

2025/02/26 [for version 5.0]

# How to setup MaterApps LIVE!

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MaterApps LIVE! Development Team





# MateriApps — a Portal Site for Materials Science Simulation

- Aiming at the community formation through the promotion of application



since May 2013

- Introducing **320+ materials science applications and tools**
- Finding applications
  - search tags: features, targets, calculation methods/algorithms
- Information of applications
  - brief introduction, link to official pages, information installation, usage, etc
- Information of hands-on sessions, software update, etc
- Glossary of keywords, Concierge, Reviews
- 17,000+ page views / month, 6,500+ unique visitors / month (FY2023)

# Applications on MateriApps

- Introducing **320+ materials science applications and tools**

## DFT

AkaiKKR★

OpenMX★

xTAPP★

ABINIT★

...

## Quantum Chemistry

FMO★

SMASH★

GAMESS★

DC★ ...

## Molecular Dynamics

MODYLAS★

Gromacs★

ERmod★

MDACP...

## Lattice Models

ALPS★

DSQSS

BLOCK

DMRG++

## Continuum Simulation

ANSYS Multiphysics

Octa ...

## Data Analysis

CLUPAN★

phonopy★

## Visualization

fu★

TAPIOCA★

Database, Integrated Environment  
Machine Learning, Quantum Computing

★ included in MateriApps LIVE!



# Current status in computational materials science

- From developers' viewpoint
  - New algorithms should be implemented and used. Or, it will be forgotten ever existed.
  - It cost much to write and update documents
  - Development of software itself is hardly considered as scientific achievements
- From users' viewpoint
  - What kind of applications? Who develop them?  
Which application should I use for my problem?
  - Manual and documentation are not well prepared.
  - How to evaluate the accuracy of results?
- Goal of MateriApps project
  - **Forming of community in the field of computational materials science through the promotion of open source software**



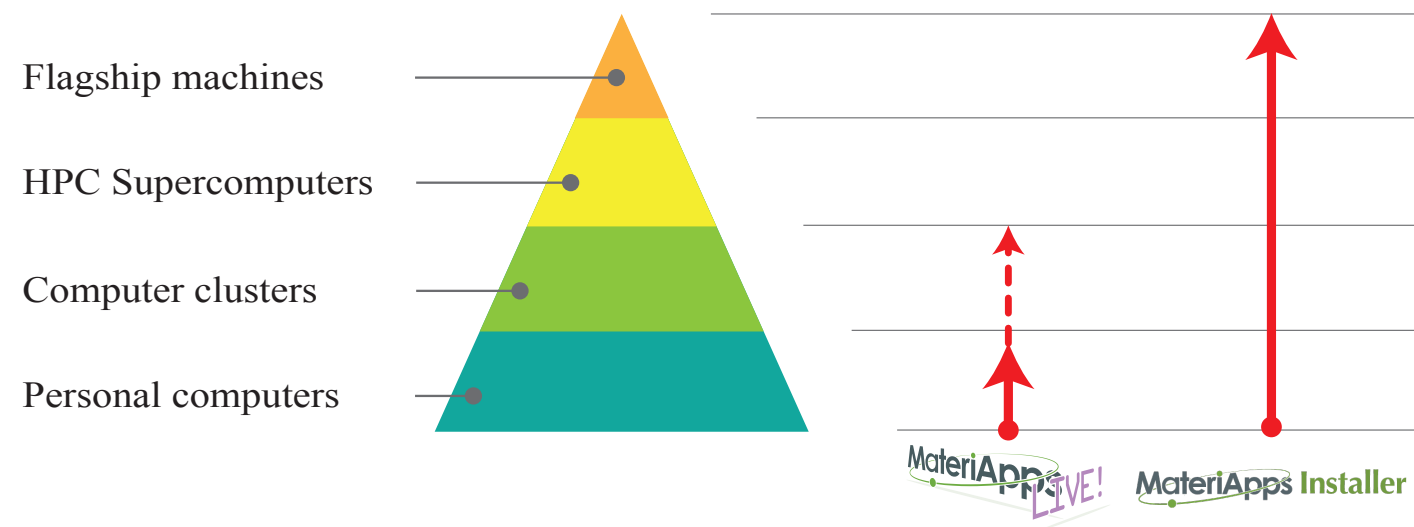
開発者



利用者

# What MateriApps will provide

- To find and learn application software
  - catalog of application/tool on **MateriApps web**
- To start using application software
  - **MateriApps LIVE!**
- To active use application software
  - pre-installation to Fugaku, supercomputers, etc: **MateriApps Installer**

