

The exact Hohenberg-Kohn functional for a lattice model

Tanja Dimitrov¹(dimitrov@fhi-berlin.mpg.de), Heiko Appel¹, Angel Rubio^{1,2}

[1] Fritz-Haber-Institut der Max-Planck-Gesellschaft, Faradayweg 4-6, 14195 Berlin, Germany
[2] NanoBio Spectroscopy group and ETSF, Universidad del País Vasco, San Sebastián, Spain



MAX-PLANCK-GESELLSCHAFT

Introduction

For a discretized soft-Coulomb lattice model, we investigate the exact solution of the many-body Schrödinger equation in Fock space. Using quadratic optimization with quadratic constraints, or alternatively exact diagonalization, we explicitly construct the exact Hohenberg-Kohn functional and the mapping from densities to wavefunctions. We analyze the resulting exact Hohenberg-Kohn functional and draw conclusions for the construction of approximate functionals.

Levy-Lieb constraint search (M. Levy 1979 [1], E. Lieb 1983 [2])

Expand eigenfunctions in a complete basis set (energy eigenfunctions, Slater-Determinants, etc.)

$$|\Psi[n]\rangle = \sum_{j=1}^M \alpha_j[n] |\phi_j\rangle, \quad M \text{ number of sites}$$

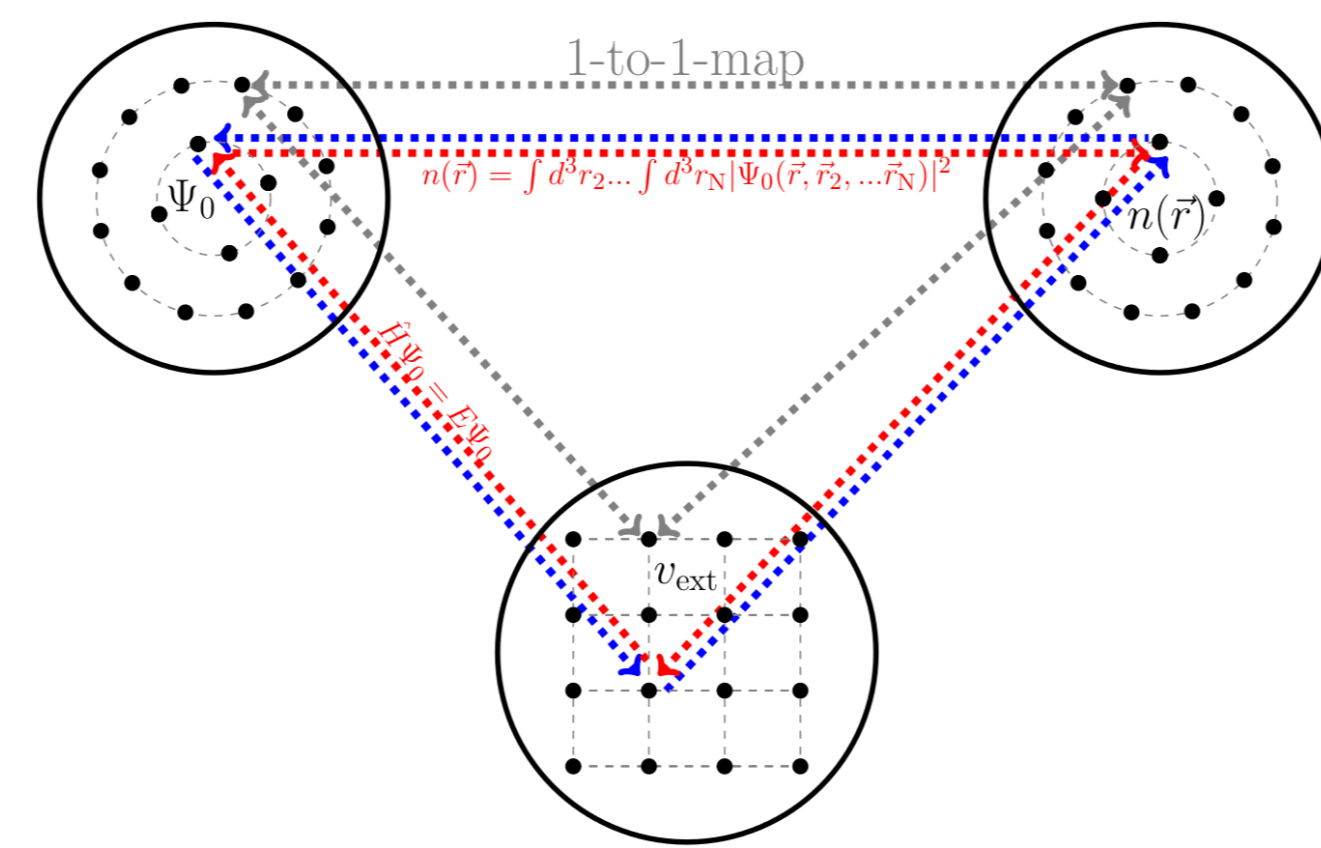
Hohenberg-Kohn functional

$$F_{\text{HK}}(\alpha_1, \dots, \alpha_M)[n] = \min_{\Psi \rightarrow n} \langle \Psi[n] | \hat{T} + \hat{W} | \Psi[n] \rangle \\ = \min_{\Psi \rightarrow n} \sum_{j,k=1}^M \alpha_j^* \alpha_k [n] \langle \phi_j | \hat{T} + \hat{W} | \phi_k \rangle$$

