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Education

- 08.2007 – 04.2011 Dr. rer. nat, *summa cum laude*
Leibniz Universität Hannover, Germany
Institut für Festkörperphysik (Prof. Herbert Pfnür)
· Adsorption of organic molecules on wide band-gap insulators
- 07.2005 – 08.2007 Research assistant, Fudan University
Institute of Microelectronic
· High-k dielectrics in nonvolatile memory
- 09.2002 – 07.2005 MSc. in Electrical Engineering, Fudan University
Institute of Microelectronics
· DFT study on atomic layer deposition of high-k gate dielectrics
- 09.1998 – 07.2002 BSc., Fudan University, Shanghai, China

Research Experience

- 06.2016 – present Post-doctoral researcher, UCLouvain, Belgium
Institute of Condensed Matter and Nanosciences
Profs. Gian-Marco Rignanese and Geoffroy Hautier
· Multicomponent alloys
· Emerging kesterite solar absorbers
· Many-body perturbation theory
· Nonempirical hybrid density functional
· High-throughput computational screening
- 04.2011 – 03.2016 Post-doctoral researcher, Ecole Polytechnique Fédérale de Lausanne (EPFL)
Chair of Atomic Scale Simulation
Prof. Alfredo Pasquarello
· Defects in semiconductors and insulators
· Interfaces in semiconductor heterojunctions
· Advanced electronic-structure methods: *GW* approximation and hybrid functionals
· Electronic structure of liquid water: many-body and nuclear quantum effects

Coding skills

FORTRAN, C, PYTHON, GNU BASH, MPI.

Active developer for [ABINIT](#) and [QUANTUM ESPRESSO](#).

Code developments

ABINIT	Bootstrap exchange-correlation kernel for accurate <i>GW</i> quasiparticle energies, Gygi-Baldereschi auxiliary function for the treatment of Coulomb singularity
QUANTUM-ESPRESSO	Range-separated hybrid density functional
FNV	PYTHON class for finite-size-corrections of periodic charged defects (https://github.com/wayn3/FNV)

Teaching Activities

03.2019 – 06.2019	Masters course: Atomistic and Nanoscopic Simulations Teaching assistance, UCLouvain
04.2012 – 04.2016	Masters course: Computational Simulation and Physical Systems I & II Teaching assistance, EPFL
06.2013 – 04.2014	Masters thesis supervisor Project: Band-offset of lattice matched semiconductor heterojunctions (K. Steiner)

Recent Talks and Seminars

09.2021	E-MRS Fall Meeting 2021 Talk: “Origin of Low Conversion Efficiency of Cu ₂ ZnSnS ₄ Kesterite Solar Absorber”
05.2021	MRS Spring Meeting 2021 Talk: “Origin of Low Conversion Efficiency of Cu ₂ ZnSnS ₄ Kesterite Solar Absorber”
09.2018	Swiss Physics Society Annual Meeting 2018, Lausanne, Switzerland Invited talk: “Electronic structures through <i>GW</i> and hybrid functionals”
08.2017	29th International Conference on Defects in Semiconductors (ICDS), Matsue, Japan Invited talk: “Towards accurate determination of defect levels in semiconductors”
05.2017	ABINIT Developer Workshop 2017, Frèjus, France Invited talk: “Accurate band gaps via efficient vertex corrections in <i>GW</i> ”
10.2015	Université catholique de Louvain, Belgium Invited seminar: “Efficient vertex corrections in <i>GW</i> ”
09.2015	PSI-K 2015 conference, San Sebastian, Spain Talk: “Accurate band gaps via efficient vertex corrections in <i>GW</i> ”
07.2015	International Conference on Defects in Semiconductors (ICDS 15), Espoo, Finland Talk: “Determination of defect energy levels through <i>GW</i> ”
04.2015	“Nothing is Perfect” workshop, Ascona, Switzerland Invited Talk: “First-principles determination of defect energy levels through <i>GW</i> ”
08.2014	International conference on the physics of semiconductors (ICPS 14), Austin, USA Talk: “Band offset of lattice-matched semiconductor heterojunctions”
07.2013	International Conference on Defects in Semiconductors (ICDS 13), Bologna, Italy Talk: “Defect energy levels: Hybrid functionals vs <i>GW</i> ”

Awards

05.2021	Best oral presentation, MRS Spring Meeting 2021
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Miscellaneous

Referee for *Phys. Rev. Lett.*, *Phys. Rev. B*, *Appl. Phys. Lett.*, *J. Phys. Chem. Lett.*, *J. Phys. Chem.*, *npj Comput. Mater.*

