Density functional theory and optimal transportation with Coulomb and Riesz cost.

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Outline

Informal introduction to Quantum mechanics/DFT

- All materials systems we study essentially consist of electrons and nuclear charge.
- Mechanical, electronic, magnetic etc. properties are due to electrons and their interaction with other electrons.
- In order to define electrons and their interaction we use Schrodinger equation.
- It allows to predict, e.g., binding energies, equilibrium geometries, intermolecular forces
- Quantum mechanics for electrons reduces to a PDE (the Schroedinger equation)

Density Functional Theory (DFT)

- If Schrodinger equation for the many electrons problem could be solved accurately and efficiently then almost any property of the materials could be determined determined accurately.
- Unfortunately, there is neither an accurate nor an efficient method to solve these problems.
- DFT is a simplified version of quantum mechanics (QM), widely used in molecular simulations in chemistry, physics, materials science
- **Main idea**: describe complicated N-particle system (a probability on $\mathbb{R}^{3N}$) using only its one-particle marginal $\rho(x)$ on $\mathbb{R}^3$.
- Feasible system size: up to a million atoms

Some history of DFT

- Thomas-Fermi: 1920s simple model
- Hohenberg-Kohn-Sham (1963-1964): practical method based on semi-empirical functions of ρ
- Levy (1979), Lieb (1983): mathematical justification
- 1970s: popular in solid state physics, but not so accurate
- 1990s: explosion in quantum chemistry, due to increase of computational resources + discovery of efficient semi-empirical functionals of ρ
- 1998 Nobel Prize for 'founding father' Walter Kohn
- About 15 000 papers per year with the keyword 'density functional theory'

Quantum mechanics-Formal definition

- The solution for this PDE is the wave function $\Psi(x_1, s_1, \ldots, x_N, s_N) \in L^2((\mathbb{R}^3 \times \mathbb{Z}_2)^N; \mathbb{C})$
- N - number of electrons, x_i position of electron i, s_i spin of electron i
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$$|\Psi(x_1, s_1, \ldots, x_N, s_N)|^2 = \text{probability density that the electrons are at positions } x_i \text{ with spins } s_i.$$

Ψ is an anti-symmetric function, which makes $|\Psi|^2$ a symmetric (N-exchangeable) probability measure.

- $$\mathcal{A}_N = \{\Psi \in L^2((\mathbb{R}^3 \times \mathbb{Z}_2)^N; \mathbb{C}) \,|\, \nabla\Psi \in L^2,\ \Psi \text{ antisymmetric},\ ||\Psi||_{L^2} = 1\}$$

- Key quantum mechanics quantity is the ground state energy E_0

$$E_0 = \inf_{\Psi \in \mathcal{A}_N} E[\Psi]$$

where

$$E[\Psi] = T_h[\Psi] + V_{ee}[\Psi] + V_{ne}[\Psi]$$

- Kinetic energy:

$$T_h[\Psi] = \frac{h^2}{2} \int |\nabla \Psi(x_1, s_1, \ldots, x_N, s_N)|^2 dz_1 dz_2 \ldots dz_N$$

- Electron-electron energy:

$$V_{ee}[\Psi] = \int \sum_{1 \le i < j \le N} \frac{1}{|x_i - x_j|} |\Psi|^2 dz_1 \ldots dz_N$$

- Nuclei-electron energy:

$$V_{ne}[\Psi] = \sum_{i=1}^{N} \int v(x_i) |\Psi(z_1, z_2, \ldots, z_N)|^2 dz_1 \ldots dz_N$$

- N-electrons density

$$\rho_N^{\Psi}(x_1,..,x_N) = \sum_{s_1,..,s_N \in \mathbb{Z}_2} |\Psi(x_1,s_1,\ldots,x_N,s_N)|^2.$$

- Pair electrons density

$$\rho_2^{\Psi}(x_1,x_2) = \binom{N}{2} \int_{\mathbb{R}^{3(N-2)}} \rho_N^{\Psi}(x_1,\ldots,x_N) dx_3 \ldots dx_N$$

- Single electron density

$$\rho^{\Psi}(x_1) = N \int_{\mathbb{R}^{3(N-1)}} \rho_N^{\Psi}(x_1,\ldots,x_N) dx_2 \ldots dx_N.$$

- $\mathcal{R}_N := \{\rho : \mathbb{R}^3 \to \mathbb{R} \,|\, \rho \text{ is the density of some } \Psi \in \mathcal{A}_N\}$

Full Scrod. eqn. can be reformulated as a hierarchy of eqn: for ρ in terms of of the pair electrons density ρ_2, for ρ_2 in terms of ρ_3 etc.

