

arXiv.org > physics > arXiv:1205.0296

Physics > Chemical Physics

Dephasing by a Continuous-Time Random Walk Process

Daniel M Packwood, Yoshitaka Tanimura

(Submitted on 2 May 2012)

Stochastic treatments of magnetic resonance spectroscopy and optical spectroscopy require evaluations of functions like <exp(i int_0^t Q(s)ds)>, where t is time, Q(s) is the value of a stochastic process at time s, and the angular brackets denote ensemble averaging. This paper gives an exact evaluation of these functions for the case where Q is a continuous-time random walk process. The continuous time random walk describes an environment that undergoes slow, step-like changes in time. It also has a well-defined Gaussian limit, and so allows for non-Gaussian and Gaussian stochastic dynamics to be studied within a single framework. We apply the results to extract qubit-lattice interaction parameters from dephasing data of P-doped Si semiconductors (data collected elsewhere), and to calculate the two-dimensional spectrum of a three level harmonic oscillator undergoing random frequency modulations.

Comments: 25 pages, 4 figures

Subjects: Chemical Physics (physics.chem-ph); Quantum Physics (quant-ph) Cite as: arXiv:1205.0296 [physics.chem-ph] (or arXiv:1205.0296v1 [physics.chem-ph] for this version)

Submission history

From: Daniel Packwood [view email] [v1] Wed, 2 May 2012 00:40:36 GMT (868kb)

Which authors of this paper are endorsers?

Link back to: arXiv, form interface, contact.

Search or Article-id

All papers 🚽 Go!

(Help | Advanced search)

Download:

• PDF only

Current browse context: physics.chem-ph < prev | next > new | recent | 1205

Change to browse by:

physics quant-ph

References & Citations

NASA ADS

Bookmark(what is this?)

