

 winmostar tutorial

# Quantum ESPRESSO

## Fermi Surface

V11.6.5

5 April 2024

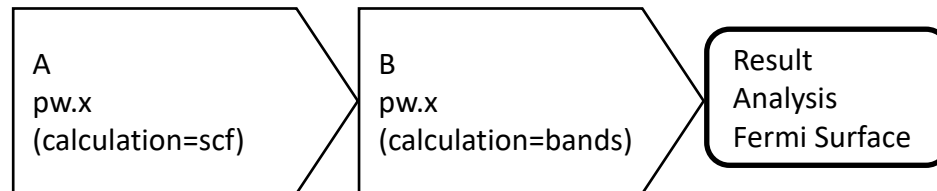
X-Ability Co., Ltd.

# About This Manual

- This manual is a tutorial demonstrating use cases for Winmostar V11.
- For those using Winmostar V11 for the first time, please consult [Beginner's Guide](#).
- For those who wish to explore the details of each feature, please refer to [Winmostar User Manual](#).
- Those who wish to practice the contents of this manual are encouraged to attend a training session.
  - [Winmostar Introductory Training Session](#): This guide only introduces the operation methods of the Basic Tutorial.
  - [Winmostar Basic Training Session](#): We will cover the theoretical background, explanations on interpreting results, operational methods of the Basic Tutorial, and procedures for some tutorials beyond the basic level.
  - [Individual Training Session](#): You can freely customize the training content according to your preferences.
- If you are unable to proceed with the operations as outlined in this manual, please first consult [Frequently asked questions](#).
- If your issue is not resolved through the Frequently Asked Questions, for the purpose of information accumulation and management, please contact us using [Contact page](#). Attach files generated at the time of the issue and provide steps to reproduce the problem.
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# Overview

- Conduct an SCF calculation for a Cu crystal, then calculate the electronic states at each k-point to display the Fermi surface. (These steps are run in sequence on Winmostar).



Note :

- The choice of k-points, number of bands, type of pseudopotential, cutoff energy, and smearing settings can affect the calculation results. In this tutorial, settings with reduced accuracy are used to obtain results more quickly.
- ◆ For a detailed explanation of Quantum ESPRESSO's calculation methods and settings, please see the following article from our company: [https://qiita.com/xa\\_member](https://qiita.com/xa_member)

# Preference of Operating Environment

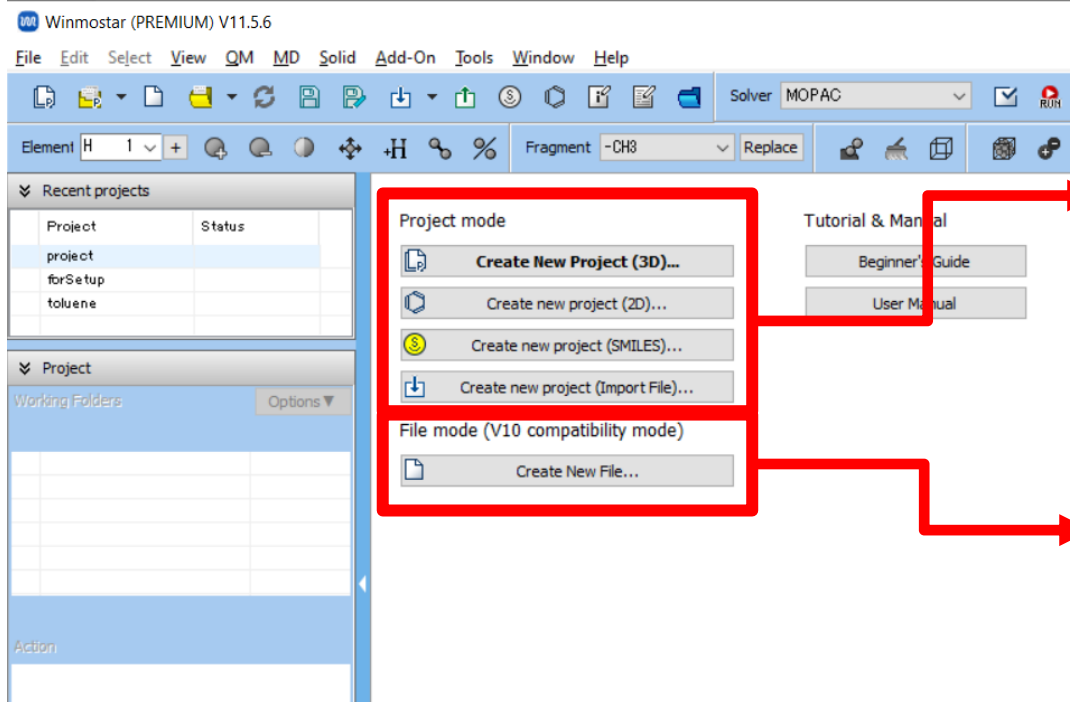
- For users of Winmostar version V11.5.0 or later on a 64-bit environment, please install and configure [CygwinWM version 2023/04/05 or later](#).
  - The CygwinWM version after 2023/04/05 includes the recommended version of 64-bit Quantum ESPRESSO.
- If the above does not apply to you or if you wish to use a version of Quantum ESPRESSO other than [the recommended one](#), you will need to install and configure [Windows version of Quantum ESPRESSO](#) separately.

