

Abstract:

Radio frequency (RF) surface electrode traps where all the RF and static control electrodes lie in a single plane are being developed for scaling ion trap quantum information processing to larger numbers of ions [1]. In such traps, ions are confined a small distance above the electrode surface. The RF ponderomotive potential is anharmonic, rising steeply for ion excursions towards the electrode surface but much more gradually for excursions away from the surface. Multipole RF traps can exhibit a significant reduction in the effective trapping depth from the calculated psuedo-potential well depth due to non-adiabatic ion motion in the anharmonic potential [2]. We simulate whether anharmonic contributions to surface electrode trap potentials can reduce the effective well depth of surface electrode traps. Specifically we simulate the motion of a charged particle in a 4wire surface electrode trap. By starting the particle in the center of the trap with different energies we determine a safe or effective well depth as a function of the q parameter of the trap. We find significant reduction in the effective well depth for q > 0.3, resulting in a maximum well depth for q < 0.3.

1. Microfabricated Chip Traps for Ions, J. M. Amini, J. Britton, D. Leibfried, and D.J. Wineland. Chapter in the upcoming book "Atom chips" edited by J. Reichel and V. Vuletic (to be published by WILEY-VCH); J.H. Wesenberg, Phys. Rev. A 78, 063410 (2008).

2. J. Mikosh et al., Phys. Rev. Lett. 98, 223001 (2007).





RF amplitude V [V]

Simulation of non-Adiabaticity in Surface Electrode Traps

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Can adiabaticity limit the effective well depth of surface electrode traps?



Definition of adiabaticity

Dieter Gerlich, Advances in Chemical Physics: State-Selected and State-to-State Ion-Molecule reaction Dynamics, Part 1, Experiment, Volume 82, 1992

defines adiabaticity as validity of expansion:

 $\vec{E}_{rf}(\vec{R}_0 - \vec{a}\cos(\Omega t)) \approx \vec{E}_{rf}(\vec{R}_0) - (\vec{a} \cdot \vec{\nabla})\vec{E}_{rf}(\vec{R}_0)\cos(\Omega t)$ where ion motion $\vec{r}(t) = \vec{R}_0(t) - \vec{a}(t)\cos(\Omega t)$ $\vec{a}(t) = \frac{eE_{rf}}{m\Omega^2}$ (the micro - motion)

term in expansion is small when adiabaticity parameter $\eta <<1$

 $\eta \equiv \frac{2e\left|\vec{\nabla}E_{rf}\right|}{m\Omega^2}, E_{rf} \equiv \left|\vec{E}_{rf}\right|, \text{ safe confinement for } \eta < 0.3$

for quadrupole trap $\eta = q$; for higher order multipoles, $\eta = \eta(r)$

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An interesting example:

²⁴Mg⁺ ion $\frac{\Omega}{2\pi} = 87 \text{ MHz}$ $V_{rf} = 103.2 \text{ V}$ $d = 40 \ \mu m$ $q_{4W} = \frac{2eV_{rf}}{m\Omega^2(\pi d^2)} \approx 0.55$

psuedo - potential well depth $E_{well} \approx (3.6 \times 10^{-3})q_{4W}V_{rf} \sim 204 \text{ meV}$ simulated adiabatic well depth $E_{safe} \sim 20 \text{ meV}$

Simulation of reduced well depth in a 4-wire trap (preliminary results)

Procedure for determining E_{safe}

•Set an initial guess for the energy E •Set the initial positions (x_i, y_i) at the center of the trap •Calculate the magnitude of initial velocity $v_0 = \sqrt{2E_i/m}$ •Vary the angle ψ of the velocity from 0 to $15\pi/16$ with an interval of $\pi/16$: $v_{xi} = v_0 \cos \psi$; $v_{yi} = v_0 \sin \psi$ • Choose the initial phase for the electric field ϕ from 0 to $15\pi/16$ with an interval of $\pi/16$ •Update the position, velocity and energy at each time step by solving the equations of motion

•If the energy is larger than full well depth, the particle is unbounded, go back to the very beginning with a slightly lower E_i , (i.e. 10^{-5} eV) •If the ion trajectory is trapped for a sufficiently long time (10⁵ rf cycles, i.e. more than 1 ms), we assume

this initial condition gives a trapped orbit •Repeat the above steps for all the initial conditions (16 velocity angles Ψ and 16 electric field phases ϕ) •For each initial condition, record the boundary energy to safely confine the trap

•From these boundary energies, choose the minimal value as E_{safe}

Numerical Simulation







the reduction in effective well depth as expected from adiabaticity theory, but also at a fine level of detail a "periodic" dependence which we currently do not have a theory for.



These simulations involving a single ion show the importance of non-adiabatic effects is surface electrode traps. Simulations with two ions can investigate whether the Coulomb interaction can amplify non-adiabatic effects. For example, experiments have observed significantly shorter lifetimes with two ions than a single ion. Simulations can investigate whether this is due to non-adiabatic effects and whether these effects can be minimized with particular trap operating parameters.



Seidelin et al., PRL 96 (2006)



 E_{safe} as a function of $q \equiv \eta(0, d)$: (a) vary V_{rf} ; (b) vary $1/md^2\Omega^2$. The simulation clearly shows