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It lays out the foundations of quantum condensed phase dynamics: quantum mechanics, statistical mechanics, liquid theory, condensed matter theory, stochastic processes, and spectroscopy, and then builds upon these to present a clear and precise picture of how chemical physicists and physical chemists view chemical reactions.

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Using the hydrated electron as an example, application of this method finds that quantum decoherence times are on the order of a few femtoseconds for condensed phase chemical systems and that they play a direct role in determining nonadiabatic transition rates.

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Energy, charge, and chemical dynamics in condensed environments: Our research group investigates a variety of dynamic events in condensed media. The investigations range from basic chemical reaction dynamics in solutions such as ionic liquids to energy and charge transfer dynamics in new materials targeted at applications in solar energy conversion.

[David A. Blank | Department of Chemistry | College of](#)

In particular, our studies deal with chemical processes involving interactions between light and matter, chemical reactions in condensed phases and at interfaces and transport phenomena in complex systems, focusing mainly on the following directions: Energy transfer processes in molecular systems. Molecular dynamics in condensed phases.