# Operating Modes of Winmostar V11

V11 offers two operating modes: **Project Mode** and **File Mode**.

This manual focuses on operations in Project Mode.

For operations in File Mode, please refer to [Quantum ESPRESSO tutorial for version 10](#).



## Project Mode **New Features in V11**

Users can manage jobs without having to manage individual files.

We generally recommend using this mode.

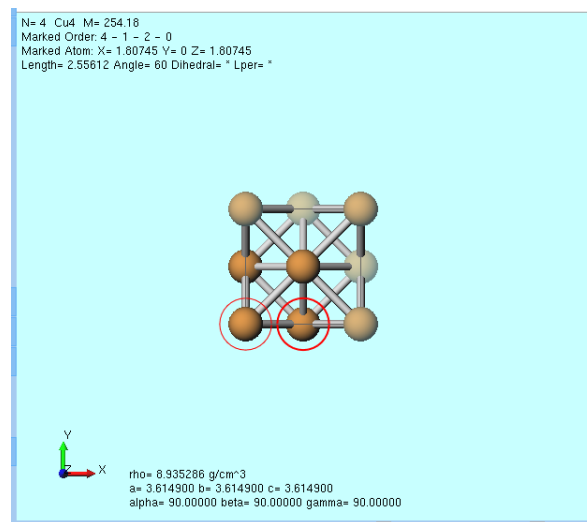
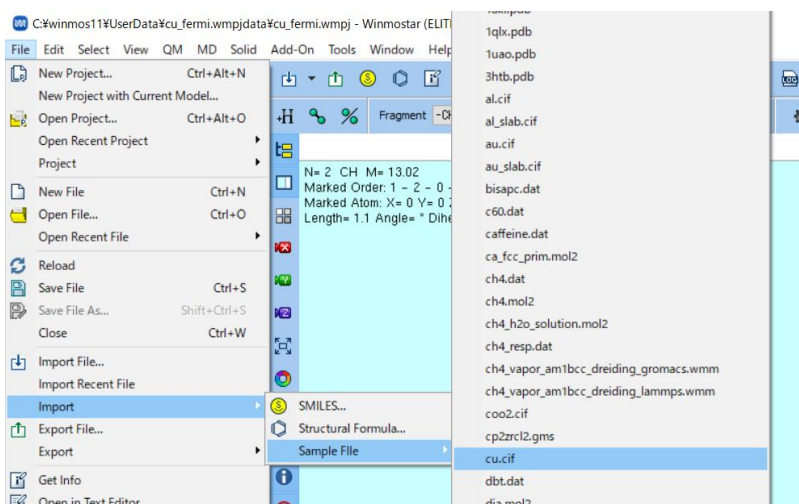
## File Mode

Users explicitly create and manage individual files.


The operational procedure is the same as from V10 and earlier versions.

