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The FullProf Suite (for Windows, Linux and macOS) is formed by a set of crystallographic programs (FullProf, WinPLOT, EdPCR, GFourier, etc...) mainly developed for Rietveld analysis (structure profile refinement) of neutron (constant wavelength, time of flight,...

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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

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Summary of information for working with symmetry modes in FullProf. Below we provide a part of the file "fp2k.inf" distributed with the FullProf Suite, in which all the modifications to the FullProf manual are provided. This part concerns the modifications of the PCR file introduced by the use of symmetry modes.

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Each isolated atom has the same definition of parameters as described in the Fullprof manual (see lines 11.4.1 and 11.4.3 of the Fullprof guide). Current information about "free atoms" is printed to the screen when parameter P6 (see below) is set to any negative value.

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#FullProf is the most used software for #refinement X-ray diffraction (#XRD) results. FullProf is considered to be the most authenticate one. As it is very powerful software and it has lots of ...

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Fullprof Tutorial: How to do Profile fitting of XRD data of Tetragonal BaTiO₃

If this settings file doesn't exist, the following dimensions will be used by default: Max_Patterns = 15 ! number of patterns Max_Points = 15000 ! number of points. WinPLOTR uses a settings file in order to adapt the capabilities of the program to the different needs and wishes of the user.

Y2O3 Fullprof - Durham University

Here I am uploading a video created by me to find lattice constants, Indexing and profile fitting of powder diffraction data of unknown phase (e.g. Tetragonal BaTiO₃) using FullProf.

FullProf user's guide - CCP14

Fullprof Refinement - Y₂O₃. Files needed: y2o3.dat; y2o3.cif Learning Outcomes: This example shows a simple Rietveld refinement of Y₂O₃ in fullprof. You may have used the same dataset in topas/gsas.

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