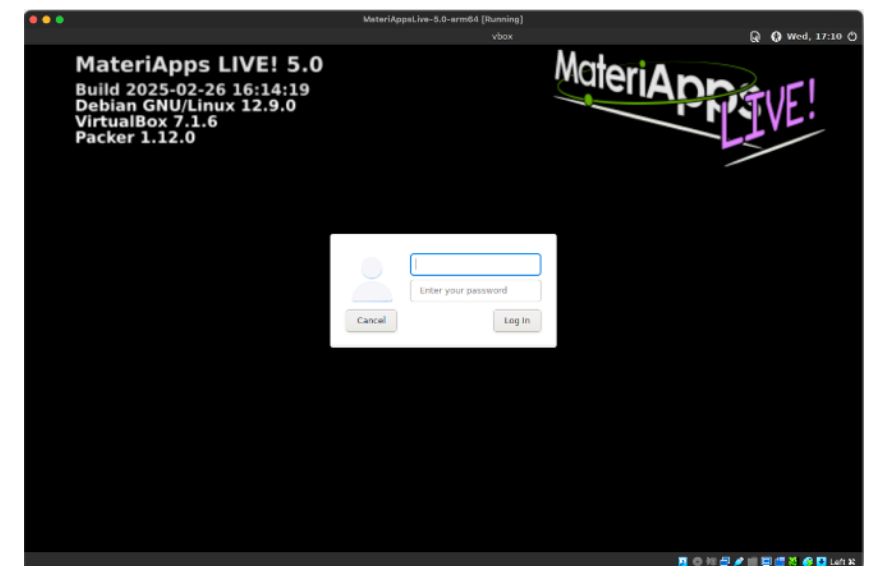


- Infrastructure for easily starting materials science simulations for theoreticians, experimentalists, researchers in companies, students, and more...

# What's MateriApps LIVE!



- Debian Linux running on a virtual machine (VirtualBox/Docker)
  - works on Windows and Mac (Intel/Apple Silicon)
  - just boot and get ready for materials science simulations without installation
- Version 5.0 was released in February 2025
- Pre-installed applications and tools
  - abinit, AkaiKKR, ALAMODE, ALPS, CONQUEST, Feram, DCore, DSQSS, HΦ, LAMMPS, mVMC, OCTA, OpenMX, Quantum ESPRESSO, PHYSBO, SMASH, TeNeS, xTAPP, etc
  - OVITO, ParaView, Tapioca, VESTA, XCrysDen...
  - GUI installer for CASINO, GAMESS, and VMD
- Available from MateriApps LIVE! webpage
  - c.a. 15,000+ copies distributed since July, 2013





# MateriApps LIVE! is useful for ...

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- Hands-on sessions using MateriApps LIVE!
  - MateriApps LIVE! Tutorials
  - H $\Phi$ , xTAPP, ALPS, DCore, mVMC, ALAMODE, DDMRG, DSQSS, SALMON, CASINO, TeNeS, etc.
- Practices in lectures
  - Computational Physics
  - Computer Experiments (UNIX + C, LaTeX, VCS)
- Used by experimentalists, researchers in private companies
- Used by researchers in the field of computer science
- Easy setup (c.a. 15min) without no troubles
- Useful for operation check, trouble shooting, user support

