

# Ion Coulomb Crystals



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Physics with Trapped Charged Particles  
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# Outline:

- I. Conditions for Coulomb crystallization
- II. Structural properties under various trapping conditions
- III. Metastable long-range-ordered Coulomb clusters
- IV. Control of crystal structures by light induced forces

# **I. Conditions for Coulomb crystallization**

# Conditions for Coulomb crystallization

## Condition for the formation of Coulomb crystals

### 3D, infinite case:

$$\Gamma = E_{\text{Coul}}/E_{\text{kin}} = Q^2/(4\pi\epsilon_0 a k_B T) > \sim 170 \quad ,$$

classical  
↙

or equivalent

$$a/\Delta r_{\text{kin}} > \sim 10$$

with  $a = (3/(4\pi n))^{1/3}$  is the Wigner-Seitz radius

J.-P. Hansen, Phys. Rev. A **8**, 3096 (1973)

### 3D, finite case:

$$\Gamma = \sim 200 \text{ (~10.000 ions)} - 500 \text{ (~100 ions)}$$

J. P Schiffer, PRL **88**, 205003(2002)

# Historical background

## Wigner 1934:

Electrons in metals at low temperatures:

Normally,  $\Delta r \sim a$  due to quantum mechanics.

By *lowering* the electron density one would, however, get Coulomb crystals or so-called

**Wigner-crystals.**

Still not observed in 3D, but....

~1980:

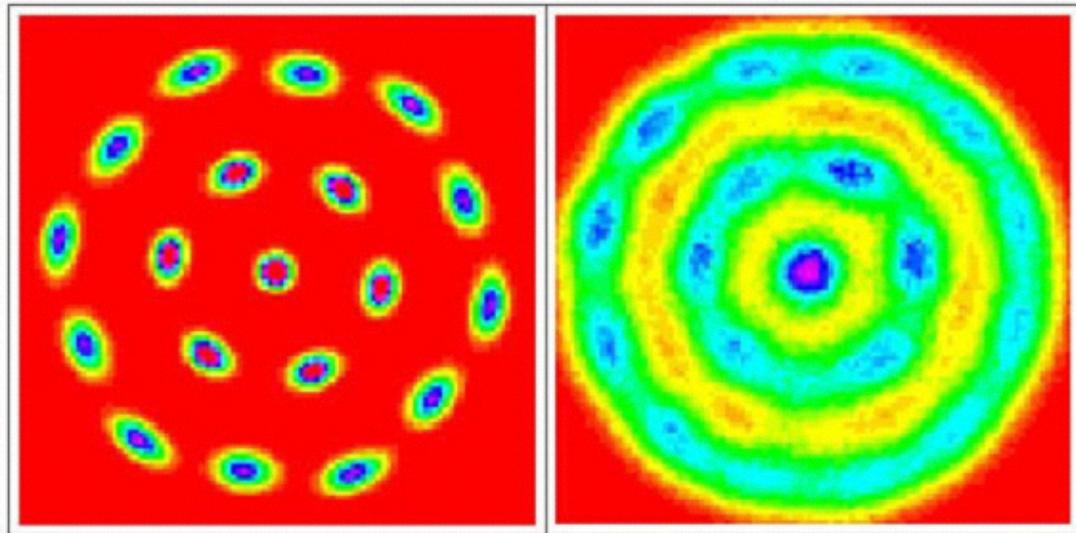
2D Wigner-crystals on the surface  
of superfluid helium

~1990:

2D Wigner-crystals in GaAs/GaAlAs  
quantum well structures

~2000:

Simulations of few electrons in a quantum dot:



# What about ions?

~1980:

Dense stellar objects: White Dwarfs and Neutron Stars

Criterion for avoiding “quantum melting”:

$$a > 10^4 \times h^2 \epsilon_0 / (Q^2 m)$$

Interior of WD:  $C^{6+}, O^{8+} \Rightarrow a > \sim 10^{-12} \text{ m}$

Crust of NS:  $Fe^{26+} \Rightarrow a > \sim 10^{-14} \text{ m}$

I.e., at sufficient “low” temperatures ( $\sim 10^6 \text{ K}$ ) ion Coulomb crystals are expected to be present in these exotic objects !