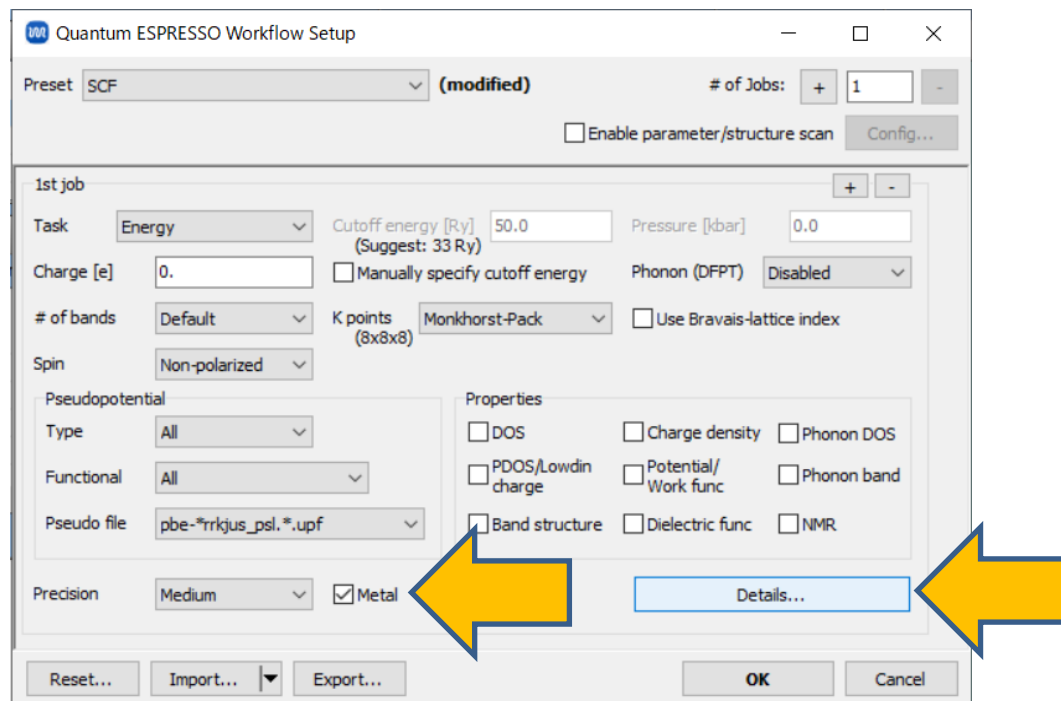
# A. Modeling of the System

- Please refer to [QE Basic Tutorial](#) for the basic operation method.
  - For detailed instructions on creating the initial structure, please refer to [Winmostar User Manual section 5, 'Structure Building'](#).
- A. Launch Winmostar and click on **Create New Project (3D)**. If Winmostar is already running, click **File | Close** first.
- B. Enter 'cu\_fermi' in **Project name** and click **Save**.
- C. Click **File | Import | Sample File | cu.cif**.
- If you wish to load a different file at this stage, use **File | Import File** instead.
- D. In **Import File** dialog, click **Discard and import**.