# Materials Science Simulation by MateriApps LIVE!

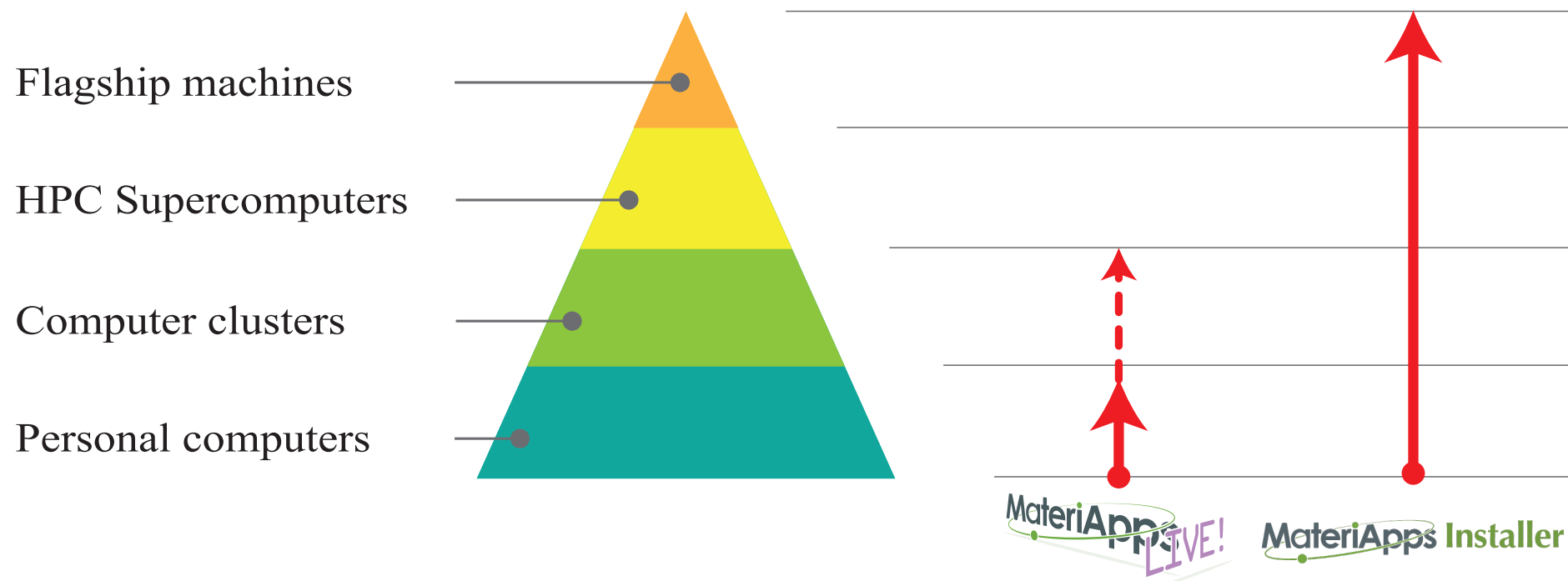
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- Introduction / Setup
- First-principles band calculation (OpenMX / Quantum ESPRESSO / xTAPP)
- Simulation of solution by molecular dynamics (LAMMPS / Gromacs)
- Lattice model simulation (ALPS / H $\Phi$  / mVMC)
- Quantum chemistry calculation (in preparation)
- Hands-on materials are available at <https://github.com/cmsi/MateriAppsLive/wiki/MaLiveTutorial> (currently only in Japanese)



# Wanna larger-scale simulations?

- For Debian/Ubuntu Linux workstations
  - Debian package for MateriApps LIVE! can be used
  - <https://github.com/cmsi/MateriAppsLive/wiki/UsingMateriAppsInDebian>
  - (Can also be installed on Google Colab)
- For PC clusters, clouds, supercomputers at ISSP and IT centers, Fugaku, etc
  - MateriApps Installer <https://ma.issp.u-tokyo.ac.jp/app/268>



# VirtualBox edition vs. Docker edition

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- VirtualBox edition
  - Pros
    - Many visualization tools (Ovito, Paraview, Vesta, etc.) run on the virtual machine
    - Note: Vesta only works on Windows and macOS Intel
  - Cons
    - Uses much memory
- Docker edition
  - Pros
    - The startup is faster after the second time
    - Uses less memory
  - Cons
    - Some visualization tools do not work



# MateriApps LIVE! (VirtualBox edition)

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- Files
  - setup-en.pdf (English), setup.pdf (Japanese)
    - this document
    - <https://speakerdeck.com/wistaria/how-to-setup-materiapps-live>
  - README-en.html, README.html
    - <https://github.com/cmsi/MateriAppsLive/wiki/MateriAppsLive-ova-en>
  - VirtualBox 7.1 Installer: VirtualBox-\*-Win.exe (Windows), VirtualBox-\*-OSX.dmg (macOS Intel), VirtualBox-\*-OSXArm64.dmg (macOS Apple Silicon)
    - available at <https://www.virtualbox.org/wiki/Downloads>
  - MateriApps LIVE! VirtualBox Diskimage: MateriAppsLive-\*-amd64.ova (Windows / macOS Intel), MateriAppsLive-\*-arm64.ova (macOS Apple Silicon)
    - available at <https://github.com/cmsi/MateriAppsLive/wiki/download>

