Divide-&-Conquer Maxwell-Ehrenfest-Surface Hopping (DC-MESH)

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Goal: To describe multiscale light-matter interaction ranging from atto-to-nano seconds & pico-to-micro meters



Linker et al., Science Adv. 8, eabk2625 (2022) Razakh et al., PDSEC (IEEE, '24)



Dawn of Attosecond Physics

The Nobel Prize in Physics 2023



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Pierre Agostini

Prize share: 1/3



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Ferenc Krausz

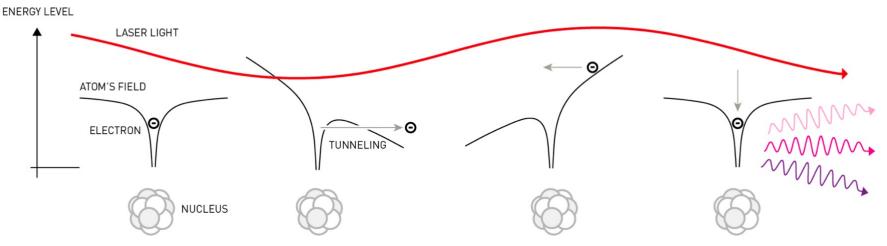
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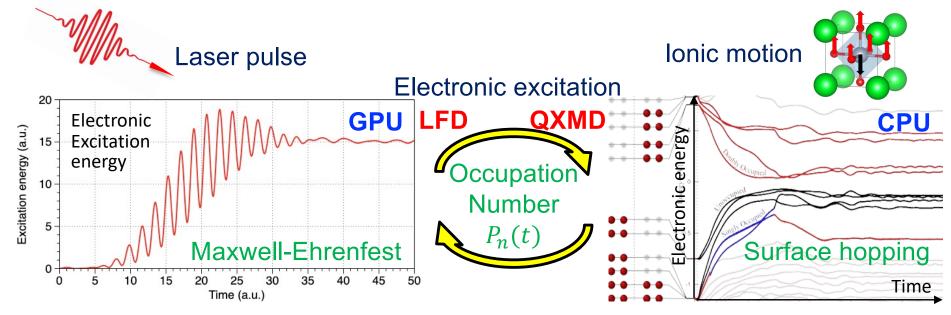
Anne L'Huillier
Prize share: 1/3

The Nobel Prize in Physics 2023 was awarded to Pierre Agostini, Ferenc Krausz and Anne L'Huillier "for experimental methods that generate attosecond pulses of light for the study of electron dynamics in matter"



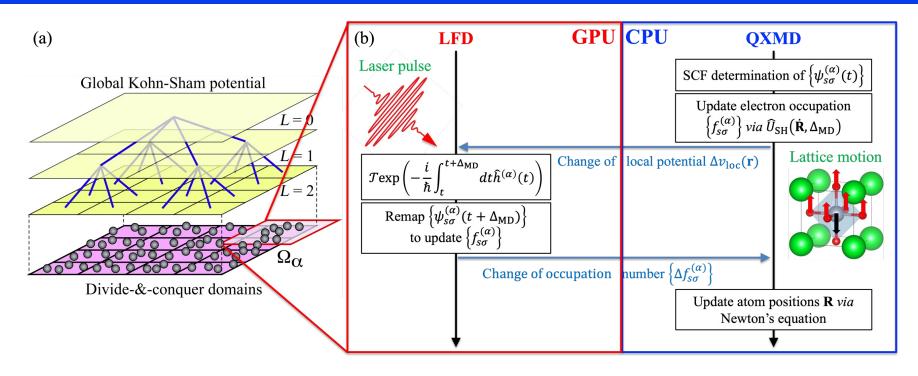
Nonadiabatic Quantum MD: DC-MESH

- DC-MESH (divide-&-conquer Maxwell + Ehrenfest + surface-hopping): O(N) algorithm to simulate photo-induced quantum materials dynamics
- LFD (local field dynamics): Maxwell equations for light & real-time time-dependent density functional theory equations for electrons to describe light-matter interaction
- QXMD (quantum molecular dynamics with excitation): Nonadiabatic coupling of excited electrons & ionic motions based on surface-hopping approach
- "Shadow" LFD (GPU)-QXMD (CPU) handshaking via electronic occupation numbers with minimal CPU-GPU data transfer
- GSLD: Globally sparse (interdomain Hartree coupling *via* multigrid) & locally dense (intradomain nonlocal exchange-correlation computation *via* BLAS) solver



