10/16 (Tue) 10:40-11:00 [1 minute x 19]

	Presenter	Affiliation	Title of poster presentation
1	Kohji Yoshikawa	University of Tsukuba	Combined effects of SIDM and stellar feedback on the structure of dwarf halos
2	Ataru Taniakwa	The University of Tokyo	Numerical study of thermonuclear explosions of white dwarf stars
3	Naoya Ukita	University of Tsukuba	2+1 Flavor Lattice QCD with the Physical Quark Masses
4	Jun Terasaki	Institute of experimental and applied physics	Examination and improvement of nuclear matrix elements of double-β decay in QRPA approach
5	Yasutaka Taniguchi	National Institute of Technology, Kagawa College	Alpha-cluster correlations in a ground state of 48Ti
6	Daisuke Takahashi	University of Tsukuba	Implementation of Parallel FFTs on Cluster of Intel Xeon Phi Processors
7	Yu-ichiro Matsushita	Tokyo Institute of Technology	Possible interface/near-interface states at SiC/SiO2
8	Norikazu Yamada	KEK	Strong CP problem and axion on the lattice
9	Eigo Shintani	RIKEN Center for Computational Science	Nucleon form factors on a (10.8 fm)^4 lattice at the physical point in 2+1 flavor QCD
10	Issaku KANAMORI	Hiroshima University	Lattice QCD code with Bridge++ for AVX-512 instruction sets
11	Eigo Shintani	RIKEN Center for Computational Science	Precision test of the standard model and search for the new physics in lattice QCD
12	Takashi Kaneko	KEK	Test of new physics models through B meson semileptonic decays
13	Miwako Tsuji	RIKEN Center for Computational Science	A performance projection method using simple benchmarks for real applications
14	Makito Abe	University of Tsukuba	Structure Formation in the Early Universe using Radiation Hydrodynamic Simulations
15	Kazuyuki Kanaya	University of Tsukuba	Theromodynamic quantities in (2+1)-flavor QCD using gradient flow
16	Shimpei Saito	University of Tsukuba	Lattice Boltzmann modeling and simulation of liquid jet breakup
17	Atsushi Yamada	University of Tsukuba	Electron Dynamics Simulation Study of Interactions between Pulse Light and Matter using SALMON software
18	Rintaro Fujikawa	University of Tsukuba	Molecular dynamics simulation for the clarification of the molecular diffusion behavior in the CO2 hydrate
19	Mitsuo Shoji	University of Tsukuba	Large scale molecular simulations for the elucidation of bimolecular structural changes and the catalytic reaction mechanisms