

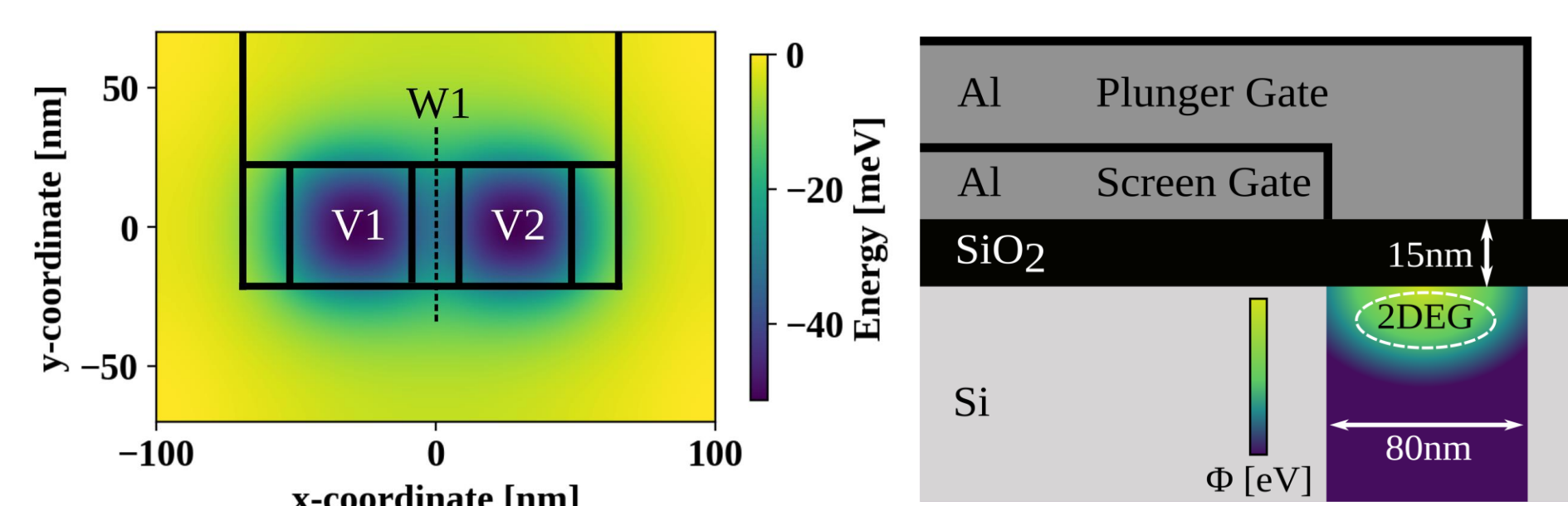
Simulated Control of Spin Qubits in MOSFET Quantum Dot Linear Arrays



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Introduction

We present a comprehensive simulator of electron spin qubits in electrostatically-defined quantum dots (QDs) to address challenges in designing quantum information processors.



Finite element solutions to Poisson's equation of realistic Silicon MOS are leveraged:

- Determine charge stability regions for various voltage configurations
- Engineer voltage pulses for spin qubit control
- Simulate gate operations on spin qubits in quantum circuits

Hubbard Hamiltonian

Hubbard parameters calculated from electrostatic potential [1]:

- Chemical potential
- Tunnel coupling
- Coulomb repulsion: onsite & interdot

$$H_{Hubbard} = H_\mu + H_t + H_U + H_u$$

$$H_\mu = - \sum_{\sigma=\uparrow,\downarrow} \sum_{i=1}^N \mu_i \hat{n}_{i,\sigma}$$

$$H_t = - \sum_{\sigma=\uparrow,\downarrow} \sum_{i<j} t_{ij} (\hat{c}_{i,\sigma}^\dagger \hat{c}_{j,\sigma} + h.c.)$$

$$H_U = \sum_{i=1}^N U_{ii} \hat{n}_{i,\uparrow} \hat{n}_{i,\downarrow}$$

$$H_u = \sum_{\sigma_1, \sigma_2 \in \{\uparrow, \downarrow\}} \sum_{i<j} U_{ij} \hat{n}_{i,\sigma_1} \hat{n}_{j,\sigma_2}$$

$$\left[-\frac{\hbar^2}{2m} \nabla_{2D}^2 + V_i(\vec{r}) \right] \psi_i(\vec{r}) = \epsilon_i \psi_i(\vec{r}) \quad \text{2D Schrödinger}$$