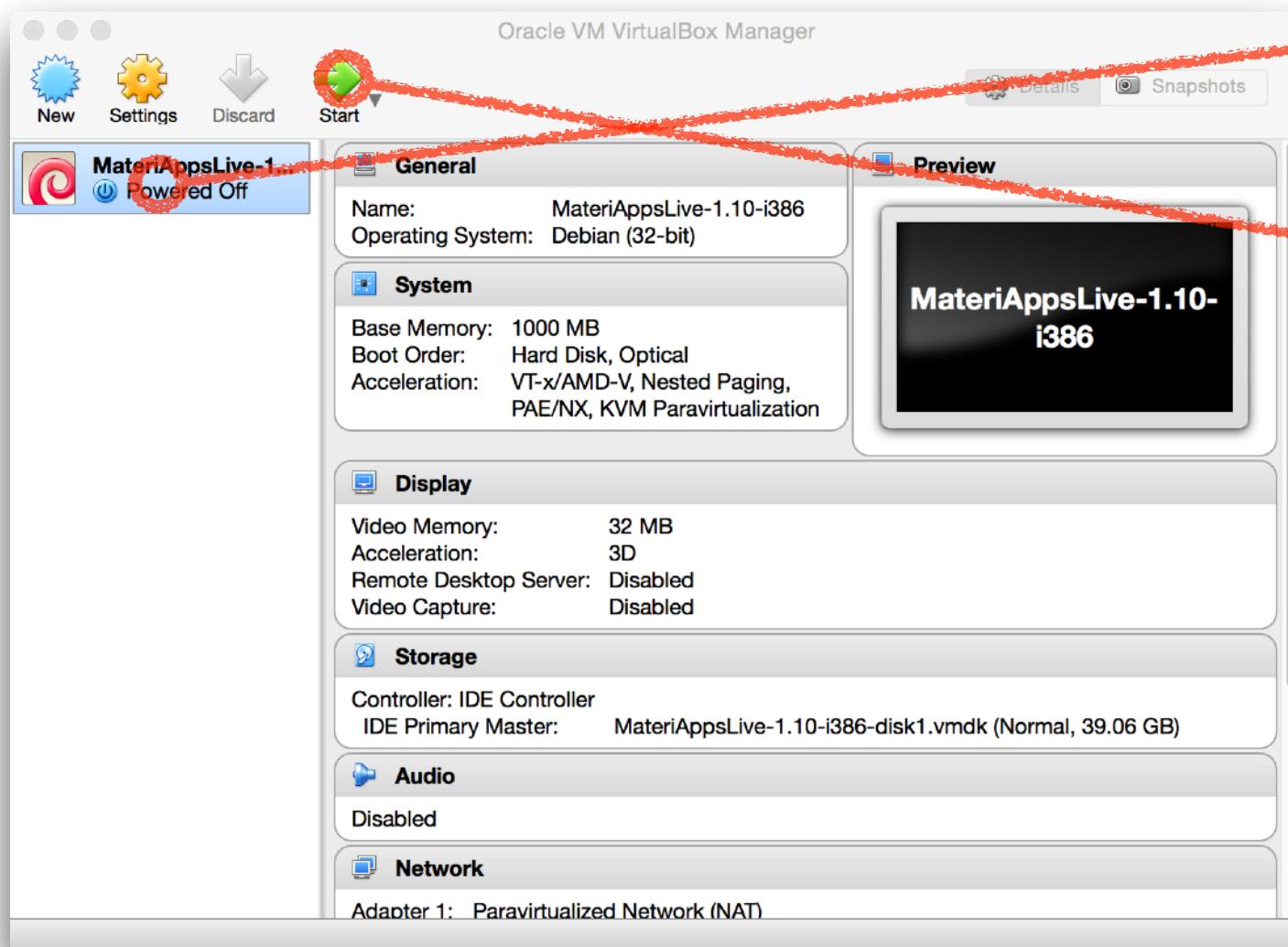
# Let's get started (VirtualBox edition)