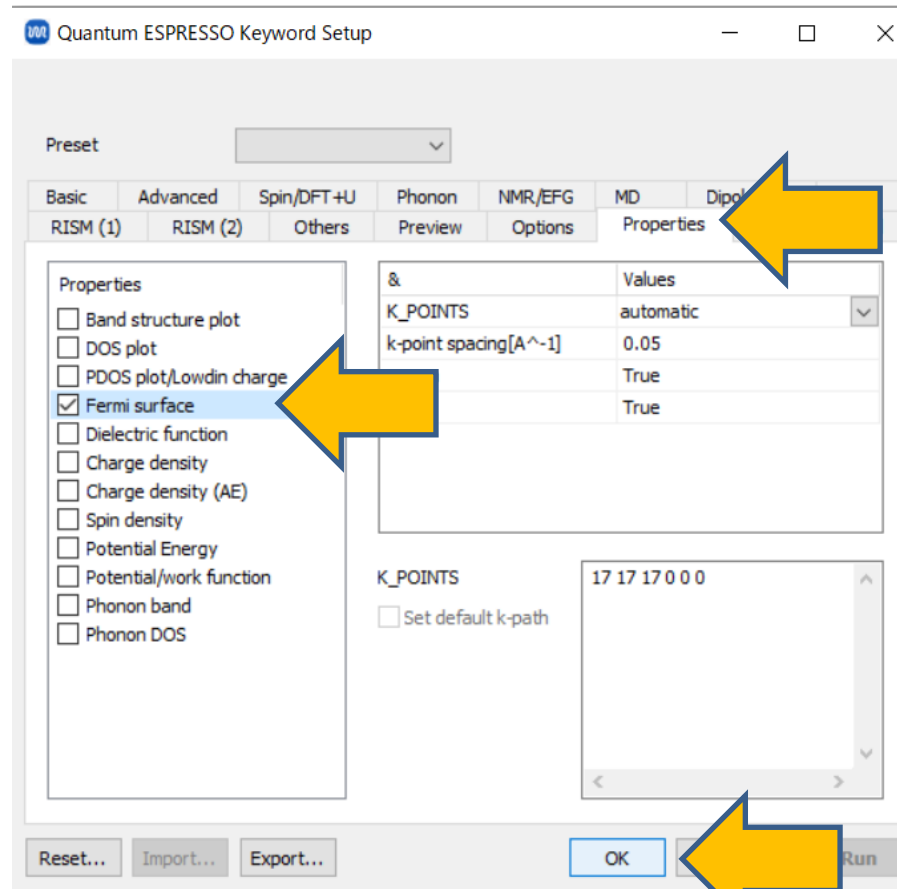
## B. Execution of Calculation

- Select **Quantum ESPRESSO** from **Solver** in Toolbar and click  (**Workflow Setup**).
- To reduce calculation time, click **Yes** when asked if you want to convert to a primitive cell. The converted structure will appear in Viewport. Click **OK** when 'Successfully converted lattice.' is displayed.
- In **Quantum ESPRESSO Workflow Setup** window, check **Metal** and click **Details**.