Effective Spin Hamiltonian

Spin Hamiltonian in rotating frame [2]:

- Stark shift: g-factor deviation
- Exchange interaction: Heitler-London/Hund-Mulliken approximation

$$H_{Spin} = H_Z + H_J$$

$$H_Z = \hbar \sum_{j=1}^N \frac{1}{2} \left[\left(1 + \frac{\delta g_j(t)}{2} \right) \omega - \omega_{RF} \right] Z_j$$

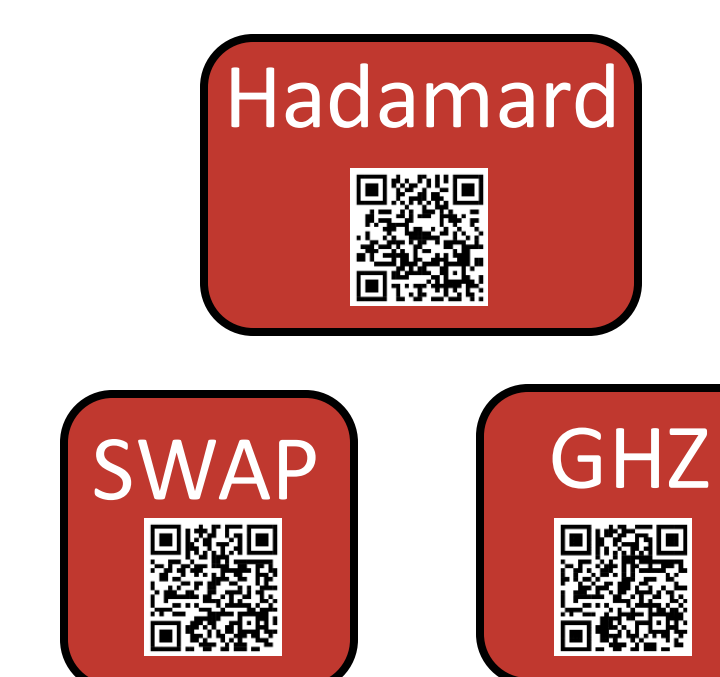
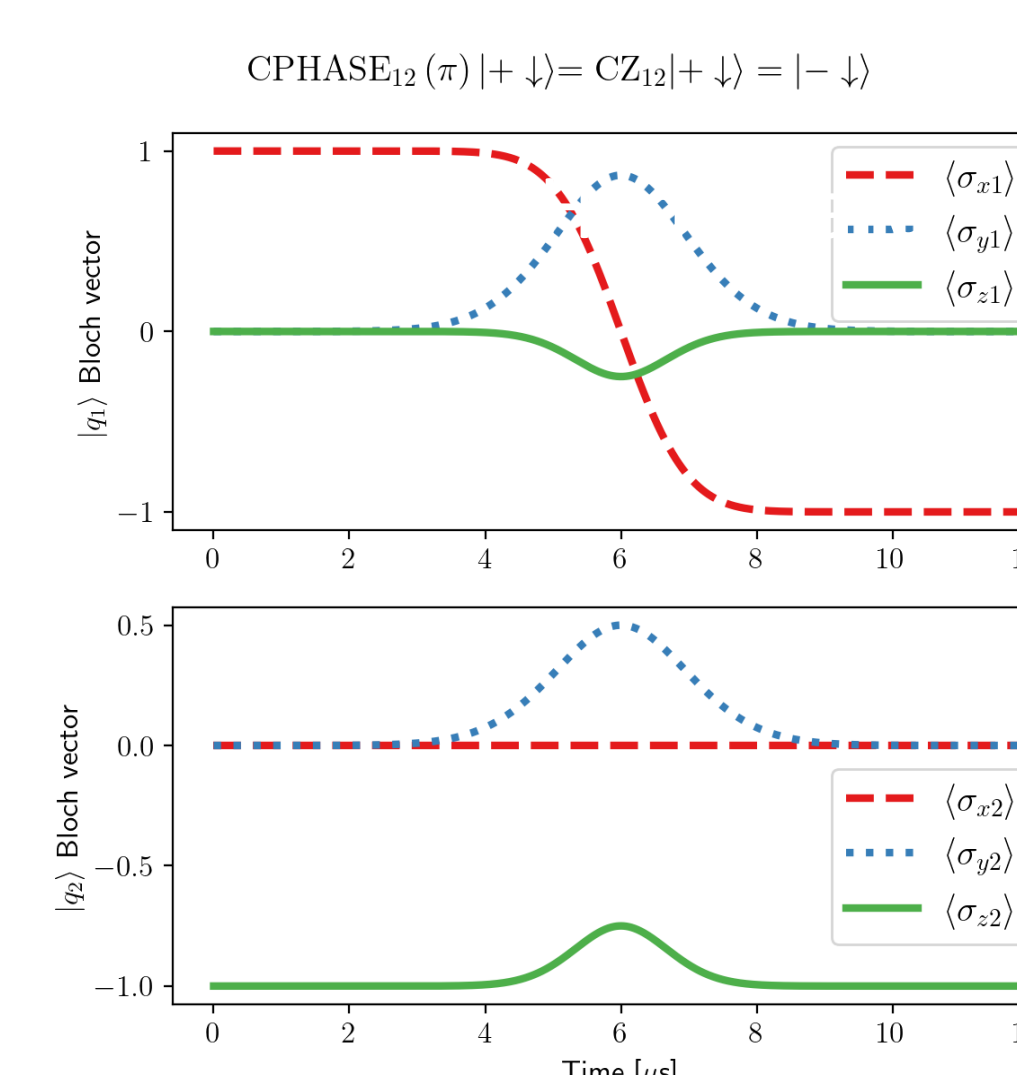
$$+ \frac{\Omega(t) \hbar}{2} (\cos \phi(t) X_j + \sin \phi(t) Y_j)$$