Variational formulation of density functional theory

(Hohenberg/Kohn 1964, M. Levy 1979, E. Lieb 1983)
For any external potential v, the exact Schroedinger eqn. satisfies

- $$E_0 = \inf_{\rho \in \mathcal{R}_N} \left\{ F_h[\rho] + N \int_{\mathbb{R}^3} v(x)\, \rho(x) dx \right\}$$

  with

- $$\begin{aligned} F_h[\rho] : &= \inf_{\Psi \in \mathcal{A}_N, \Psi \mapsto \rho} \left\{ T_h[\Psi] + V_{ee}[\Psi] \right\} \\ &= \inf_{\Psi \in \mathcal{A}_N, \Psi \mapsto \rho} \left\{ T_h[\Psi] + \int_{\mathbb{R}^6} \frac{1}{|x-y|} \rho_2^{\Psi}(dx, dy) \right\}, \end{aligned}$$

  $F_h[\rho]$ is the famous Hohenberg-Kohn functional.

- Not useful for computations (definitely still contains the big space of $\Psi(x_1, s_1, \ldots, x_N, s_N)$). But useful starting point for model reduction in asymptotic limits.

Correlations in DFT

- Mathematical structure: Minimize an approximate energy functional $F[\rho]$ which depends on the electron density $\rho(x)$, a function on $\mathbb{R}^3$.
- Catch: exact QM energy requires knowledge of electron-pair density $\rho_2(x,y)$, a function on $\mathbb{R}^6$, which entails correlations.
- Standard way out: start by assuming independence, add semi-empirical corrections to $F_h[\rho]$ accounting for correlations. Often but not always accurate/reliable.

Popular functionals

All functionals used in practice are of form

$$\text{Mean field} \ + \ \text{additive corrections.}$$

Why mean field? Interactions not weaker than single-particle terms.

- The mean field approximation:

$$\int_{\mathbb{R}^6} \frac{1}{|x-y|} \rho_2^{\Psi}(dx,dy) = \frac{1}{2} \int_{\mathbb{R}^6} \frac{1}{|x-y|} \rho(dx)\rho(dy) =: J[\rho].$$

- Local Density Approximation approximation:

$$\int_{\mathbb{R}^6} \frac{1}{|x-y|} \rho_2^{\Psi}(dx,dy) = J[\rho] - \frac{4}{3}(3/\pi)^{1/3} \int_{\mathbb{R}^3} \rho(x)^{4/3} dx.$$

- Quantum mechanics is becoming so unbelievably complex that it is taking longer and longer to train a quantum theorist.
- It is taking so long, in fact, to train him to the point where he understands the nature of physical problems that he is already too old to solve them. (*Eugene Wigner*)
- Most cited physicist of all time is a designer of DFT models, J.Perdew.

- ρ_N measure in $\mathbb{R}^{Nd}$, ρ measure in $\mathbb{R}^d$
- $0 < s < d$
- Minimize the transportation cost

$$\int_{\mathbb{R}^{Nd}} \left(\sum_{1 \le i < j \le N} \frac{1}{|x_i - x_j|^s} \right) d\rho_N(x_1, \ldots, x_N)$$

subject to the constraint

$$\int_{\mathbb{R}^{(N-1)d}} \rho_N(x_1, \ldots, x_N) dx_2 \ldots dx_N = \rho(x) \ldots$$

$$\int_{\mathbb{R}^{(N-1)d}} \rho_N(x_1, \ldots, x_{N-1}, x_N) dx_1 \ldots dx_{N-1} = \rho(x).$$

- We can symmetrise the measure ρ_N-finite exchangeable.
- Coulomb: $s = 1, d = 3$