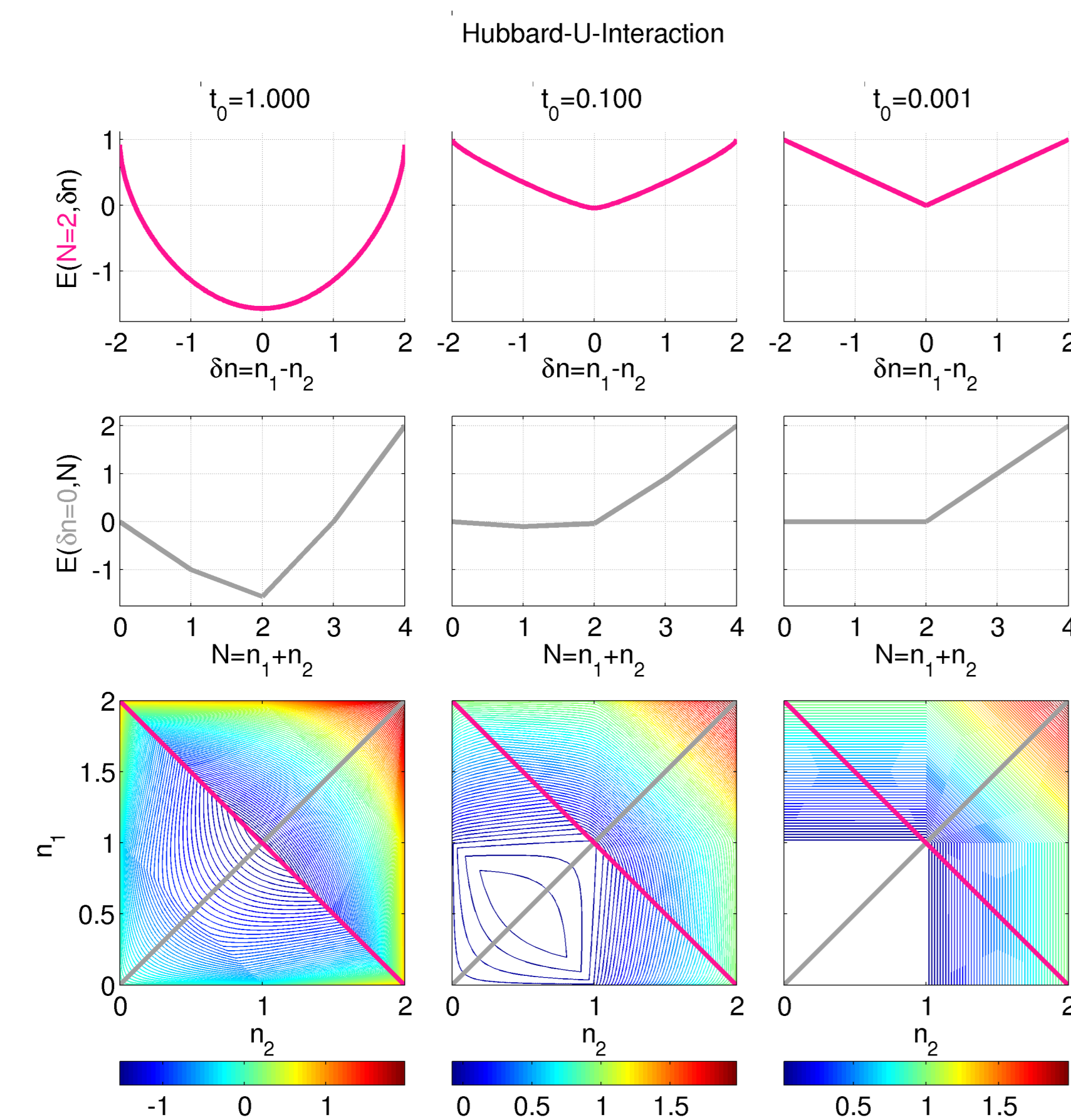
Two-site soft-Coulomb model

$$\hat{H} = \hat{T} + \hat{W} + \hat{V}, \quad \hat{T} = -t_0 \sum_{l,\sigma} (\hat{c}_{l,\sigma}^\dagger \hat{c}_{l+1,\sigma} + \hat{c}_{l+1,\sigma}^\dagger \hat{c}_{l,\sigma}) + 2t_0 \sum_{l,\sigma} \hat{n}_{l,\sigma}, \quad t_0 = \frac{\hbar^2}{2m_e \Delta^2}, \\ \hat{W}_{\text{II}} = U \sum_l \hat{n}_{l,\uparrow} \hat{n}_{l,\downarrow}, \quad \hat{W}_{\text{SC}} = \sum_{l,m,\sigma,\sigma'} \frac{e^2 \hat{c}_{l,\sigma}^\dagger \hat{c}_{m,\sigma'}^\dagger \hat{c}_{m,\sigma'} \hat{c}_{l,\sigma}}{2\sqrt{(l\Delta - m\Delta)^2 + 1}}, \quad \hat{V} = \sum_{l,\sigma} \hat{n}_{l,\sigma} \cdot v_{l,\sigma}, \quad \hat{n} = \sum_{j,\sigma} \hat{c}_{j,\sigma}^\dagger \hat{c}_{j,\sigma}$$

