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Bachelor Thesis in Physics
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Electric Field Optimization of a Rydberg Atom Experiment

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Abstract

Modern experiments with ultracold Rydberg atoms with application to many body physics and quantum information science, demand a high level of experimental sophistication to precisely control experimental parameters like external electric fields, as Rydberg atoms are very polarizable. In the experiment this is achieved by (a structure hosting) ~ 10 individually controllable electrodes. However, the task of finding the optimal control voltages for these is complicated by incomplete knowledge of the charge distributions, including possible patch fields (making it particularly time consuming). To overcome this challenge we have applied evolutionary algorithms, a group of powerful search heuristics, to optimize the overall performance of our experiment. With particular focus on electric field control we assess the performance of several algorithms, on competing requirements of noise robustness and fast convergence, in solving two problems: cancellation of electric fields and optimum guiding of field ionized Rydberg atoms to a MCP detector. Additionally Foreseeable applications to controlling quantum state evolution and engineering strongly correlated many body systems of interacting Rydberg atoms will be considered.

Introduction

Introduction introduction introduction

Contents

| | | |
|----------|--|----------|
| 1 | Concept | 3 |
| 2 | Theoretical Background | 4 |
| 3 | Optimization | 5 |
| 3.1 | Mathematical Formulation | 5 |
| 3.2 | Background: Why look at multiple algorithms? | 5 |
| 3.3 | Evolutionary Algorithms | 5 |
| 3.3.1 | Population Size | 6 |
| 3.3.2 | Exploration//Exploitation | 6 |
| 3.4 | Swarm Intelligence | 6 |
| 3.5 | Evolutionary Computation | 6 |
| 4 | Implementation | 7 |
| 4.1 | Test Function | 7 |
| 4.2 | Comparison of Algorithms | 7 |
| 4.3 | Appendix | 8 |

Chapter 1

Concept

The aim of this thesis is the application of optimization methods to a Rydberg atom experiment. In the course of making one measurement several steps have to be completed sequentially. These start with collecting Rubidium atoms into a MOT, pre-cooling then and then loading them into a dipole trap where they are excited into Rydberg atoms. After this has been done the atoms can be imaged by a camera system. These steps form a cycle that is repeated every few seconds with variation of parameters in order to make measurements of the change in behaviour of the Rydberg atoms depends on the varied parameters. While many properties of the experimental setup are physically fixed, such as the position of the lasers, a relatively large number (77) that can be controlled by the software controlling the experiment, such as the timing of activating these lasers. The aim of this project is to add information feedback to this cycle, so that the experiment can choose new experimental parameters based the results of previous evaluations. The schematic for this is shown in figure 1.1

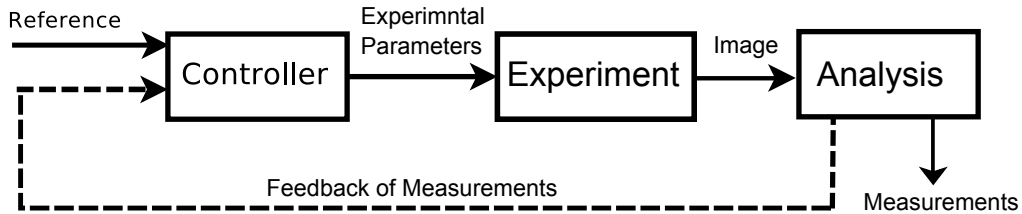


Figure 1.1: Information Feedback

Due to the discrete nature of experimental cycle evaluations and the large number of parameters as well as considerable measurement noise an optimization approach was chosen that generates and evaluates new experimental parameters. Using one can define the goal of the optimization as a function of several of the measurement results calculated by the analysis step.

Of particular interest is the control of experimental electrodes surrounding the Rydberg atoms. One of the experimental goals is creating minimal electric fields as the predictions that are made assume a field-free environment. Additionally it is important to generate a homogeneous field as this prevents position dependent energies. When working with Rydberg atoms one mostly wants to increase the principal quantum number n in order to scale with it the physical effects one is trying to measure, however the polarizability of Rydberg atoms scales with n^7 requiring fine control over the electric field if one wants to work with larger Rydberg atoms.