The 2-marginal Optimal Transport Problem with Coulomb Cost

- ρ_2 measure in $\mathbb{R}^{2d}$, ρ measure in $\mathbb{R}^d$
- Minimize the transportation cost

$$\int_{\mathbb{R}^{2d}} \frac{1}{|x-y|} d\rho_2(x,y)$$

subject to the constraint

$$\int_{\mathbb{R}^d} \rho_2(x,y)dy = \rho(x) \quad \text{and} \quad \int_{\mathbb{R}^d} \rho_2(x,y)dx = \rho(y).$$

- General pattern: $c : \mathbb{R}^d \times \mathbb{R}^d \to \mathbb{R} \cup \{+\infty\}$,with $c(x,y) := l(|x-y|)$, such that $l \geq 0$ is strictly convex, strictly decreasing and C^1 on $(0,\infty)$, $l(0) = +\infty$.

Theorem

(C, Friesecke, Klueppelberg - CPAM 2013) *Let $c(x,y) := l(|x-y|)$, such that $l \geq 0$ is strictly convex, strictly decreasing and C^1 on $(0,\infty)$, $l(0) = +\infty$, ρ absolutely continuous with respect to the Lebesgue measure. Then*

- *There exists a unique optimizing measure ρ_2 with*

$$\rho_2 = (id, T)_{\#\rho},$$

where the optimal map $T : \mathbb{R}^d \to \mathbb{R}$ is unique. Moreover $\rho \circ T^{-1} = \rho$.

For Coulomb cost, we have also $T(x) = x + \frac{\nabla v(x)}{|\nabla v(x)|^{3/2}}$ for some convex potential $v : \mathbb{R}^3 \to \mathbb{R}$

- Physical meaning 1: $T(x) =$ position of the 2nd electron if the first electron is at x.
- Physical meaning 2: the graph of T is the support of the electron pair density ρ_2.

The Method

- Adaptation of **Gangbo and McCann:** The geometry of optimal transportation, *Acta Math.* **177**, 113-161 (1996).
- Check that formula (originally for increasing costs) generalizes to decreasing costs with singularity on the diagonal.
- Explicit Solution: For $d = 1$, for all marginals. As simple example, take ρ to be the uniform measure on $[0, 1]$. Then $\rho_2 = (id, T)_{\#\rho}$, and T rigidly switches right and left half of $[0, 1]$.
- More precisely, we have $T(x) = x + 1/2$ for $x < 1/2$, and $T(x) = x - 1/2$ for $1/2 < x < 1$.

- For $d \geq 2$, explicit solution for all symmetric marginals. Then for ρ_1 and ρ_2 densities of μ, ν, with $\rho_1(x) = \lambda_1(|x|)$ and $\rho_2(x) = \lambda_2(|x|), x \in \mathbb{R}^d$.
- Then T is of form: $T(x) = x\frac{g(|x|)}{|x|}, x \in \mathbb{R}^d$, with $g : [0, \infty) \to \mathbb{R}$. Moreover $g \leq 0$, and g is an increasing function with $g(0_+) = -\infty$ and $g(+\infty) = 0$.
- Physical interpretation: 2nd electron is in the opposite direction of first.

Theorem

(C, Frieescke, Klueppelberg - CPAM 2013) *Suppose that $\mu = \nu$. Let $t \in (0, \infty)$ and let*

$$F_1(t) = |S^{d-1}| \int_0^t \lambda(s) s^{d-1} ds$$

and

$$F_2(-t) = |S^{d-1}| \int_t^\infty \lambda(s) s^{d-1} ds.$$

Then

$$g(t) = F_2^{-1}(F_1(t)).$$

Let

- $$E_0 = \inf_{\Psi \in \mathcal{A}_N} \left\{ T_h[\Psi] + V_{ne}[\rho^\Psi] + V_{ee}[\rho_2^\Psi] \right\}$$

  and

- $$E_0^{OT} = \inf_{\Psi \in \mathcal{A}_N} \left\{ T[\Psi] + V_{ne}[\rho^\Psi] + E_{OT}^N[\rho^\Psi] \right\},$$

  where recall

$$E_{OT}^N[\rho^\Psi] = \inf_\gamma \sum_{1 \le i < j \le N} \int \frac{1}{|x_i - x_j|} d\gamma(x_1, x_2, \ldots, x_N),$$

subject to equal marginals ρ^Ψ.