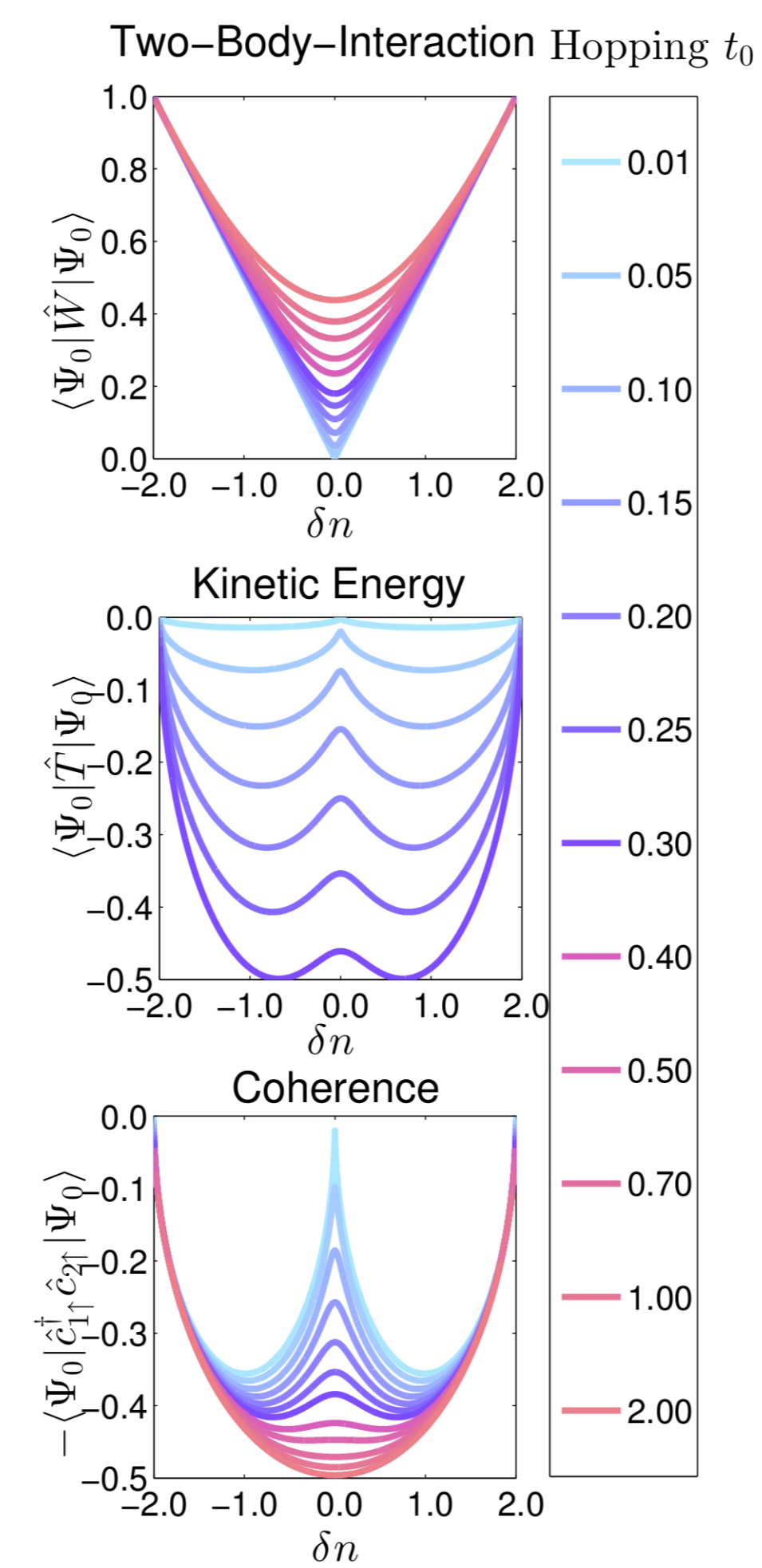
We consider different particle numbers by including a chemical potential μ . Spacing Δ