Publications in Chronological Order

- [1] J. M. de Almeida, N. L. Nguyen, N. Colonna, W. Chen, C. Rodrigues Miranda, A. Pasquarello, and N. Marzari, “Electronic structure of water from koopmans-compliant functionals”, *J. Chem. Theory Comput.* **17**, 3923 (2021).
- [2] G. Bokas, W. Chen, A. Hilhorst, P. Jacques, S. Gorsse, and G. Hautier, “Unveiling the thermodynamic driving forces for high entropy alloys formation through big data ab initio analysis”, *Scr. Mater.* **202**, 114000 (2021).
- [3] W. Chen, D. Dahliah, G.-M. Rignanese, and G. Hautier, “Origin of the low conversion efficiency in Cu₂ZnSnS₄ kesterite solar cells: the actual role of cation disorder”, *Energy Environ. Sci.* **14**, 3567 (2021).
- [4] D. Evans, J. Chen, G. Bokas, W. Chen, G. Hautier, and W. Sun, “Visualizing temperature-dependent phase stability in high entropy alloys”, *npj Comput. Mater.* **7**, 151 (2021).
- [5] M. Markov et al., “Ferroelectricity and multiferroicity in anti-Ruddlesden-Popper structures”, *Proc. Natl. Acad. Sci.* **118**, e2026020118 (2021).
- [6] A. Tal, W. Chen, and A. Pasquarello, “Vertex function compliant with the Ward identity for quasiparticle self-consistent calculations beyond *GW*”, *Phys. Rev. B* **103**, L161104 (2021).
- [7] X. Gonze et al., “The ABINIT project: impact, environment and recent developments”, *Comput. Phys. Commun.* **248**, 107042 (2020).
- [8] T. Bischoff, J. Wiktor, W. Chen, and A. Pasquarello, “Nonempirical hybrid functionals for band gaps of inorganic metal-halide perovskites”, *Phys. Rev. Materials* **3**, 123802 (2019).
- [9] W. Chen, J. George, J. B. Varley, G.-M. Rignanese, and G. Hautier, “High-throughput computational discovery of In₂Mn₂O₇ as a high Curie temperature ferromagnetic semiconductor for spintronics”, *npj Comput. Mater.* **5**, 72 (2019).
- [10] S. Hadke, W. Chen, J. M. R. Tan, M. Guc, V. Izquierdo-Roca, G.-M. Rignanese, G. Hautier, and L. H. Wong, “Effect of Cd on cation redistribution and order-disorder transition in Cu₂(Zn, Cd)SnS₄”, *J. Mater. Chem. A* **7**, 26927 (2019).
- [11] L. A. Burton, F. Ricci, W. Chen, G.-M. Rignanese, and G. Hautier, “High-throughput identification of electrides from all known inorganic materials”, *Chem. Mater.* **30**, 7521 (2018).
- [12] W. Chen, G. Miceli, G.-M. Rignanese, and A. Pasquarello, “Nonempirical dielectric-dependent hybrid functional with range separation for semiconductors and insulators”, *Phys. Rev. Materials* **2**, 073803 (2018).
- [13] W. Chen and A. Pasquarello, “Comment on ‘Fundamental resolution of difficulties in the theory of charged point defects in semiconductors’”, *Phys. Rev. Lett.* **120**, 039603 (2018).
- [14] Z. Guo, F. Ambrosio, W. Chen, P. Gono, and A. Pasquarello, “Alignment of redox levels at semiconductor–water interfaces”, *Chem. Mater.* **30**, 94 (2018).
- [15] G. Miceli, W. Chen, I. Reshetnyak, and A. Pasquarello, “Nonempirical hybrid functionals for band gaps and polaronic distortions in solids”, *Phys. Rev. B* **97**, 121112(R) (2018).
- [16] W. Chen and A. Pasquarello, “Accuracy of *GW* for calculating defect energy levels in solids”, *Phys. Rev. B* **96**, 020101(R) (2017).
- [17] A. Faghaninia et al., “A computational assessment of the electronic, thermoelectric, and defect properties of bournonite (CuPbSbS₃) and related substitutions”, *Phys. Chem. Chem. Phys.* **19**, 6743 (2017).