# Trap experiments

<1980:

Trap ion densities of  $n \sim 10^8 \text{ cm}^{-3}$  and temperatures of  $>1 \text{ K}$

$$\Rightarrow \Gamma < 1$$

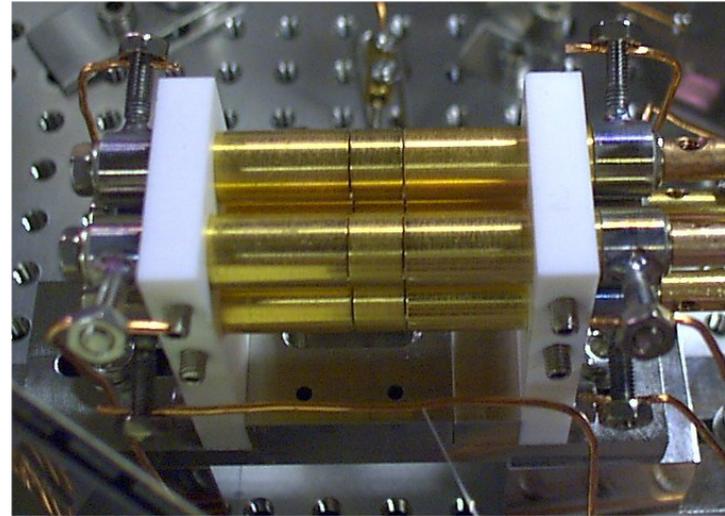
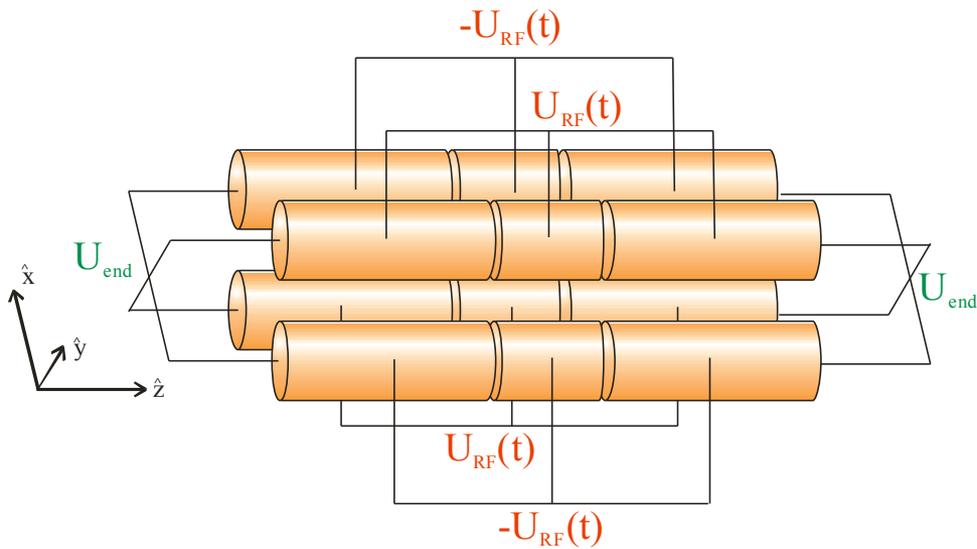
~1980:

First demonstration of laser cooling of ions. (NIST, Hamburg)

$$\Rightarrow \Gamma \sim 1000$$

~1980 ->: Non-neutral plasma physics, statistical physics, non-linear dynamics, metrology, quantum optics, cold molecules, nuclear physics, solid state physics...