$$\delta g_j(t) = \eta \langle \psi_i(t) | \mathcal{E}_z(t)^2 | \psi_i(t) \rangle$$

$$H_J = \sum_{i<j} \frac{J_{ij}(t)}{4} \vec{\sigma}_i \cdot \vec{\sigma}_j$$

A novel, custom control method maps experimental control pulses from the expected effective parameter behavior:

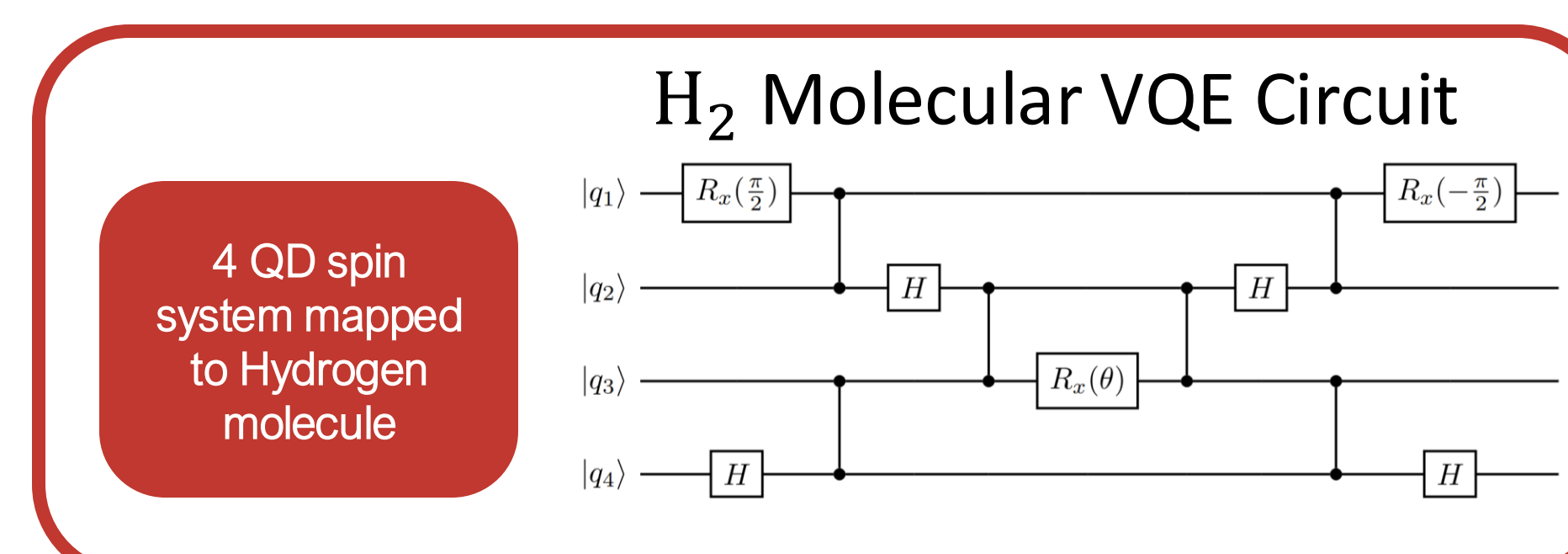
- Incorporation of realistic device geometries
- Account for electrostatic cross-talk between QDs
- Universal set of gates & quantum algorithms