Chapter 2

Theoretical Background

Chapter 3

Optimization

3.1 Mathematical Formulation

A n dimensional real value optimization problem can be stated in the form

$$\min f(\mathbf{x})$$

for

$$f : S \rightarrow \mathbb{R}$$

where

$$\mathbf{x} \in S \subseteq \mathbb{R}^n$$

A point $\mathbf{x}^* \in D$ is a global minimum if $f(\mathbf{x}^*) \leq f(\mathbf{x}) \forall \mathbf{x} \in S$

A point $\mathbf{x}^* \in D$ is a local minimum if in the surrounding $U \subseteq S$ of \mathbf{x}^* it holds that $f(\mathbf{x}^*) \leq f(\mathbf{x}) \forall \mathbf{x} \in U$

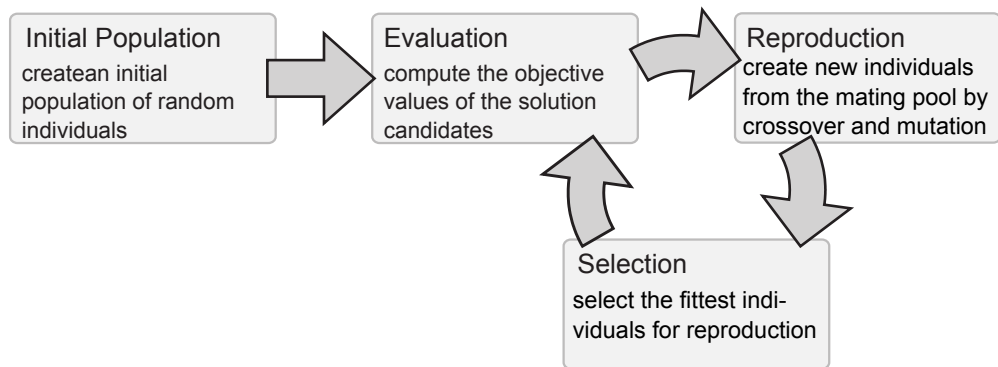
3.2 Background: Why look at multiple algorithms?

There are a large number of optimization algorithms that have been developed in order to solve different problems. The key to applying optimization algorithms then is to find one suitable to solving kind of problems one is working with. Abstractly this is the converse to the No Free Lunch Theorem which states that the average performance of an optimization algorithms, when applied to the set of all optimization problems, does not outperform any other optimization algorithm or random walk. An algorithm is thus only, comparatively, suitable if it is able to utilize some form of problem specific information coupling the choice of algorithm to the problem at hand.

Due to the kind of problems to which optimization is to be applied to only the subset of Monte Carlo probabilistic global optimization algorithms will be considered. These algorithms sacrifice both evaluating the entire search space and the necessity of evaluating the function exactly in favour of a shorter runtime. In these algorithms the choice of which candidates to evaluate is made by a heuristic which makes an induction based on previous evaluations. It is in this heuristic that is represents the problem specific information of the optimization algorithm.

3.3 Evolutionary Algorithms

Evolutionary Computation represents a subset of heuristic based approaches in which a set of possible solution candidates is maintained which the algorithm tries to refine over a number of generations. In the following sections we will introduce the different categories of algorithms used as well as the specific implementations of algorithms from these categories.



3.3.1 Population Size

3.3.2 Exploration//Exploitation

3.4 Swarm Intelligence

3.5 Evolutionary Computation

Chapter 4

Implementation

The application of optimization directly to the Rydber experiment posed a number of difficulties in that the parameters for optimization on the physical experiment were very different from those which numerical optimization is usually applied to, while it is possible to evaluate a numerical function or simulation very quickly and or in parallel there was a strong constraint on the number of evaluations of solution candidates. In order to still be able to achieve convergence of the optimization algorithm this required a reduced dimensionality of the problem that we were attempting to solve. In addition to this physical measurements are also affected by noise. These conditions were used to choose an appropriate algorithm that could be applied to the experiment.

