

Abstract Submitted
for the MAR10 Meeting of
The American Physical Society

Resonance Lifetimes from Complex Densities¹ DANIEL WHITENACK, Department of Physics, Purdue University, ADAM WASSERMAN, Department of Chemistry, Purdue University — The *ab-initio* calculation of resonance lifetimes of metastable anions challenges modern quantum-chemical methods. The exact lifetime of the lowest-energy resonance is encoded into a complex “density” that can be obtained via complex-coordinate scaling. Using one and two electron examples, we illustrate a method for extracting the lifetime from the complex density and explore a Kohn-Sham analog for resonances.

¹Support from the Petroleum Research Fund grant No.PRF # 49599-DNI6.

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Date submitted: 25 Nov 2009

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