Application

Variational Quantum Eigen-Solver (VQE)

VQE sets an upper bound on ground state energy for a molecular electronic system. Approximating the multi-electron wavefunction is crucial in capturing correlation features in a many-electron system. With proper choice of Ansatz and optimization routine, a multi-electron wavefunction can be computed efficiently and accurately [3].

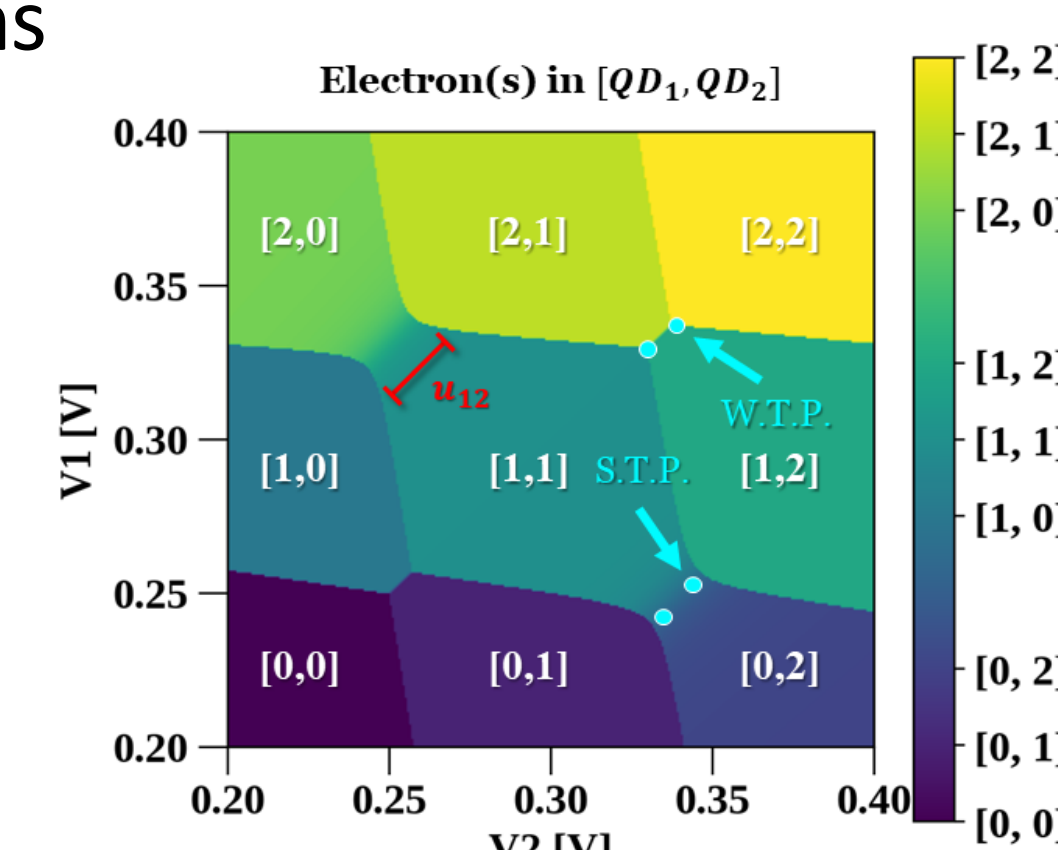