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- ✓ Download distribution files
- ✓ Install VirtualBox by double-clicking the installer
  - For Windows: VirtualBox-\*-Win.exe
  - For macOS: VirtualBox-\*-OSX.dmg (Intel), VirtualBox-\*-OSXArm64.dmg (Apple Silicon)
- ✓ Import MateriApps LIVE!
  - double-click MateriAppsLive-\*.ova
  - VirtualBox will start automatically and import window will open. Then press “import” button
  - VirtualBox Manager window will appear in two or three minutes
- Host (host OS): operating system (Windows, Mac OS X, etc) on which VirtualBox is running
- Virtual machine (guest OS): operating system (= MateriApps LIVE!) running on VirtualBox




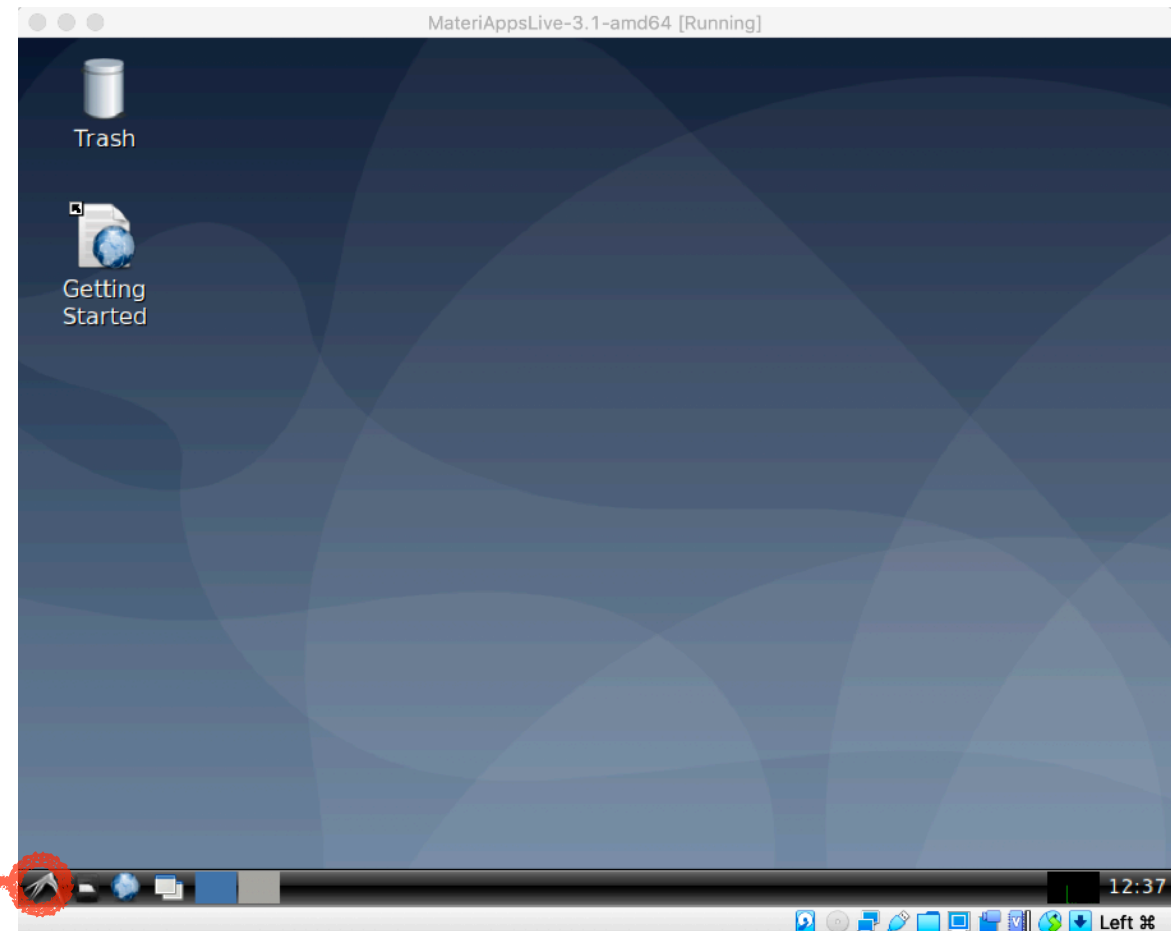
# Booting virtual machine (VirtualBox edition)