# The linear Paul trap



Sinusoidal RF potential:  $U_{RF}(t) = U_{RF} \sin(\Omega t)$

Effective oscillation freq.'s:

$$\omega_r = 1/2 \beta \Omega, \quad \beta = (1/2 q^2 + a)^{1/2}$$

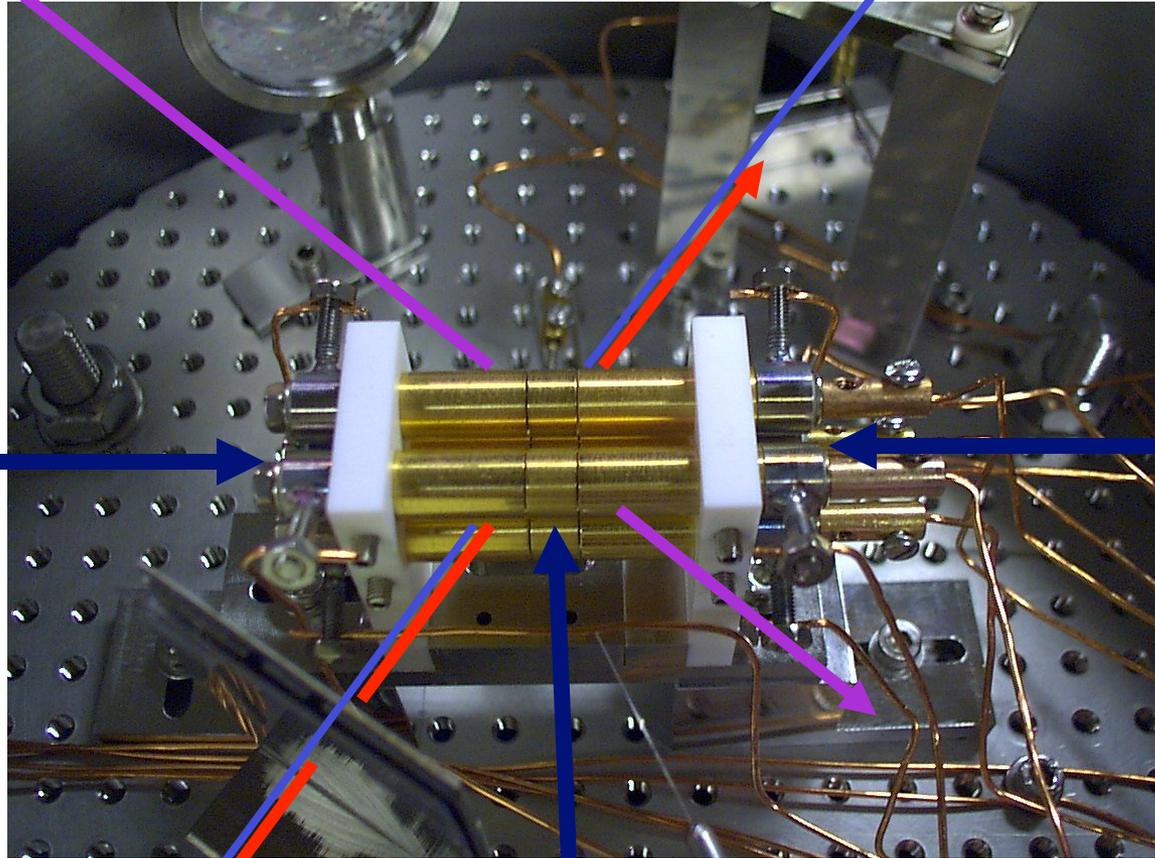
$$\omega_z = (-1/2 a)^{1/2} \Omega$$

$$q = \frac{4Q U_{RF}}{m \Omega^2 r_0^2} \quad a = - \frac{\alpha Q U_{end}}{m \Omega^2 r_0^2}$$