Results

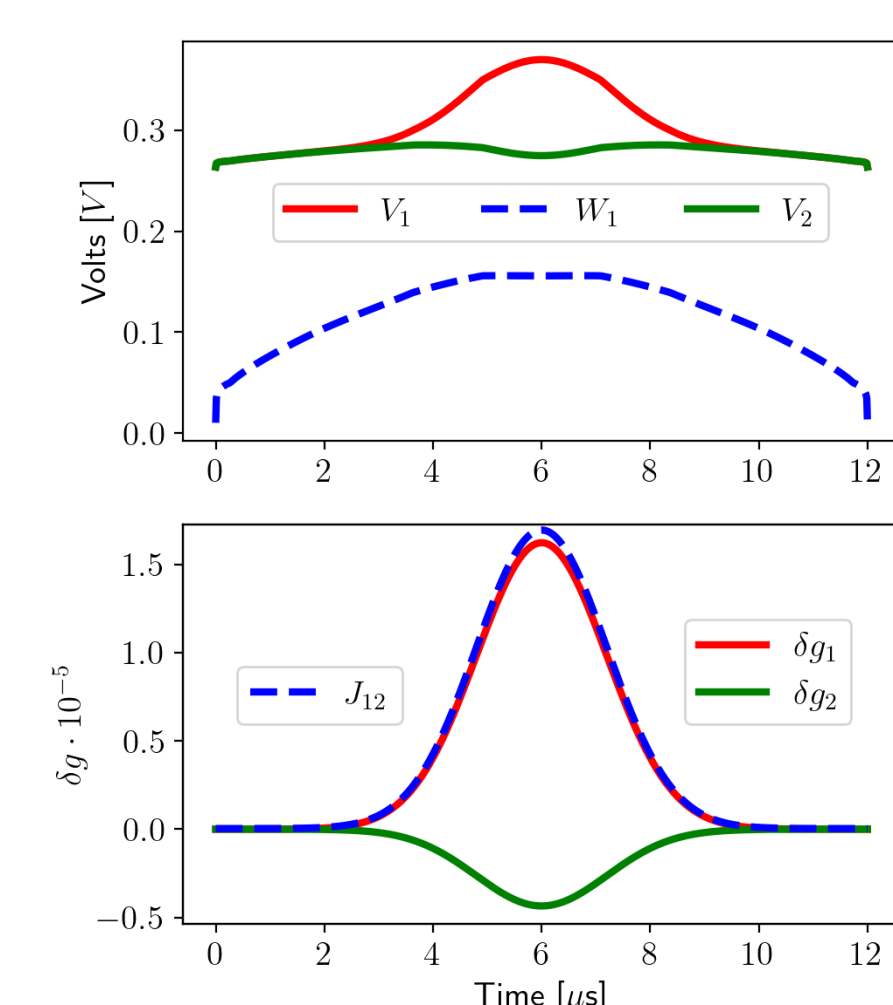
Charge Stability Regions

Charge stability regions for a double QD device. Coulomb repulsion and tunnel coupling features:

- Inform optimal control: pulse engineering & experiment



Qubit Gate Operations

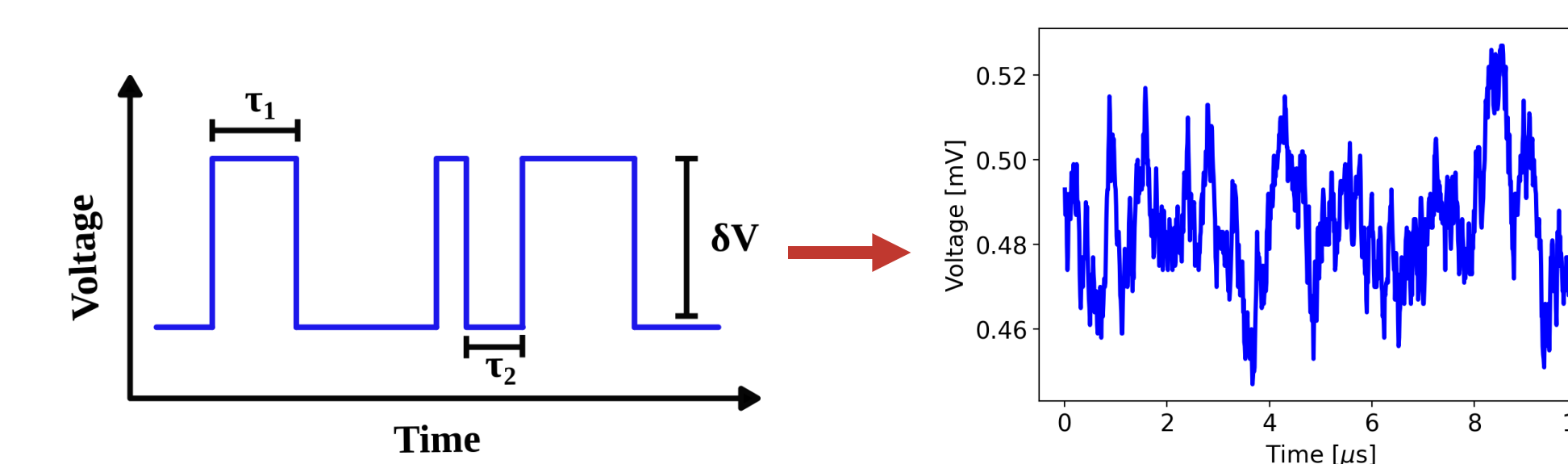


Global Electron Spin Resonance (ESR) field with gate voltage control enables effective parameter control:

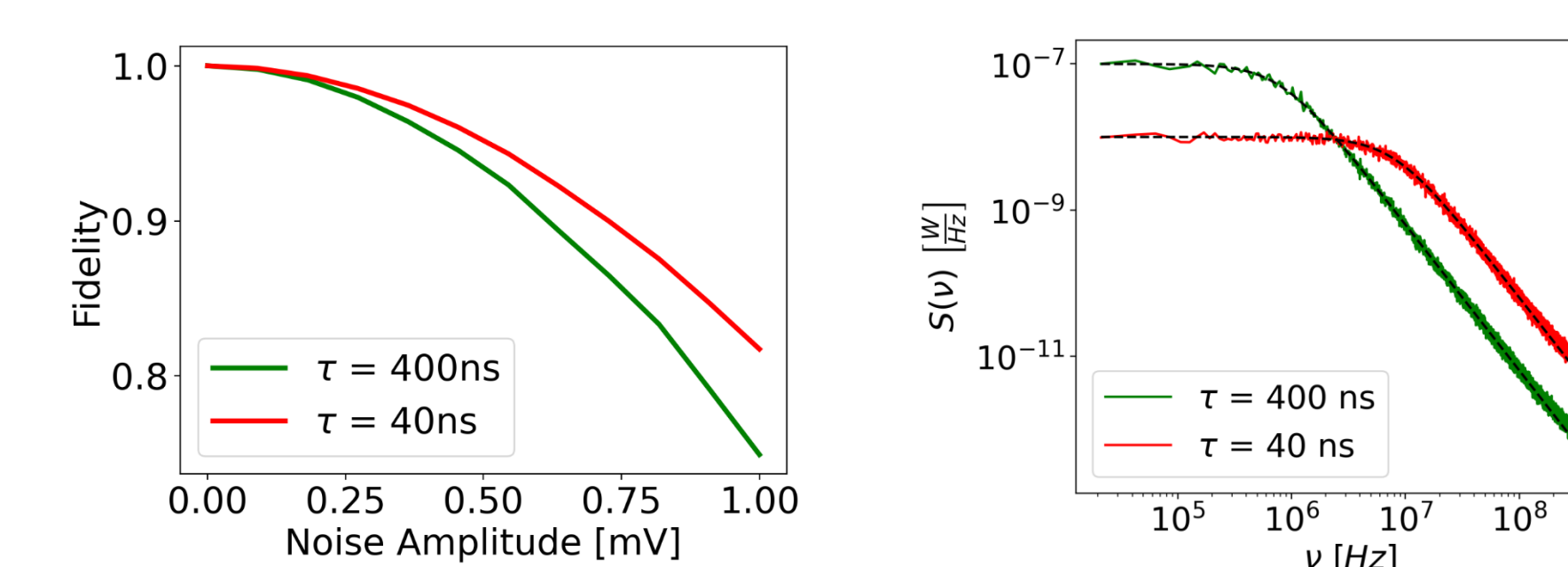
- ESR 1-qubit gates
- Voltage-only 2-qubit gates

Noisy Intermediate Scale Quantum (NISQ) Devices

Simulation of electrical noise generated by an ensemble of Random Telegraph Noise (RTN) fluctuators for varied switching times τ . A Coupled Cluster (CC) designed VQE is applied to a 4-qubit QD linear array, which optimizes parameter θ for a quantum circuit with gate $R_x(\theta)$ to estimate the ground state wavefunction.



The impact on process fidelity of variational algorithm: RTN with varying electrical amplitudes and switching times \rightarrow mimic that of experimentally observed values in real QD devices.



[1] Merino et al. AMMCS (2024), URL: <https://arxiv.org/abs/2402.15499>
 [2] Khromets et al. AMMCS (2024), URL: <https://arxiv.org/abs/2402.08146>
 [3] Yuxuan Du et al. npj QI (2022), DOI: 10.1038/s41534-022-00570-y.

Acknowledgments: This research was undertaken thanks in part to funding from NSERC and Canada First Research Excellence Fund (TQT)