

# Developing a Quantitative q-EELS Framework for Probing Collective Electron Dynamics

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## ABSTRACT

With a background in applied physics and materials science, I am motivated to develop experimental techniques that quantitatively probe how electrons respond to external perturbations. Momentum-dependent electron energy-loss spectroscopy (q-EELS) is particularly well suited for this purpose because it directly measures the loss function  $\text{Im}[-1/\epsilon(\omega, q)]$  and reveals the energy-momentum dispersion of collective excitations. This capability enables the extraction of key electronic parameters such as the effective mass  $m^*$ , Fermi velocity  $v_F$ , and carrier density  $n$ —quantities that fundamentally govern screening, transport, and dynamical behavior in solids. To establish a reliable analysis workflow, I first applied q-EELS to silicon as an illustrative benchmark, demonstrating that plasmon dispersion fitting yields  $m^*$  and  $v_F$  values consistent with the literature. In a separate example, I prepared and structurally verified monolayer and bilayer WSe<sub>2</sub> using substrate transfer and atomic-resolution STEM, providing representative low-dimensional samples for potential future studies. While these materials demonstrate the versatility of q-EELS, they are not the focus of the research itself; rather, they serve as practical test cases supporting the development of a broadly applicable technique. Moving forward, my primary objective is to refine q-EELS into a robust, general framework for extracting electronic dynamical parameters across diverse material systems, with optional extensions toward 2D materials when appropriate. Ultimately, I aim to establish a quantitative methodology that links atomic structure to collective electron dynamics through plasmon dispersion analysis.

## REFERENCES

1. Geim, A. K.; Grigorieva, I. V. *Van der Waals heterostructures*. Nature 2013, 499, 419–425.
2. Chou, S. A.; Chang, C.; Wu, B. H.; Chuu, C. P.; Kuo, P. C.; Pan, L. H.; et al. *Large-scale alkali-assisted growth of monolayer and bilayer WSe<sub>2</sub> with a low defect density*. Nature Communications 2025, 16, 2777.
3. Wang, I. T.; Chou, T. L.; Hsu, C. E.; Lei, Z.; Wang, L. M.; Lin, P. H.; et al. *The growing charge-density-wave order in CuTe lightens and speeds up electrons*. Nature Communications 2024, 15, 9345.
4. Shekhar, P.; Malac, M.; Gaiand, V.; Dalili, N.; Meldrum, A.; Jacob, Z. *Momentum-resolved electron energy loss spectroscopy for mapping the photonic density of states*. ACS Photonics 2017, 4, 1009–1014.
5. Nerl, H. C.; Guerrero-Felipe, J. P.; Valencia, A. M.; Elyas, K. F.; Höflich, K.; Koch, C. T.; Cocchi, C. *Mapping the energy-momentum dispersion of hBN excitons and hybrid plasmons in hBN–WSe<sub>2</sub> heterostructures*. npj 2D Materials and Applications 2024, 8, 68.