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~~Refinement using FullProf Software Part I |  
Electroceramics Lab Plot FullProf  
Rietveld Refinement Graph in Origin -  
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How to do labelling of XRD pattern via  
Rietveld Refinement using FullProf suite  
software Rietveld Refinement and Crystal  
Structure of CuFe<sub>2</sub>O<sub>4</sub> using FullProf Suite  
and VESTA Software. ~~Easiest way to do  
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Art of the Book PART ONE Introduction  
to Rietveld Analysis by Prof Ashish Garg,  
IIT Kanpur

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Tutorial of VESTA software for creating  
crystal structures

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Simulate Powder XRD PATTERN using VESTA- [TUTORIAL #3]Xrd Analysis using x'pert Highscore Search and Match  
Whats new in Match! version 3.3 **Quick Reitveld Refinement using Match | Electroceramics Lab** How to estimate Bond Structure via Rietveld Refinement using FullProf Software. XRD Analysis for Phase Determination using X'Pert HighScore Plus *Fitting Si NIST 640e Standard Data from NOMAD in Fullprof*  
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**Part 2 Multiphase or Dual Phase**

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## **Rietveld Refinement of BiCaFeMnO(R3c+Pbnm) using FullProf Software**

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How to do Reitveld Refinement using  
X'pert Highscore Plus | Electroceramics  
LabOR61: Optimising the Teaching of OR  
– Yesterday and Tomorrow

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### Guide Of Fullprof Fullprof Manual

The manual begins with a short description of the way to obtain FullProf from the anonymous ftp-area. It follows with a chapter where a brief description of the purpose of the program and the list of input and output files. The second chapter is dedicated to the description of strategies for using the program and the description of examples.

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AN TOTHE PROGRAM... AN  
INTRODUCTION TO THE PROGRAM  
FullProf Manual Users' Guide of FullProf

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FullProf : Crystal and magnetic structure refinement, powder/single crystals, polarised neutrons, constant wavelength, TOF, energy dispersive, multiple patterns,

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simulated annealing for solving crystal and magnetic structures (integrated and profile intensities).

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Introduction to the programs of the  
FullProf Suite

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The crystal structure has been refined by the Rietveld method, by using the program fullprof. All the samples are tetragonal and were indexed to the space group P42/mnm.

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