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Anställning

Biträdande universitetslektor
Matematisk fysik
Lunds universitet
Lund, Sverige
2021 feb. 9 → present

Profilområdesmedlem

LTH profilområde: Nanovetenskap och halvledarteknologi
Lunds universitet
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2022 aug. 26 → present

Principal Investigator

eSSENCE: The e-Science Collaboration
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2023 jan. 25 → present

Principal Investigator

NanoLund: Centre for Nanoscience
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2023 juni 2 → present

Profilområdesmedlem

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2023 okt. 16 → present

Forskningsoutput

Downfolding from ab initio to interacting model Hamiltonians: comprehensive analysis and benchmarking of the DFT+cRPA approach

Chang, Y., van Loon, E. G. C. P., Eskridge, B., Busemeyer, B., Morales, M. A., Dreyer, C. E., Millis, A. J., Zhang, S., Wehling, T. O., Wagner, L. K. & Rösner, M., 2024 dec., I: npj Computational Materials. 10, 1, 129.

Nb₃Cl₈: a prototypical layered Mott-Hubbard insulator

Grytsiuk, S., Katsnelson, M. I., Loon, E. G. C. P. V. & Rösner, M., 2024 dec., I: npj Quantum Materials. 9, 1, 8.

Second-order phase transitions and divergent linear response in dynamical mean-field theory

Van Loon, E. G. C. P., 2024 juni 15, I: Physical Review B. 109, 24, L241110.

Dual Bethe-Salpeter equation for the multiorbital lattice susceptibility within dynamical mean-field theory

Van Loon, E. G. C. P. & Strand, H. U. R., 2024 apr., I: Physical Review B. 109, 15, 155157.

Unconventional Charge-Density-Wave Gap in Monolayer NbS₂

Knispel, T., Berges, J., Schobert, A., van Loon, E. G. C. P., Jolie, W., Wehling, T. O., Michely, T. & Fischer, J., 2024 jan. 31, I: Nano Letters. 24, 4, s. 1045-1051 6 s.

Ab initio electron-lattice downfolding: Potential energy landscapes, anharmonicity, and molecular dynamics in charge density wave materials

Schobert, A., Berges, J., van Loon, E. G. C. P., Sentef, M. A., Brener, S., Rossi, M. & Wehling, T. O., 2024, I: SciPost Physics. 16, 2, 046.

Larmor precession in strongly correlated itinerant electron systems

van Loon, E. & Strand, H., 2023 okt. 13, I: Communications Physics. 6, 9 s., 289.

Coulomb engineering of two-dimensional Mott materials

van Loon, E. G. C. P., Schüler, M., Springer, D., Sangiovanni, G., Tomczak, J. M. & Wehling, T. O., 2023, I: npj 2D Materials and Applications. 7, 1, 47.

Two-particle correlations and the metal-insulator transition: Iterated perturbation theory revisited

Van Loon, E. G. C. P., 2022 juni 6, I: Physical Review B. 105, 24, 245104.

Degenerate plaquette physics as key ingredient of high-temperature superconductivity in cuprates

Danilov, M., van Loon, E. G. C. P., Brener, S., Iskakov, S., Katsnelson, M. I. & Lichtenstein, A. I., 2022, I: npj Quantum Materials. 7, 1, 50.

A full gap above the Fermi level: the charge density wave of monolayer VS2

Van Efferen, C., Berges, J., Hall, J., Van Loon, E., Kraus, S., Schobert, A., Wekking, T., Huttmann, F., Plaar, E., Rothenbach, N., Ollefs, K., Arruda, L. M., Brookes, N., Schönhoff, G., Kummer, K., Wende, H., Wehling, T. & Michely, T., 2021 dec. 1, I: Nature Communications. 12, 1

Downfolding the Su-Schrieffer-Heeger model

Schobert, A., Berges, J., Wehling, T. & van Loon, E., 2021 okt., I: SciPost Physics. 11, 4, 079.

Downfolding approaches to electron-ion coupling: Constrained density-functional perturbation theory for molecules

Van Loon, E. G. C. P., Berges, J. & Wehling, T. O., 2021, I: Physical Review B. 103, 20

Efficient fluctuation-exchange approach to low-temperature spin fluctuations and superconductivity: From the Hubbard model to $\text{Na}_x\text{CoO}_2\cdot y\text{H}_2\text{O}$
Witt, N., Van Loon, E. G. C. P., Nomoto, T., Arita, R. & Wehling, T. O., 2021, I: Physical Review B. 103, 20

Random phase approximation for gapped systems: Role of vertex corrections and applicability of the constrained random phase approximation
Van Loon, E. G. C. P., Rösner, M., Katsnelson, M. I. & Wehling, T. O., 2021, I: Physical Review B. 104, 4

Undervisning

2022	Grundläggande statistisk fysik och kvantstatistik (FYSB 23)
2021	Solid State Theory (FYST25)