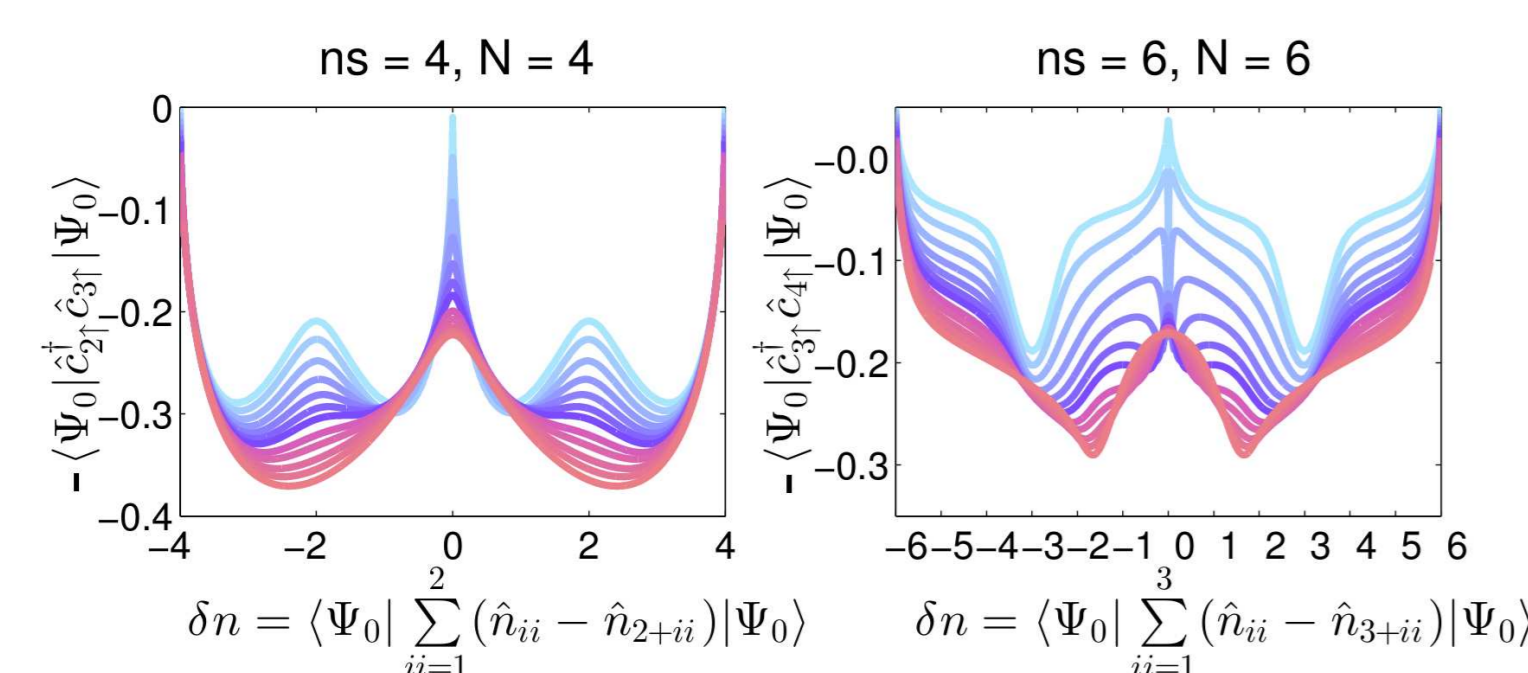
Exact Hohenberg-Kohn-Functional



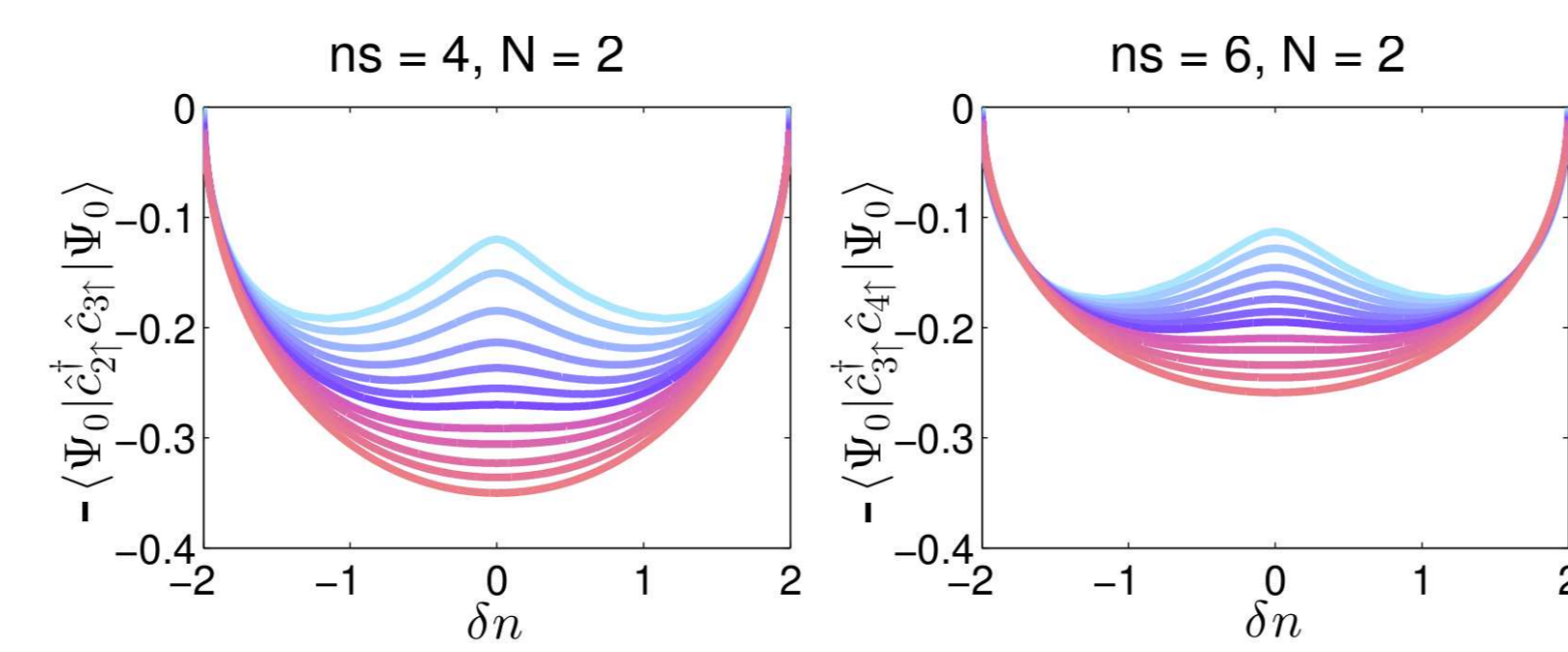
Exact Coherences



Coherences for ns = 4 and ns = 6 (half-filling)



Coherences for ns = 4 and ns = 6 (N=2)



Soft-Coulomb molecules in 1D

Hamiltonian

