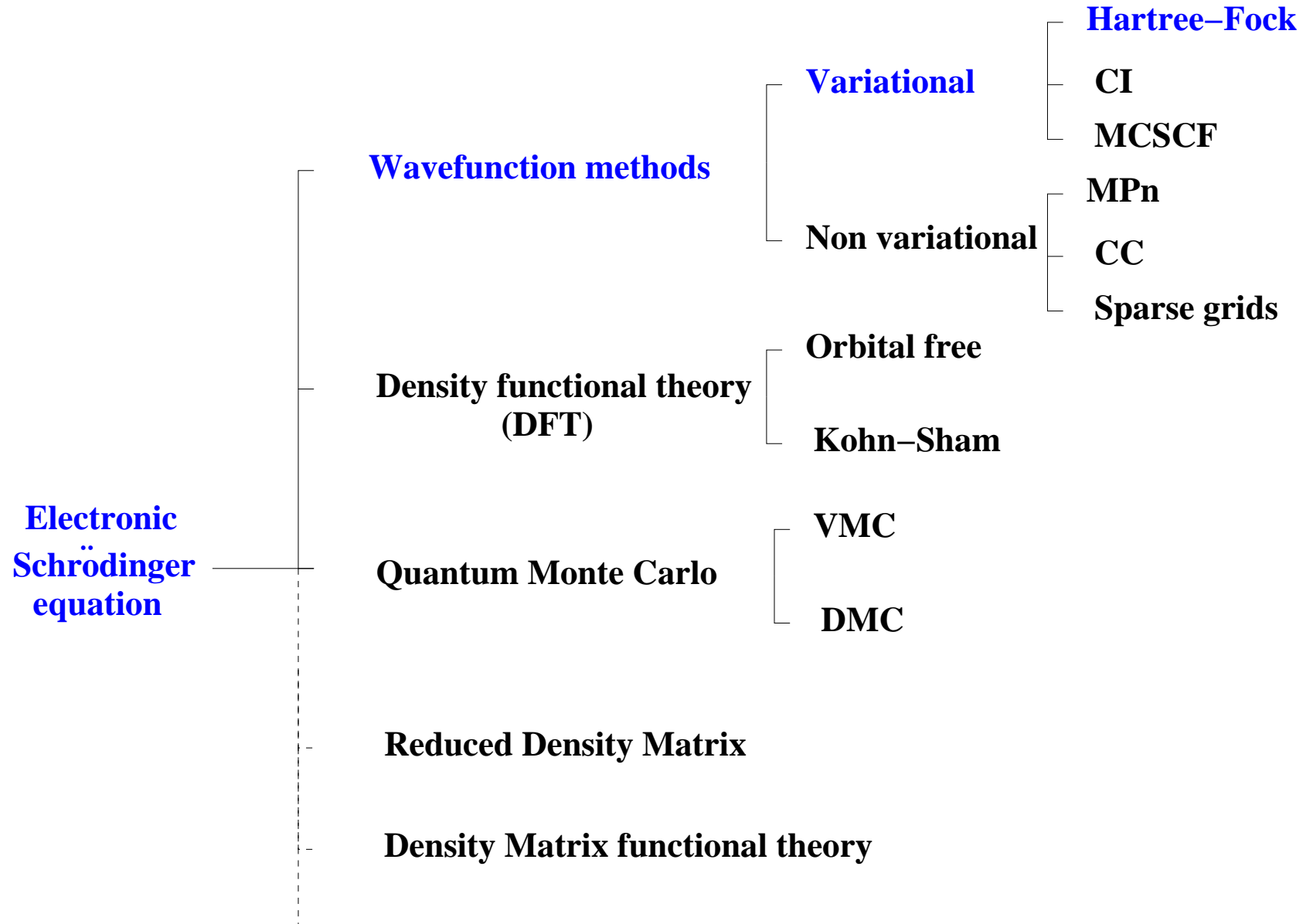


Mathematical modelling of electronic structures

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IMA, Minneapolis, September 2008



1 - Motivation: Ab initio molecular dynamics