1. Choose “MateriAppsLive...”
2. Press “Start” button.
3. Wait until login window will appear.

# Login to MateriApps LIVE! (VirtualBox edition)

- Login by using
  - User name (login): *user*
  - Password: *live*
- Desktop (right) will appear
- *start menu* 
- How to open a terminal window  
start menu ⇒ “System Tools” ⇒ “LXTerminal”
- How to shutdown the virtual machine  
start menu ⇒ “Logout” ⇒ “Shutdown”





# Keyboard, resolution setting (VirtualBox edition)

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## ✓ Using Japanese keyboard

- start menu ⇒ “System Tools” ⇒ “Switch to Japanese Keyboard Layout”
- check if “@” key works correctly

## ✓ Changing resolution and scale of the virtual machine window

- Resolution (number of pixels): “800 x 600” by default
- Scale (pixel density): “200%” by default
- The resolution and scale can be changed from the “View” menu ⇒ “Virtual screen 1” of the host OS
- When using visualization software such as VESTA, OVITO, ParaView, set the resolution to “1024 x 768” or higher
- If characters are small and difficult to see, increase the scale

# File sharing, copy & paste (VirtualBox edition)

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- ✓ File sharing between host OS and virtual machine
  - shutdown the virtual machine, if it is running
  - choose MateriAppsLive-\* in VirtualBox Manager window, then “Settings”
  - open “Shared Folders” tab and click “+” on the right
  - click “v” on the right of “Folder Path”, choose “Other...”, and select the folder to be shared
  - check “Auto-mount” box and press “OK”. Then press “OK” again
  - the folder specified above can be accessed as /media/sf\_... after restarting the virtual machine
- ✓ Copy & Paste: How to paste strings copied from a PDF file on host OS?
  - right click on terminal window ⇒ “Paste”
  - or press “V” with “shift” and “control” keys
  - right click ⇒ “Copy”, or “shift + control + C” to copy a string



# MateriApps LIVE! (Docker edition)

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- Files
  - setup-en.pdf (English), setup.pdf (Japanese)
    - this document
    - <https://speakerdeck.com/wistaria/how-to-setup-materiapps-live>
  - README-en.html, README.html
    - <https://github.com/cmsi/MateriAppsLive/wiki/MateriAppsLive-docker-en>
  - Docker Desktop installer (for macOS)
    - available at <https://www.docker.com/>
  - XQuartz installer (for macOS)
    - available at <https://www.xquartz.org/>
  - MateriApps LIVE! Docker script
    - <https://malive.s3.amazonaws.com/malive.sh>

# Let's get started (macOS Docker edition)

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- ✓ Download and install Docker Desktop
- ✓ Download, install, and setup XQuartz
- ✓ Set up the shared folder
- ✓ Download and execute MateriApps LIVE! Docker script
  - `curl -L -O https://sf.net/projects/materiappslive/files/docker/malive.sh`
  - `sh malive.sh`
- Ref. <https://github.com/cmsi/MateriAppsLive/wiki/GettingStartedDocker-en>



# Let's get started (Windows Docker edition)

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- ✓ Install WSL2 (Windows Subsystem for Linux 2)
- ✓ Set up the Docker repository and execute apt-get
- ✓ Set up the shared folder
- ✓ Download and execute MateriApps LIVE! Docker script
  - `curl -L -O https://sf.net/projects/materiappslive/files/docker/malive.sh`
  - `sh malive.sh`
- Ref. <https://github.com/cmsi/MateriAppsLive/wiki/GettingStartedDocker-en>

# MateriApps planning & production

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- Administration:
  - Center for Computational Materials Science, Institute for Solid State Physics, University of Tokyo (ISSP-CCMS)
- MateriApps Development Team
  - Kota Ido (ISSP), Shusuke Kasamatsu (Dept. of Phys., Yamagata Univ.), Takeo Kato (ISSP), Naoki Kawashima (ISSP), Hikaru Kouta (ISSP), Syngge Todo (Dept. of Phys., Univ. of Tokyo/ISSP), Masahito Fukuda (ISSP), Kanako Yoshizawa (RIST), Jun Yamazaki (ISSP)
  - (contract) Academeia (Yusuke Konishi, Masashi Noda, Gotai Yamada)
- Cooperation:
  - Research Organization for Information Science and Technology (RIST)
  - Professional development Consortium for Computational Materials Scientists (PCoMS)
  - Data creation and utilization-type MaTerial R&D project (DxMT)