$$\hat{H}(\alpha) = \hat{T} + \hat{W} + \hat{V}(\alpha) \\ \hat{T} = \sum_{j=1}^2 -\frac{\hbar^2}{2m_e} \frac{d^2}{dx_j^2}, \quad \hat{W} = \frac{1}{2} \sum_{j \neq k} \frac{1}{\sqrt{(x_j - x_k)^2 + 1}} \\ \hat{V}(\alpha) = \sum_{j=1}^2 \frac{Z_j(\alpha)}{\sqrt{(x_j - d)^2 + 1}} + \frac{Z_2(\alpha)}{\sqrt{(x_2 + d)^2 + 1}} \\ Z_1(\alpha) = -\alpha, \quad Z_2(\alpha) = -(2 - \alpha), \quad \alpha \in [0, 2], \quad d = 3, 8 \text{ Bohr}$$

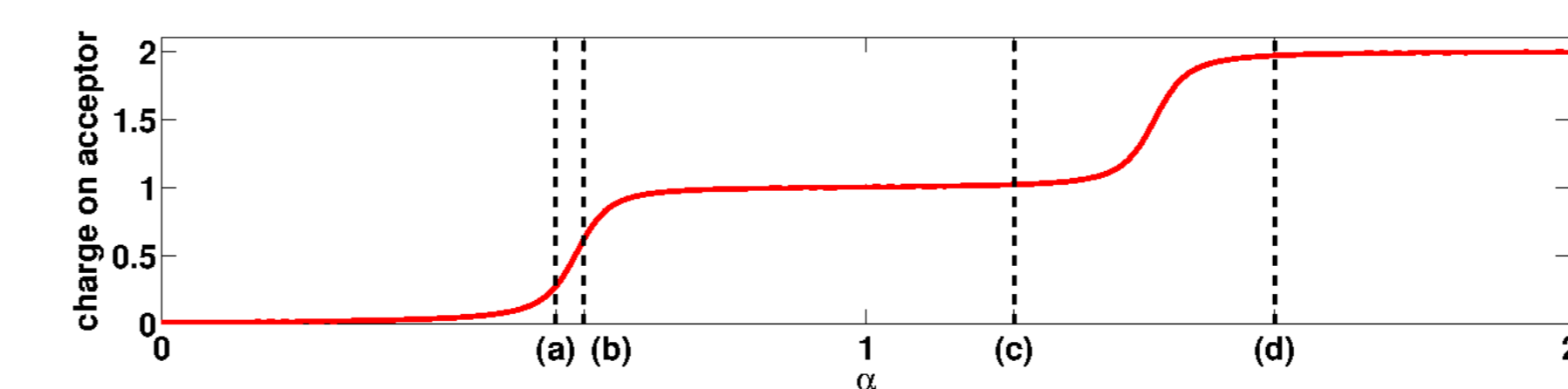
Exact Kohn-Sham potential for two electrons in spin singlet configuration (Helbig et al. 2009 [3])

$$v_{\text{KS}}(x) = \frac{1}{2} \frac{\nabla^2 \sqrt{n(x)}}{\sqrt{n(x)}} + \epsilon_1$$