Theorem

(C, Friesecke, Klueppelberg - CPAM 2013) *Fix $\rho \in \mathcal{R}_N$. Let $N = 2$. Then*

$$\lim_{h \to 0} F_h[\rho] = E^N_{OT}[\rho]$$

for every $\rho \in \mathcal{R}_N$, where recall that

$$F_h[\rho] := \inf_{\Psi \in \mathcal{A}_N, \Psi \mapsto \rho} \Big\{ T_h[\Psi] + V_{ee}[\Psi] \Big\}.$$

- De Pascale - Bindini (2017): extension to $N = 3$.
- C, Friesecke, Klueppelberg (in preparation): extension to $N \geq 4$

Theorem

(C,Friesecke, Klueppelberg - CPAM 2013) *For every N and every v*

$$E_0 \geq E_0^{OT}.$$

Seidl'99, Seidl/Perdew/Levy 1999, Seidl/Gori-Giorgi/Savin 2007

Many-marginals optimal transport problem

- Problem coincides with infimum over Monge states "strongly correlated electrons" (Colombo-Di Marino 2015)
- Existence and uniqueness of Monge for $N \geq 2$ in $d = 1$ (Colombo-De Pascale-Di Marino 2013)
- Buttazzo, De Pascale & Gori-Giorgi (2012); Pass (2013); Benamou, Carlier & Nenna (2015); Di Marino, Gerolin & Nenna (2015)

The infinite Optimal Transportation marginal problem

- Let γ be an infinite dimensional measure, γ symmetric (exchangeable), ρ probability measure in $\mathbb{R}^d$.

$$F_{OT}^{\infty}[\rho] = \inf_{\gamma} \lim_{N\to\infty} \frac{1}{\binom{N}{2}} \int_{\mathbb{R}^{dN}} \sum_{1\le i<j\le N} \frac{1}{|x_i - x_j|^{d-2}} d\gamma(x_1, .., x_N),$$

subject to the constraint

$$\int_{\mathbb{R}\times\mathbb{R}\times\ldots} \gamma(x_1, x_2, \ldots, x_N, \ldots) dx_2 dx_3 \ldots = \rho(x_1).$$

Theorem

(C, Friesecke, Pass - Calc Var PDEs 2015)

$$\lim_{N\to\infty} F_{OT}^{N}[\rho] = F_{OT}^{\infty}[\rho] = \frac{1}{2} \int_{\mathbb{R}^{2d}} \frac{1}{|x-y|^{d-2}} \rho(x)\rho(y) dxdy.$$

- Proof by use of de Finetti's Theorem and Fourier transforms
- De Finetti theorem: Let γ_∞ be a symm. inf. dim. Borel measure. Then there exists a unique Borel prob. measure ν such that

$$\gamma_\infty = \int_{\mathcal{P}(\mathbb{R}^d)} Q^{\otimes\infty} d\nu(Q).$$

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$$F^\infty_{OT}[\rho] = \int_{\mathbb{R}^{2d}} \ell(x-y)\,\rho(dx)\,\rho(dy) + \int_{\mathbb{R}^d} \hat{\ell}(z)\left(\mathrm{var}_{\nu(dQ)} Re(\hat{Q}(z)) + \mathrm{var}_{\nu(dQ)} Im(\hat{Q}(z))\right) dz$$

Let γ be an infinite dimensional measure, γ symmetric (exchangeable), ρ probability measure in $\mathbb{R}^d$ and let $0 < s < d$.

$$F^{\infty}_{OT,s}[\rho] = \inf_{\gamma} \lim_{N\to\infty} \frac{1}{\binom{N}{2}} \int_{\mathbb{R}^{dN}} \sum_{1\le i<j\le N} \frac{1}{|x_i - x_j|^s} d\gamma(x_1, .., x_N),$$

subject to the constraint

$$\int_{\mathbb{R}\times\mathbb{R}\times\ldots} \gamma(x_1, x_2, \ldots, x_N, \ldots) dx_2 dx_3 \ldots = \rho(x_1).$$

