

# Program

(JST = UTC + 9)

## 07 December 2022 (Wed)

### Introduction of Nuclear DFT

(Chair: Tomoya Naito)

13:20–13:30 Tomoya Naito (RIKEN iTHEMS/U. Tokyo)

[Opening](#)

13:30–15:00 Nobuo Hinohara (U. Tsukuba)

[Recent progress in nuclear DFT](#)

## 08 December 2022 (Thu)

### Time-Dependent Density Functional Theory I

(Chair: Shunsuke A. Sato)

10:30–12:00 Kazuyuki Sekizawa (Tokyo Tech)

[Nuclear TDDFT: Past, Present, and Future](#)

### Time-Dependent Density Functional Theory II

(Chair: Kazuyuki Sekizawa)

13:30–15:00 Takashi Nakatsukasa (U. Tsukuba)

[Requantizing the time-dependent density functional dynamics](#)

## 09 December 2022 (Fri)

### Many-Body Theoretical Methods I

(Chair: Osamu Sugino)

10:30–12:00 Gianluca Colò (U. Milan/INFN Milano/YITP, Kyoto U.)

[Many-body correlations and EDF-based nuclear models](#)

### Many-Body Theoretical Methods II

(Chair: Takashi Nakatsukasa)

13:30–15:00 Osamu Sugino (Institute for Solid State Physics, U. Tokyo)

[Many-body Green's Function and Density Functional Theory for Condensed Matters](#)

## 12 December 2022 (Mon)

### One-Day Workshop (15 min. talk & 5 min. discussion)

(Chair: Kenichi Yoshida)

- 10:20–10:40 **Qiang Zhao** (Center for Exotic Nuclear Studies, Institute of Basic Science)  
 (Online) Covariant Density Functional Theory with Localized Exchange Terms
- 10:40–11:00 **Hiroyuki Tajima** (U. Tokyo)  
 (Online) Towards universal energy density functionals of non-relativistic spin-1/2 Fermi gases
- 11:00–11:20 **Yusuke Tanimura** (Tohoku U.)  
 Visualization of nuclear many-body correlations with the most probable nucleon arrangement
- 11:20–11:40 **Mitsuaki Kawamura** (Institute for Solid State Physics, U. Tokyo)  
 Benchmark and application of density functional theory for superconductors
- 11:40–12:00 **Kai Wen** (U. Tsukuba)  
 Collective coordinates and collective inertial masses for nuclear reactions

### One-Day Workshop (15 min. talk & 5 min. discussion)

(Chair: Gianluca Colò)

- 13:30–13:50 **Kenta Yoshimura** (Tokyo Tech)  
 Time-Dependent Band Theory for the Inner Crust of Neutron Stars: Extension to Include Superfluidity
- 13:50–14:10 **Tomohiro Oishi** (Yukawa Institute for Theoretical Physics, Kyoto U.)  
 Magnetic excitations based on the relativistic energy-density functional theory
- 14:10–14:30 **Tomohito Amano** (U. Tokyo)  
 First-principles study of lattice dielectric function in rutile TiO<sub>2</sub>
- 14:30–14:50 **Moemi Matsumoto** (Tohoku U.)  
 Generator coordinate method with variational basis generation
- 14:50–15:10 **Akira Ohnishi** (Yukawa Institute for Theoretical Physics, Kyoto U.)  
 Density dependence of  $\Lambda$  potential in nuclear matter from heavy-ion collisions and hypernuclear spectroscopy

### One-Day Workshop (15 min. talk & 5 min. discussion)

(Chair: Takeru Yokota)

- 15:40–16:00 **Hiroyuki Sagawa** (U. Aizu/RIKEN RNC)  
 Beyond mean field approach for nuclear excitations and decays
- 16:00–16:20 **Shunsuke Yamada** (National Institutes for Quantum Science and Technology)  
 Ab initio TDDFT study of high-harmonic generation from dielectric thin films
- 16:20–16:40 **Kenichi Yoshida** (Kyoto U.)  
 Linear-response TDDFT for rotating nuclei
- 16:40–17:00 **Kohei Washiyama** (U. Tsukuba)  
 Large-amplitude collective dynamics with collective Hamiltonian derived from nuclear DFT + local QRPA
- 17:00–17:20 **William Dawson** (RIKEN R-CCS)  
 Density functional theory calculations of large systems:  
 Interplay between fragments, observables, and computational complexity
- 17:20–17:40 **Kotaro Uzawa** (Kyoto U.)  
 Microscopic description of induced fission based on the generator coordinate method

## 13 December 2022 (Tue)

### Isospin Symmetry Breaking and Its Symmetry Restoration (Chair: Ryosuke Akashi)

10:30–12:00 Koichi Sato (Kochi U.)  
Nuclear density functional calculations with proton-neutron mixing

### Orbital-Free DFT and Machine Learning (Chair: Masahiro Fukuda)

13:30–15:00 Fumihiko Imoto (Schrödinger)  
Recent Advances in Orbital-Free Density Functional Theory and Machine Learning Approaches to Develop Kinetic Energy Functionals

## 14 December 2022 (Wed)

### Relativistic DFT (Chair: Koichi Sato)

10:30–12:00 Yasuhiro Ikabata (Toyohashi Tech)  
Picture-change corrected density functional theory for relativistic quantum chemical calculations

## 15 December 2022 (Thu)

### Construction of Energy Density Functional (Chair: Fumihiko Imoto)

10:30–12:00 Tomoya Naito (RIKEN iTHEMS/U. Tokyo)  
Towards First-Principles Energy Density Functionals

### Time-Dependent Density Functional Theory III (Chair: Nobuo Hinohara)

13:30–15:00 Shunsuke A. Sato (U. Tsukuba)  
(Online) Application of real-time TDDFT simulation to nonlinear and nonequilibrium electron dynamics in matter

### Discussion (Chair: Gianluca Colò)

15:30–17:00 Discussion (Novel ideas to build EDFs & Collective coordinates and/or full CI)

## 16 December 2022 (Fri)

### Calculation Method of Local Physical Quantities (Chair: Takeru Yokota)

10:30–12:00 Masahiro Fukuda (Institute for Solid State Physics, U. Tokyo)  
Local physical quantities based on quantum electrodynamics

### Superconducting Density Functional Theory (Chair: Gianluca Colò)

13:30–15:00 Ryosuke Akashi (National Institutes for Quantum Science and Technology)  
Development of density functional theory for superconductors: recent progress

## 19 December 2022 (Mon)

---

### Machine Learning Approach

(Chair: Kenichi Yoshida)

---

- 10:30–12:00 Haozhao Liang (U. Tokyo)  
[Quantum computing for nuclear structure properties?](#)
- 12:00–12:10 Kenichi Yoshida (Kyoto U.)  
[Closing](#)

## 20 December 2022 (Tue)

---

### AM & PM Free Discussion

---