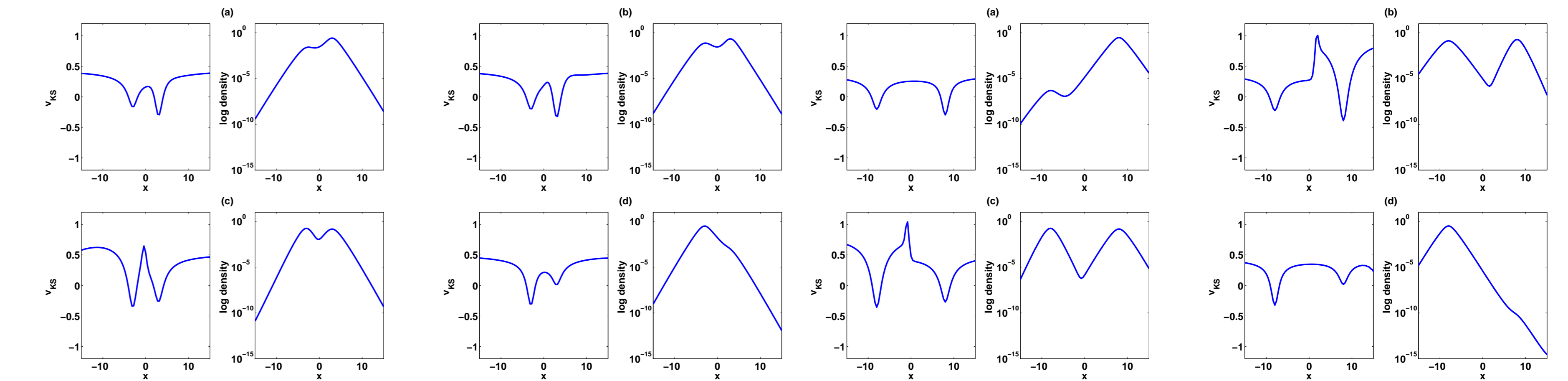
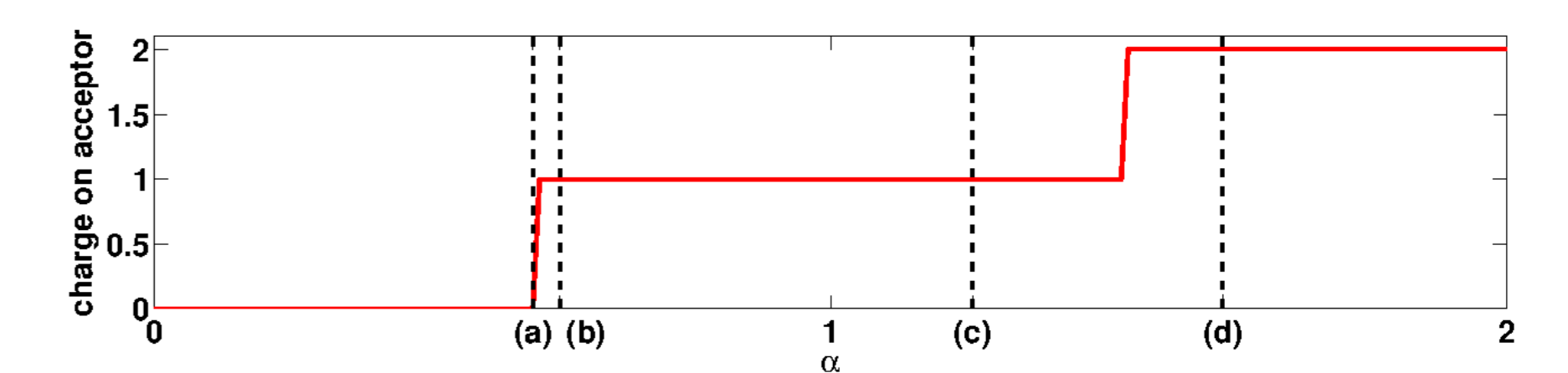
Exact solution of static two-electron Schrödinger equation with octopus (A. Castro et al. [4])