- [18] W. Chen, F. Ambrosio, G. Miceli, and A. Pasquarello, “*Ab initio* electronic structure of liquid water”, *Phys. Rev. Lett.* **117**, 186401 (2016).
- [19] W. Chen and A. Pasquarello, “Accurate band gaps of extended systems via efficient vertex corrections in *GW*”, *Phys. Rev. B* **92**, 041115(R) (2015).
- [20] W. Chen and A. Pasquarello, “First-principles determination of defect energy levels through hybrid density functionals and *GW*”, *J. Phys.: Condens. Matter* **27**, 133202 (2015).
- [21] F. A. Rasmussen and K. S. Thygesen, “Computational 2d materials database: electronic structure of transition-metal dichalcogenides and oxides”, *The Journal of Physical Chemistry C* **119**, 13169 (2015).
- [22] W. Chen and A. Pasquarello, “Band-edge positions in *GW*: Effects of starting point and self-consistency”, *Phys. Rev. B* **90**, 165133 (2014).
- [23] K. Steiner, W. Chen, and A. Pasquarello, “Band offsets of lattice-matched semiconductor heterojunctions through hybrid functionals and G_0W_0 ”, *Phys. Rev. B* **89**, 205309 (2014).
- [24] W. Chen and A. Pasquarello, “Correspondence of defect energy levels in hybrid density functional theory and many-body perturbation theory”, *Phys. Rev. B* **88**, 115104 (2013).
- [25] W. Chen and A. Pasquarello, “Band-edge levels in semiconductors and insulators: Hybrid density functional theory versus many-body perturbation theory”, *Phys. Rev. B* **86**, 035134 (2012).
- [26] W. Chen, C. Tegenkamp, H. Pfnür, and T. Bredow, “Anomalous molecular orbital variation upon adsorption on a wide band gap insulator”, *J. Chem. Phys.* **132**, 214706 (2010).
- [27] W. Chen, C. Tegenkamp, H. Pfnür, and T. Bredow, “Color centers in NaCl by hybrid functionals”, *Phys. Rev. B* **82**, 104106 (2010).
- [28] W. Chen, C. Tegenkamp, H. Pfnür, and T. Bredow, “Insight from first-principles calculations into the interactions between hydroxybenzoic acids and alkali chloride surfaces”, *J. Phys. Chem. C* **114**, 460 (2010).
- [29] W. Chen, C. Tegenkamp, H. Pfnür, and T. Bredow, “Tailoring band gaps of insulators by adsorption at surface defects: Benzoic acids on NaCl surfaces”, *Phys. Rev. B* **79**, 235419 (2009).
- [30] W. Chen, C. Tegenkamp, H. Pfnür, and T. Bredow, “The interplay of van der Waals and weak chemical forces in the adsorption of salicylic acid on NaCl(001)”, *Phys. Chem. Chem. Phys.* **11**, 9337 (2009).
- [31] M. Zhang, W. Chen, S.-J. Ding, Z.-Y. Liu, Y. Huang, Z.-W. Liao, and D. W. Zhang, “Physical and electrical characterization of atomic-layer-deposited Ru nanocrystals embedded into Al₂O₃ for memory applications”, *J. Phys. D: Appl. Phys.* **41**, 032007 (2008).
- [32] W. Chen, W.-J. Liu, M. Zhang, S.-J. Ding, D. W. Zhang, and M.-F. Li, “Multistacked Al₂O₃-HfO₂-SiO₂ tunnel layer for high-density nonvolatile memory application”, *Appl. Phys. Lett.* **91**, 022908 (2007).
- [33] W. Chen, Q.-Q. Sun, M. Xu, S.-J. Ding, D. W. Zhang, and L.-K. Wang, “Atomic layer deposition of hafnium oxide from tetrakis-(ethylmethylamino)-hafnium and water precursors”, *J. Phys. Chem. C* **111**, 6495 (2007).
- [34] W. Chen, M. Zhang, D. Zhang, S.-J. Ding, J.-J. Tan, M. Xu, X.-P. Qu, and L.-K. Wang, “Growth of high-density Ru- and RuO₂-composite nanodots on atomic-layer-deposited Al₂O₃ film”, *Appl. Surf. Sci.* **253**, 4045 (2007).
- [35] S.-J. Ding, M. Zhang, W. Chen, D. W. Zhang, and L.-K. Wang, “Memory effect of metal-insulator-silicon capacitor with HfO₂-Al₂O₃ multilayer and hafnium nitride gate”, *J. Electron. Mater.* **36**, 253 (2007).
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- [38] Q.-Q. Sun, W. Chen, S.-J. Ding, M. Xu, H.-L. Lu, H.-C. Lindh-Rengifo, D. W. Zhang, and L.-K. Wang, “Comparative study of passivation mechanism of oxygen vacancy with fluorine in HfO_2 and HfSiO_4 ”, *Appl. Phys. Lett.* **90**, 142904 (2007).
- [39] Q.-Q. Sun, W. Chen, S.-J. Ding, M. Xu, D. W. Zhang, and L.-K. Wang, “Effects of chlorine residue in atomic layer deposition hafnium oxide: A density-functional-theory study”, *Appl. Phys. Lett.* **91**, 022901 (2007).
- [40] M. Zhang, W. Chen, S.-J. Ding, X.-P. Wang, D. W. Zhang, and L.-K. Wang, “Investigation of atomic-layer-deposited ruthenium nanocrystal growth on SiO_2 and Al_2O_3 films”, *J. Vac. Sci. Technol., A* **25**, 775 (2007).
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- [44] H.-L. Lu, W. Chen, S.-J. Ding, M. Xu, D. W. Zhang, and L.-K. Wang, “Quantum chemical study of the initial surface reactions in atomic layer deposition of TiN on the SiO_2 surface”, *J. Phys.: Condens. Matter* **18**, 5937 (2006).
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- [47] J. Ren, H.-L. Lu, W. Chen, M. Xu, and D. W. Zhang, “Surface reaction mechanism of atomic layer deposition of HfO_2 on $\text{Ge}(100)$ - 2×1 : A density functional theory study”, *Appl. Surf. Sci.* **252**, 8466 (2006).
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- [50] W. Chen, D. W. Zhang, J. Ren, H.-L. Lu, J.-Y. Zhang, M. Xu, J.-T. Wang, and L.-K. Wang, “Density functional theory study of initial stage of ZrO_2 atomic layer deposition on $\text{Ge/Si}(100)$ -(2×1) surface”, *Thin Solid Films* **479**, 73 (2005).
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