Linker et al., Science Adv. 8, eabk2625 (2022); Razakh et al., PDSEC (IEEE, '24)

Divide-Conquer-Recombine (DCR)



- Treat multi-physics at appropriate scales & levels of approximation
- Hartree potential & electromagnetic field are computed globally using the scalable O(N) multigrid method & macroscopic grid, respectively
- Higher-order correlations represented by the exchange-correlation (XC) kernel in time-dependent density functional theory (TDDFT) are treated locally within each divide-&-conquer (DC) domain since they are known to be short-ranged [Nakano & Ichimaru, *Phys. Rev. B* **39**, 4930 ('89)]
- See notes on <u>dynamic correlation</u>, <u>DCR-NAQMD</u>, <u>embedded TDDFT</u>, and <u>Ehrenfest-hopping dynamics (EHD)</u>

LFD Algorithm

Hamiltonian in the α-th domain [Yabana, Phys. Rev. B 85, 045134 ('12); Jestadt, Adv. Phys. 68, 225 ('19)]

$$\hat{h}(t, \mathbf{R}(t)) = \frac{1}{2} \left(\frac{\nabla}{i} + \frac{1}{c} \mathbf{A}(\mathbf{r}_{\alpha}, t) \right)^{2} - \phi(\mathbf{r}_{\alpha}, t) + \hat{v}_{xc} + v_{ion}(\mathbf{r}, \mathbf{R}) + \Delta \dot{\mathbf{R}} \cdot \frac{\partial}{\partial \mathbf{R}} v_{ion}$$

Electromagnetic vector & scalar potentials at the α -th domain

Nonadiabatic coupling

Trotter expansion of time propagator

exp
$$\left(-i\hat{h}\Delta_{\mathrm{MD}}\right) \cong \exp\left(-i\hat{h}_{\mathrm{el-ion}}\Delta_{\mathrm{MD}}/2\right)\mathcal{T}\exp\left(-i\int_{t}^{t+\Delta_{\mathrm{MD}}}dt\hat{h}_{\mathrm{el}}(t)dt\right)\exp\left(-i\hat{h}_{\mathrm{el-ion}}\Delta_{\mathrm{MD}}/2\right)$$
QXMD

QXMD

QXMD

Self-consistent propagator [Sato, J. Chem. Phys. 143, 224116 ('15); Lian, Adv. Theo. Sim. 1, 1800055 ('18)]

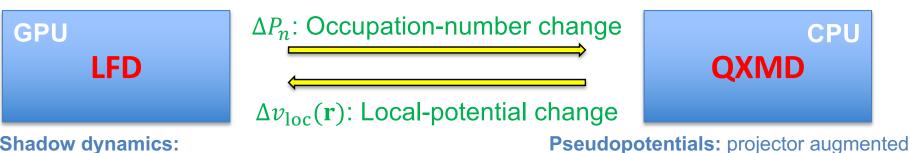
$$\mathcal{T}\exp\left(-i\int_{t}^{t+\Delta_{\mathrm{MD}}}dt\hat{h}_{\mathrm{el}}(t)\right) \cong \prod_{n=1}^{N_{\mathrm{QD}}=\Delta_{\mathrm{MD}}/\Delta_{\mathrm{QD}}}\exp\left(-i\Delta_{\mathrm{QD}}\hat{h}_{\mathrm{el}}\left(t+\left(n-\frac{1}{2}\right)\Delta_{\mathrm{QD}}\right)\right)$$

Nonlocal exchange-correlation propagator [Vicek, J. Chem. Phys. 150, 184118 ('19)]

See note on self-consistent time propagator

Reduced-Communication Shadow Dynamics

- Fundamental physics equations are local at the finest spatiotemporal scales, *i.e.*, simple partial differential equations with differential operators acting locally in a data-parallel fashion LFD fits naturally to GPU; on the other hand, coarsegrained schemes to approximately describe complex chemical interactions often come with an excessive computational cost of nonlocal operations in space and time QXMD takes advantage of complex instruction sets in CPU
- At each molecular-dynamics step, LFD informs QXMD of occupation-number change due to light-electron & electron-electron interactions
- QXMD performs excited-state quantum molecular dynamics & informs LFD of local-potential change for the next $N_{\rm QD}$ (= $\Delta_{\rm MD}/\Delta_{\rm QD}$) quantum-dynamics steps
- "Shadow" electronic wave functions in LFD are resident on GPU, while QXMD wave functions on CPU, to minimize CPU-GPU data transfers