$$\hat{H}(\alpha) \Psi_j(\alpha) = E_j(\alpha) \Psi_j(\alpha) \\ n(x) = \langle \Psi | \hat{n}(x) | \Psi \rangle \\ \hat{n}(x) = \sum_j \delta(x - x_j)$$

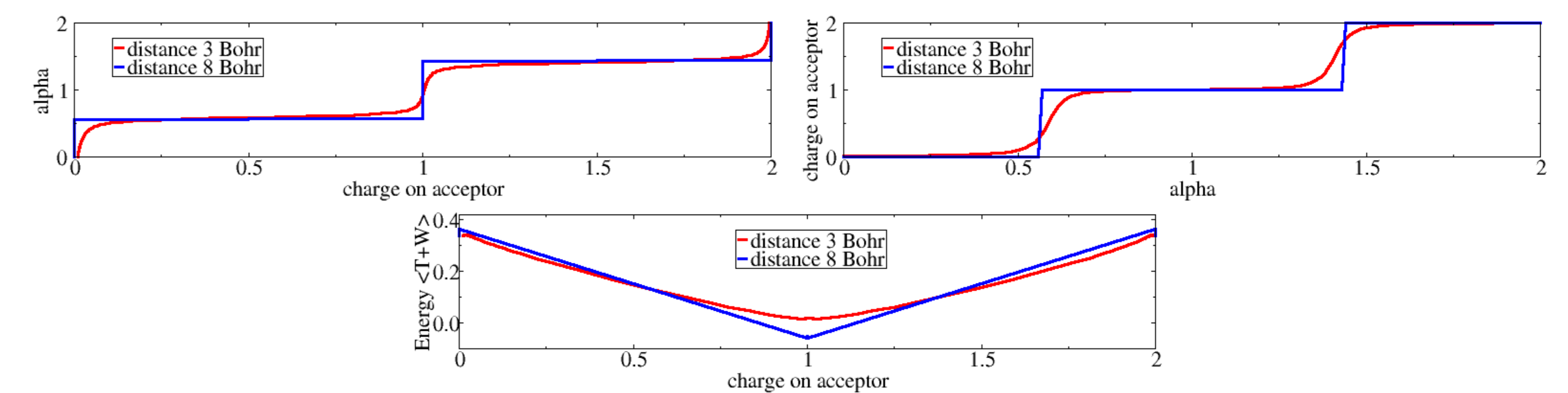
High-density limit (small distance $d = 3$ Bohr)



Low-density limit (large distance $d = 8$ Bohr)



Softened intra-system derivative discontinuity



Conclusion & Outlook

- The exact Hohenberg-Kohn functional shows a softened intra-system derivative discontinuity in the low-density limit.
- Expectation values of operators are affected by the softened intra-system derivative discontinuity.
- We observe softened intra-system derivative discontinuity also for soft-Coulomb molecules in 1D.
- We currently develop an approximate functional which incorporates the intra-system derivative discontinuity.

References

- M. Levy, Proc. Natl Acad. Sci. USA 76 6062 (1979)
- E. Lieb, Int. J. Quantum Chem. 24 24377 (1983)
- N. Helbig, I.V. Tokatly, A. Rubio, Journal of Chemical Physics 131, 224105 (2009)
- A. Castro, H. Appel, Micael Oliveira, C.A. Rozzi, X. Andrade, F. Lorenzen, M.A.L. Marques, E.K.U. Gross, and A. Rubio, Phys. Stat. Sol. B 243 2465-2488 (2006)
- John P. Perdew, Robert G. Parr, Mel Levy, Jose L. Balduz, Jr., Phys. Rev. Lett. 49, 16911694 (1982)