## B. Execution of Calculation

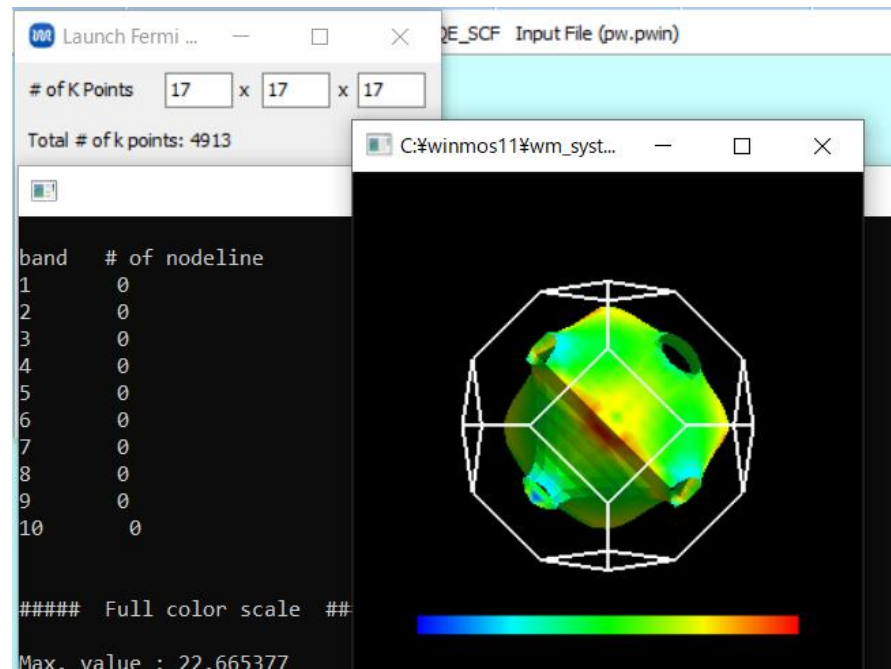
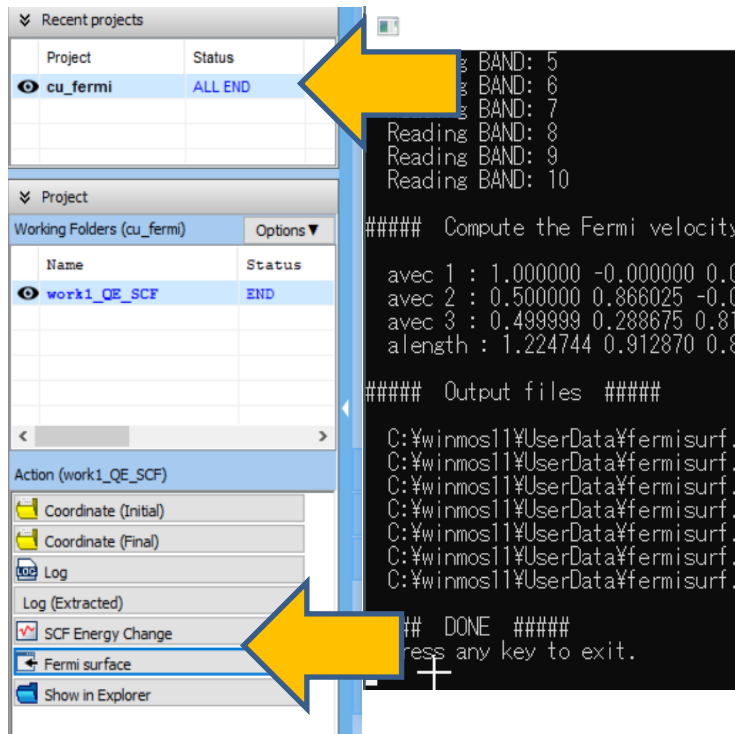
- A. In **Properties** tab, check **Fermi surface** under **Properties** and click **OK**.
- B. Click **OK** in **Quantum ESPRESSO Workflow Setup** window, then make appropriate settings in **Job Setting** window before clicking **Run**.





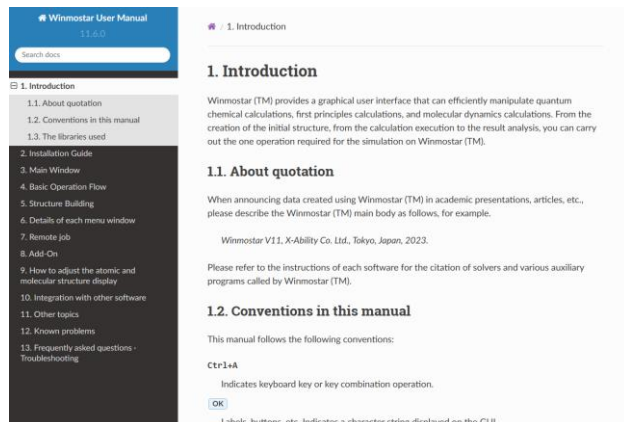
## C. Result Analysis

- After the status of **work1\_QE\_SCF** in **Working Folders** changes to **END (blue)**, click **work1\_QE\_SCF** in **Working Folders** and select **Fermi surface** in **Action**.
- A black window will appear with the message 'Press any key to exit.' Press Enter key on your keyboard.
- The Fermi surface will be displayed in a separate window. For operation instructions, please refer to [FermiSurfer website](#).



# Finally

- For detailed information on each feature, please refer to [Winmostar User Manual](#).



## [Winmostar User Manual](#)

## Scenes from [Winmostar Training Session](#)

- If you wish to practice the contents of this guide, please consider attending [Winmostar Introductory Training Session](#), [Winmostar Basic Training Session](#), or [Individual Training Session](#). (See page 2 for details.)
- If you are unable to proceed as instructed in this guide, please first consult [Frequently asked questions](#).
- If FAQs do not resolve your issue, for the purposes of information accumulation and management, please contact us through [Contact page](#), detailing the steps to reproduce the issue and attaching any generated files at that time.