Then by de Finetti's Theorem and positive-definiteness, we have

Theorem

(Petrache 2015)

$$\lim_{N\to\infty} F^{N}_{OT,s}[\rho] = F^{\infty}_{OT}[\rho] = \frac{1}{2} \int_{\mathbb{R}^{2d}} \frac{1}{|x-y|^s} \rho(x)\rho(y) dxdy.$$

C, di Marino, Lewin, Lieb, Petrache ...

$$\lim_{N\to\infty} N^{1-s/d}\left(F^N_{OT,s}[\rho] - F^\infty_{OT}[\rho]\right) = ?$$

- Lieb-Oxford bound

$$N^{1-s/d}\left(F^N_{OT,s}[\rho] - F^\infty_{OT}[\rho]\right) \geq -C_{LO}\int_{\mathbb{R}^d} g(x)^{1+s/d}dx.$$

- Trivially, we have

$$F^N_{OT,s}[\rho] - F^\infty_{OT}[\rho] \leq 0.$$

- Question: Does the limit exit?

- Exists $C(d,s,\rho) > 0$ such that

$$\lim_{N\to\infty} N^{1-s/d}\left(F^N_{OT,s}[\rho] - F^\infty_{OT}[\rho]\right) = -C(d,s,\mu)\int_{\mathbb{R}^d} g(x)^{1+s/d}dx$$

- Uniform marginal ρ (uniform electron gas): $C(d,s,\mu) = C_{unif}$
- Chemists conjecture for $s = 1, d = 3$: $C(d,s,\mu) = C_{unif}$ (Rasanen, Pittalis, Capelle & Proetto 2009)
- Exact value is unknown, although everybody thought for decades that it is approx 1.4441, related to Epstein Zeta function
- C_{unif} is exactly known for $N = 1$ (di Marino-2017).
- Numerics by Seidl-Vuckovic-Gori Giorgi 2015 ($N = 50$). $C_{unif} \geq 1.401$.

Connection with Coulomb and Riesz gases (Jellium)

- N electrons and a neutralizing background in a domain Ω with $|\Omega| = N$.
- Minimize over x_i

$$\sum_{1\leq i<j\leq N} \frac{1}{|x_i - x_j|} - \sum_{j=1}^{N} \int_\Omega \frac{1}{|x_j - y|} dy + \frac{1}{2} \int_\Omega \int_\Omega \frac{1}{|x-y|} dxdy$$

- Let minimization be $\xi(N, \Omega)$, then the limit (Lieb & Narnhofer 1975)

$$\lim_{N\to\infty} \frac{\xi(N,\Omega)}{N^{4/3}} = -C_{jel}.$$

- Wigner crystallisation conjecture: in limit $N \to \infty$, the electrons place themselves on a BCC lattice (hexagonal lattice in $d = 2$)
- Lewin-Lieb (2015): comparison with uniform electron gas constant in $d = 3$

Sandier, Serfaty, Rougerie, Petrache ...

Assume that $s \in [d-2, d[$. Let $V : X \to \mathbb{R}$ and denote by

$$H_N(x_1, \dots, x_N) := \sum_{i \neq j} \frac{1}{|x_i - x_j|^s} + N \sum_i V(x_i).$$

Let μ_V be the minimizer of

$$\mathcal{E}_V(\mu) = \int \int \frac{1}{|x-y|^s} d\mu(x) d\mu(y) + \int V(x) d\mu(x)$$

If $\mu_V \ll \mathcal{L}^d$ and the density ρ_V is smooth enough

$$\min_{x_1,\dots,x_N} H_N(x_1, \dots, x_N) - N^2 \mathcal{E}_V(\mu_V) = N^{1+\frac{s}{d}} (\min \mathcal{W}) \int_\Sigma \rho^{1+\frac{s}{d}}(x) dx + o\left(N^{1+\frac{s}{d}}\right).$$

- Abrikosov conjecture: in $2d$, the regular triangular lattice is a minimizing configuration for $\mathcal{W}$.
- Known for Coulomb case (Sandier, Serfaty 2012)
- For general dimension, the conjecture is that the minimum of $\mathcal{W}$ for all $d-2 \leq s < d$ is always achieved by some lattice.

THANK YOU!