4.1 Test Function

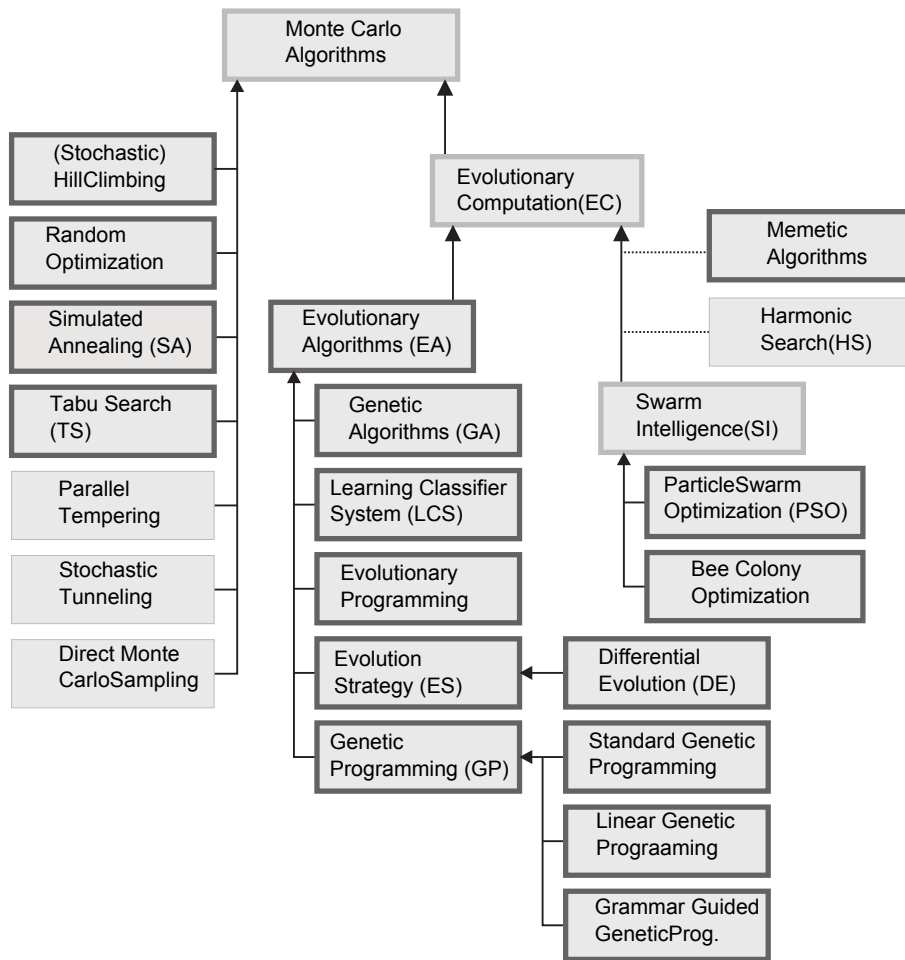
In order to evaluate the different algorithms that are available a mathematical test problem was formulated. This problem could then be evaluated by computer allowing for the large number of evaluations needed to compare the algorithm with different parameters and compare the results to other algorithms. In order to make this comparison meaningful one needs to choose a test problem that matches the actual experiment as closely as possible, to do this a n dimensional Gaussian was chosen.

To further increase similarity with the experiment three sources of noise were added: measurement noise affecting the amplitude of the function(σ_a), background noise added to the function(σ_b) and noise in the parameters that were passed to the function σ_p .

$$g(x_i, \sigma) = e^{-\frac{(x_i^* - x_i)^2}{2\sigma}}$$
$$G(\mathbf{x}, \sigma) = \prod_{i=1}^n g(x_i, \sigma)$$
$$f(\mathbf{x}) = \mathcal{N}(1, \sigma_a) * \prod_{i=1}^n g(\mathcal{N}(x_i, \sigma_p), \sigma) + \mathcal{N}(0, \sigma_b)$$

As we know only little about the problem landscape of the actual experimental parameters that we are going to optimize this test problem represents a best guess, strictly speaking we cannot induce the suitability to application in the experiment of the algorithm chosen according to this test function, however it will be attempted as nonetheless as it is the best starting point available.

4.2 Comparison of Algorithms



4.3 Appendix