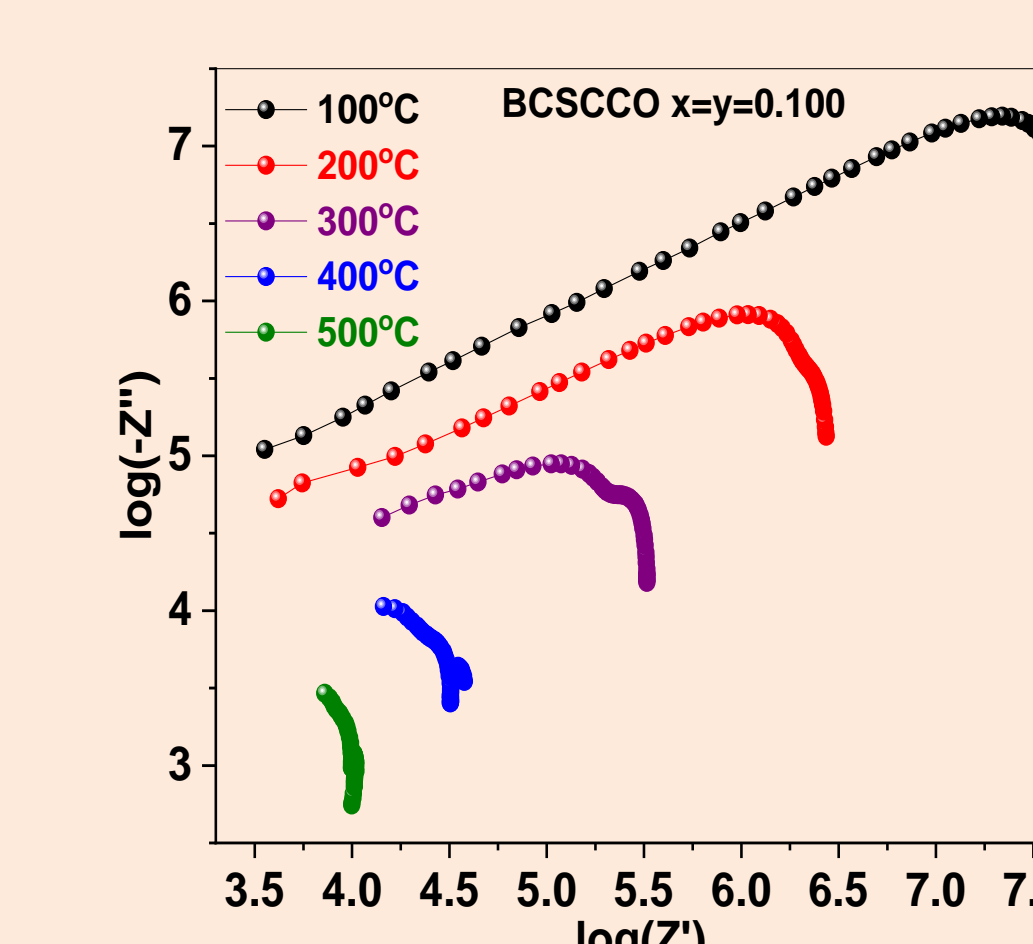
Jonscher Fitting



Variation of s and A with temperature



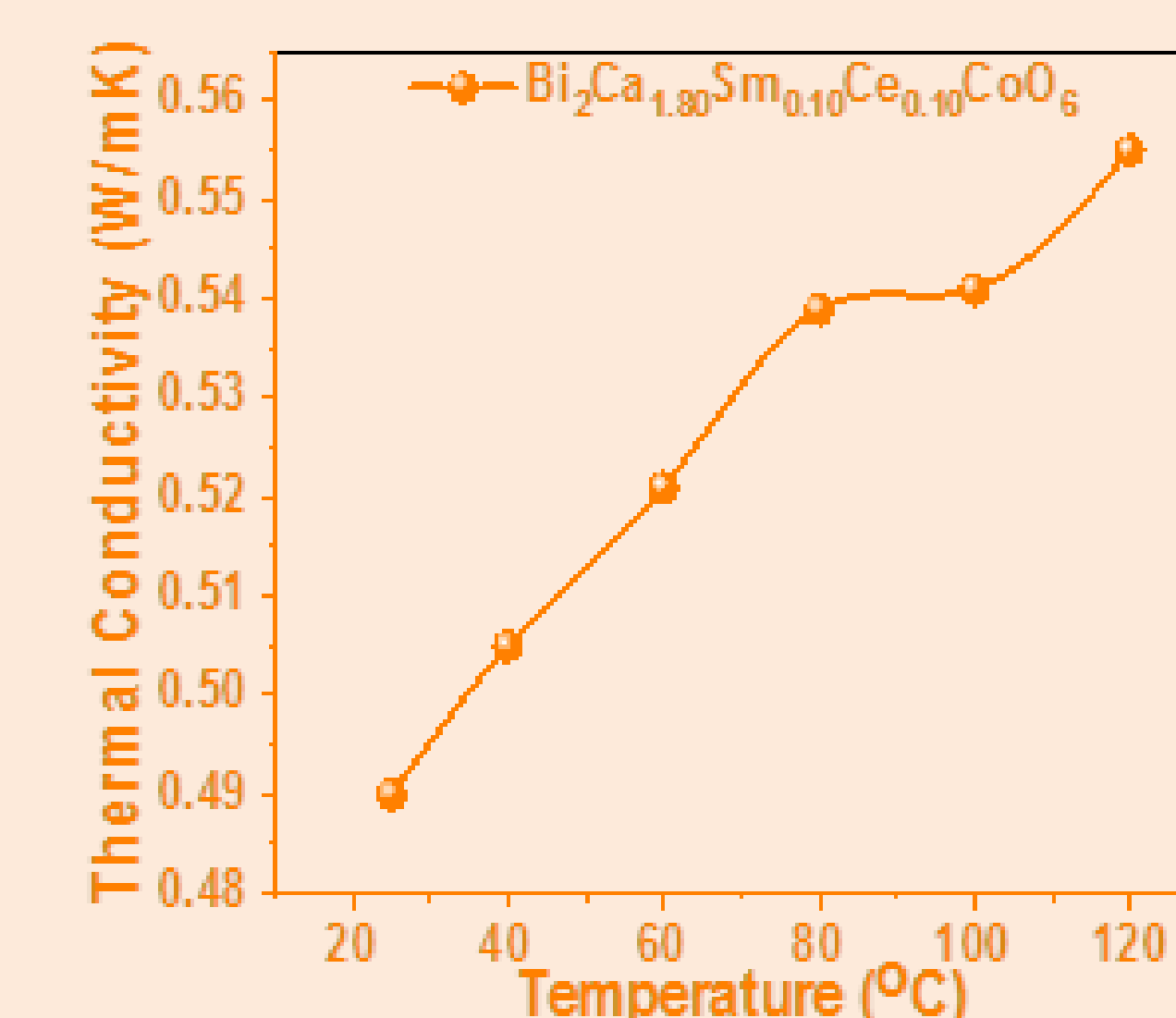
Resistive part of impedance



Cole-Cole plot

References

- [1] Muraleedharan, S., Thiruvengatam, V., Karayi, S. O., Karupiah, K., Jawahar, V., & Ashok, A. M. (2020). Investigation on temperature-dependent electrical properties of $\text{La}_{1-x}\text{A}_x\text{CoO}_3$ (A=La, Li, Mg, Ca, Sr, Ba). CrystEngComm, **22**(1), 85-94



Thermal conductivity