Shadow dynamics: real-time time-dependent density functional theory (RT-TDDFT)

Exchange-correlation functionals: metaGGA (SCAN), hybrid exactexchange (HSE *etc.*), DFT+U, DFT-D,

wave (PAW), ultrasoft (Vanderbilt)

nonlocal vdW

See note on shadow EHD

Data-Parallel & BLASified LFD

- Data-parallel local LFD: Auxiliary-field electronic time propagator for local potential [Car & Parrinello, Solid State Commun. 62, 403 ('87); Nakano et al., Comput. Phys. Commun. 83, 181 ('94)] on real-space mesh achieves high performance on GPU
- BLASified nonlocal LFD: Operation of nonlocal potential is projected onto a vector space spanned by Kohn-Sham orbitals at time 0 within the real-time scissor approximation [Wang et al., J. Phys. Condens. Mat. 31, 214002 ('19)], making it dense matrix operations implemented with highly optimized level3 (or matrix-matrix) BLAS (basic linear algebra subprogram) library on GPU

$$\hat{v}_{\rm nl}|\psi_n(t)\rangle \cong \Delta_{\rm sci} \sum_{m\geq {
m LUMO}} |\psi_m\rangle\langle\psi_m|\psi_n(t)\rangle$$

• See notes on auxiliary-field electron propagator and real-time scissor

Global Maxwell's Equations

• Global Maxwell's equations are solved on a macroscopic grid

Yabana, Phys. Rev. B 85, 045134 ('12); cf. Gabay, Phys. Rev. B 101, 235101 ('20)

$$\begin{split} \mathbf{A} &= \mathbf{A}_{\text{ext}} + \mathbf{A}_{\text{ind}} + \mathbf{A}_{\text{xc}} \\ & \left(\frac{1}{c^2} \frac{\partial^2}{\partial t^2} - \frac{\partial^2}{\partial \mathbf{R}^2} \right) \mathbf{A}_{\text{ind}} = \frac{4\pi}{c} \mathbf{J} \quad \text{Induced vector potential} \\ & \left(\frac{1}{c^2} \frac{\partial^2}{\partial t^2} - \frac{\partial^2}{\partial \mathbf{R}^2} \right) \phi = \frac{4\pi}{c} \rho \quad \text{Scalar potential} \end{split}$$

• Local domain-averaged current & charge densities

$$\mathbf{J}(\mathbf{R}_{\alpha},t) = \frac{1}{\Omega_{\alpha}} \int_{\Omega_{\alpha}} d\mathbf{r} \mathbf{j}(\mathbf{r},t)
\rho(\mathbf{R}_{\alpha},t) = -\frac{1}{\Omega_{\alpha}} \int_{\Omega_{\alpha}} d\mathbf{r} n(\mathbf{r},t)
\mathbf{j}(\mathbf{r},t) = -\sum_{n\sigma} \operatorname{Re} \left[\psi_{n\sigma}^{*}(\mathbf{r},t) \frac{\nabla}{i} \psi_{n\sigma}(\mathbf{r},t) \right] f_{n\sigma} - \frac{1}{c} \mathbf{A}(\mathbf{r},t) n(\mathbf{r},t)
n(\mathbf{r},t) = \sum_{n\sigma} \left| \psi_{n\sigma}(\mathbf{r},t) \right|^{2} f_{n\sigma}$$

• Long-range correction in time-dependent current density functional theory (TDCDFT) [Vignale, *Phys. Rev. Lett.* 77, 2037 ('96); Maitra, *Phys. Rev. B* 68, 045109 ('03); Sun, *Phys. Rev. Lett.* 127, 077401 ('21)]

$$\left(\frac{1}{c^2}\frac{\partial^2}{\partial t^2} - \frac{\partial^2}{\partial \mathbf{R}^2}\right)\mathbf{A}_{xc} = -\frac{\alpha}{c}\mathbf{J}$$
 Exchange—correlation vector potential

See note on <u>Maxwell solver</u> and <u>exciton dynamics</u>