



MLQC4Dyn - Week 3

September 19

09:30 - 10:00	Onsite registration - coffee
10:00 - 10:10	Welcome from the organizers Chair: Bruno Senjean
10:10 - 11:00	Michael Marthaler - HQS Quantum Simulations <i>Simulating open quantum systems with quantum computers</i>
11:00 - 11:50	Ada Warren - Virginia Tech <i>ADAPTive Variational Quantum Algorithms for Gibbs State Preparation</i>
12:00 - 14:00	Lunch break Chair: Florian Eich
14:00 - 14:50	David Munoz Ramo - Quantinuum <i>Using molecular symmetries to perform quantum simulations on current quantum hardware</i>
14:50 - 15:10	Coffee break
15:10 - 16:00	Victor Alessandrini - Maison de la Simulation <i>Quantum Computing Emulation</i>
16:00 - 18:00	Group discussions / Round tables

September 20

	Chair: Francesco Tacchino
09:50 - 10:40	Matthias Degroote - Boehringer Ingelheim B.V. <i>Efficient quantum computation of molecular forces and other energy gradients</i>
10:40 - 11:00	Coffee break

11:00 - 11:50	Benjamin Lasorne - CNRS Montpellier <i>Quantum dynamics on quantum computers: where do we stand, where may we go?</i>
12:00 - 14:00	Lunch break Chair: Dario Rocca
14:00 - 14:50	Ivan Rungger - National Physical Laboratory <i>Green's functions algorithms on quantum computers for materials science</i>
14:50 - 15:10	Coffee break
15:10 - 16:00	Giulia Galli (R) - The University of Chicago and Argonne National Laboratory <i>Quantum embedding methods for simulations of materials on quantum computers</i>
16:00 - 18:00	Group discussions / Round tables
19:30 - 21:30	Social dinner

September 21

	Chair: Thomas Ayril
09:00 - 09:50	Ivano Tavernelli - IBM Research Zurich <i>TBA</i>
09:50 - 10:40	Francesco Tacchino - IBM Research Zurich <i>Quantum algorithms for dynamics: from quantum simulations to quantum machine learning</i>
10:40 - 11:00	Coffee break
11:00 - 11:50	Rick van Bijnen - Innsbruck University / IQOQI <i>Variational Quantum Optimization Algorithms</i>
12:00 - 14:00	Lunch break Chair: Martin Beseda
14:00 - 14:50	Artur Izmaylov - University of Toronto <i>Algebraic techniques for dealing with difficult operators in quantum chemistry on a quantum computer</i>
14:50 - 15:10	Coffee break

15:10 - 16:00

Saad Yalouz - CNRS and Université de Strasbourg

Quantum computing to describe properties of correlated many-body systems: from quantum chemistry to condensed matter physics

16:00 - 16:50

Roman Krems (R) - University of British Columbia

Optimal quantum kernels for classification and regression problems

September 22

Chair: Matthieu Saubanère

09:50 - 10:40

Francesco Evangelista - Emory University

Projective quantum algorithms for quantum chemistry

10:40 - 11:00

Coffee break

11:00 - 11:50

Bruno Senjean - CNRS Montpellier

Towards Density Functional Theory on Quantum Computers

12:00 - 14:00

Lunch break

Chair: Saad Yalouz

14:00 - 14:50

Martin Beseda - CNRS Montpellier

Quantum dynamics with state-averaged VQE in noisy environment

14:50 - 15:40

Matthieu Saubanère - CNRS Montpellier

Alternatives to unitary coupled cluster Ansatz for variational quantum Eigensolver

16:00 - 19:00

Happy hour

September 23

Chair: Ivan Rungger

09:00 - 09:50

Giuseppe Carleo - EPFL

Variational learning in the age of machine learning and quantum computing

09:50 - 10:10

Coffee break

10:10 - 11:00

Thomas Ayril - Atos Quantum Lab

Quantifying quantum advantage through high-performance classical simulations: two examples

11:00 - 11:50

Adam Smith (R) - The University of Nottingham

Topology on quantum computers

12:00 - 14:00

Lunch

(R) denotes remote speakers