# The Aarhus linear Paul trap

Photo-ionizing laser

Electron beam



Cooling laser

Cooling laser

Atomic beam

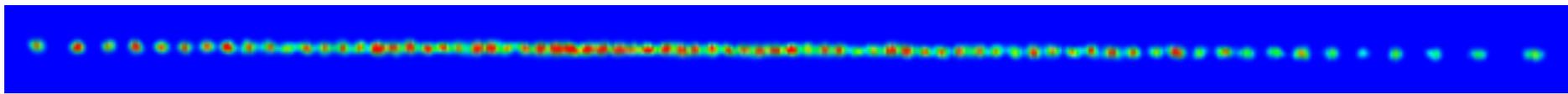
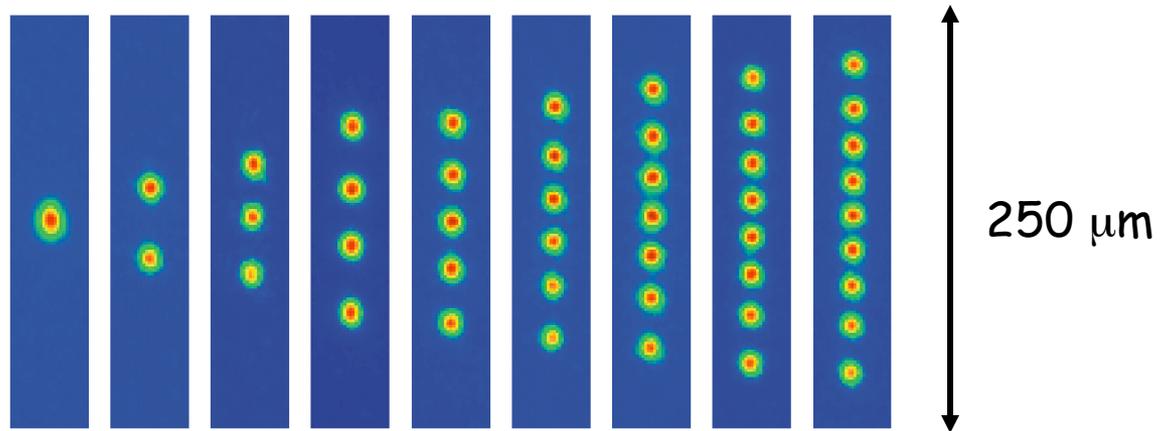
Cooling laser

5 cm

## **II. Structural properties under various trapping conditions**

# Single species ion Coulomb crystals

Smaller crystals - strings of ions



1 mm

Ion spacing:  $\sim 10 \mu\text{m}$

# How to determine the 1D linear ion structure?

Potential energy:

$$V = \sum_{m=1}^N \frac{1}{2} M v^2 x_m(t)^2 + \sum_{\substack{n,m=1 \\ m \neq n}}^N \frac{Z^2 e^2}{8\pi\epsilon_0} \frac{1}{|x_n(t) - x_m(t)|}$$

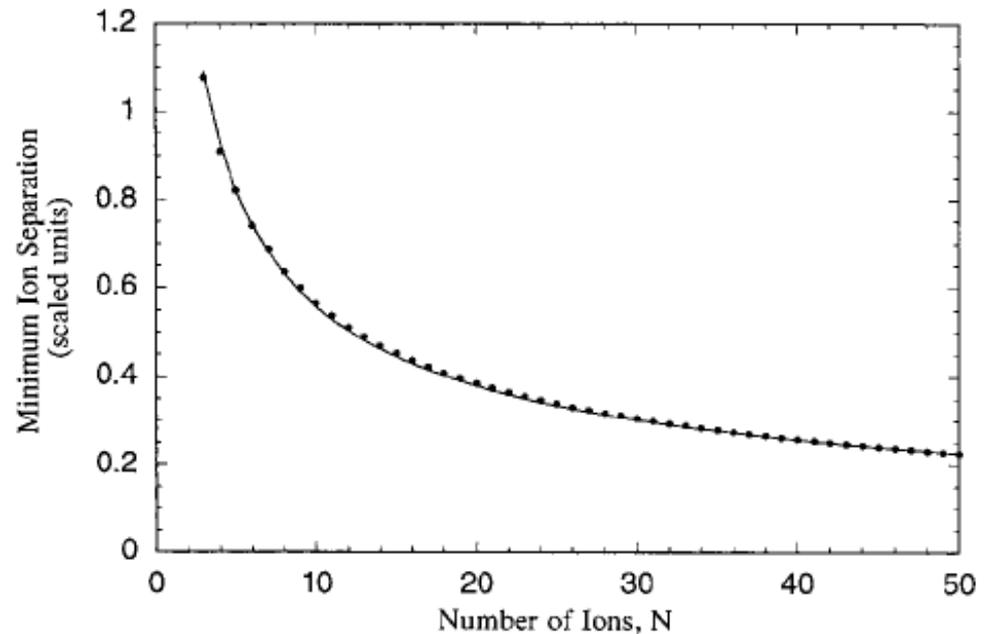
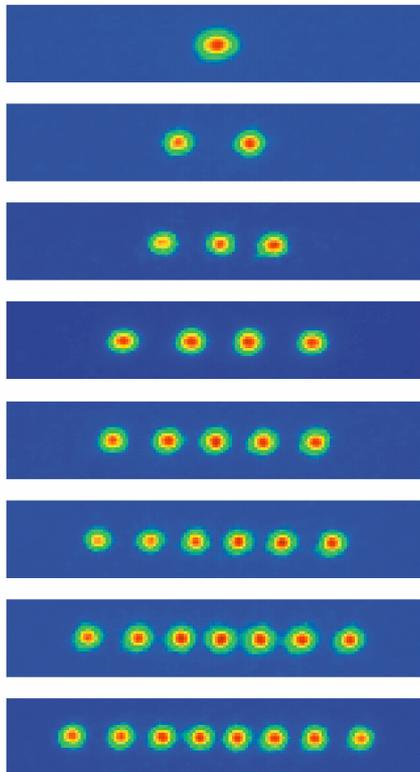
Equilibrium positions:

$$\left[ \frac{\partial V}{\partial x_m} \right]_{x_m = x_m^{(0)}} = 0$$

# Scaled equilibrium positions

N

2				-0.62996	0.62996								
3				-1.0772	0	1.0772							
4				-1.4368	-0.45438	0.45438	1.4368						
5				-1.7429	-0.8221	0	0.8221	1.7429					
6				-2.0123	-1.1361	-0.36992	0.36992	1.1361	2.0123				
7				-2.2545	-1.4129	-0.68694	0	0.68694	1.4129	2.2545			
8				-2.4758	-1.6621	-0.96701	-0.31802	0.31802	0.96701	1.6621	2.4758		
9				-2.6803	-1.8897	-1.2195	-0.59958	0	0.59958	1.2195	1.8897	2.6803	
10				-2.8708	-2.10003	-1.4504	-0.85378	-0.2821	0.2821	0.85378	1.4504	2.10003	2.8708



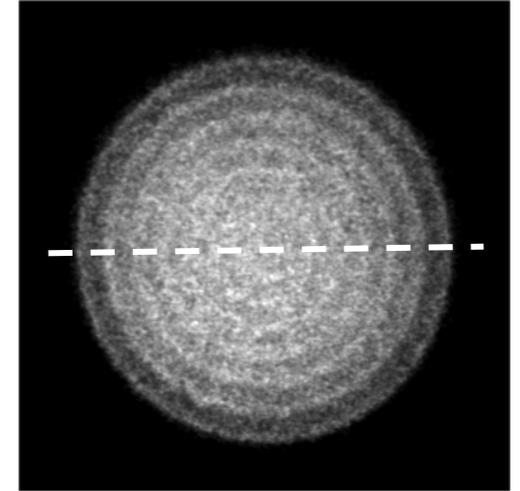
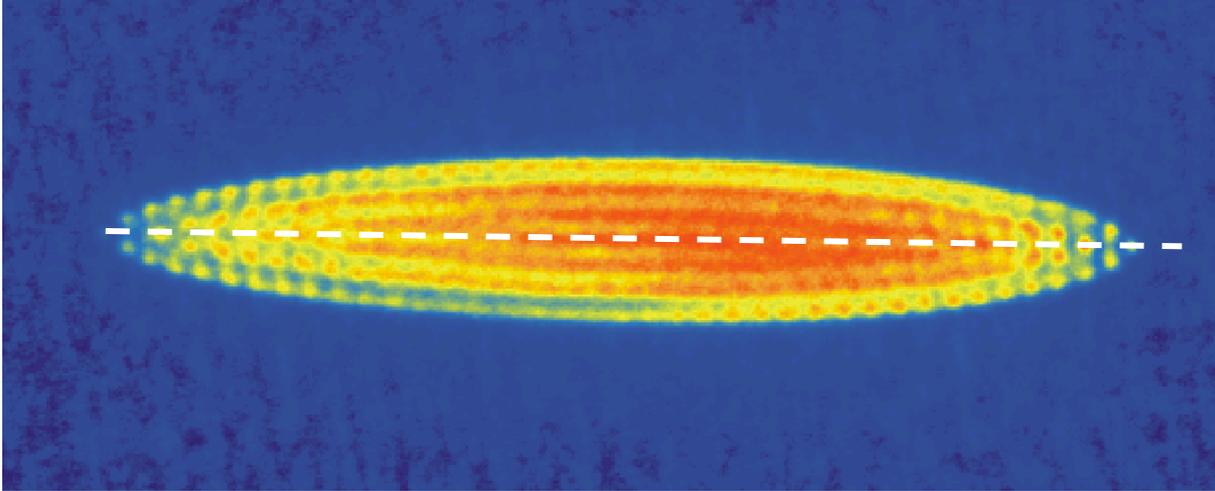
# Normal mode frequencies and eigenvectors

#ions  $\omega_{N,i}^2$ :

---

N=2	1	( 0.7071,	0.7071)					
	3	(-0.7071,	0.7071)					
N=3	1	( 0.5774,	0.5774,	0.5774)				
	3	(-0.7071,	0,	0.7071)				
	5.8	( 0.4082,	-0.8165,	0.4082)				
N=4	1	( 0.5,	0.5,	0.5,	0.5)			
	3	(-0.6742,	-0.2132,	0.2132,	0.6742)			
	5.81	( 0.5,	-0.5,	-0.5,	0.5)			
	9.308	(-0.2132,	0.6742,	-0.6742,	0.2132)			
N=5	1	( 0.4472,	0.4472,	0.4472,	0.4472,	0.4472)		
	3	(-0.6395,	-0.3017,	0,	0.3017,	0.6395)		
	5.818	( 0.5377,	-0.2805,	-0.5143,	-0.2805,	0.5377)		
	9.332	(-0.3017,	0.6395,	0,	-0.6395,	0.3017)		
	13.47	( 0.1045,	-0.4704,	0.7318,	-0.4704,	0.1045)		
N=6	1	( 0.4082,	0.4082,	0.4082,	0.4082,	0.4082,	0.4082)	
	3	(-0.608,	-0.3433,	-0.1118,	0.1118,	0.3433,	0.608)	
	5.824	(-0.5531,	0.1332,	0.4199,	0.4199,	0.1332,	-0.5531)	
	9.352	( 0.3577,	-0.5431,	-0.2778,	0.2778,	0.5431,	-0.3577)	
	13.51	( 0.1655,	-0.5618,	0.3963,	0.3963,	-0.5618,	0.1655)	
	18.27	(-0.04902,	0.2954,	-0.6406,	0.6406,	-0.2954,	0.04902)	
N=7	1	( 0.378,	0.378,	0.378,	0.378,	0.378,	0.378,	0.378)
	3	(-0.5801,	-0.3636,	-0.1768,	0,	0.1768,	0.3636,	0.5801)
	5.829	(-0.5579,	0.031,	0.3213,	0.4111,	0.3213,	0.031,	-0.5579)
	9.369	(-0.3952,	0.445,	0.3818,	0,	-0.3818,	-0.445,	0.3952)
	13.55	(-0.213,	0.5714,	-0.1199,	-0.4769,	-0.1199,	0.5714,	-0.213)
	18.32	( 0.08508,	-0.4121,	0.5683,	0,	-0.5683,	0.4121,	-0.08508)
	23.66	( 0.02222,	-0.1723,	0.4894,	-0.6787,	0.4894,	-0.1723,	0.02222)

# Medium size crystals ( $\sim 100$ - $\sim 1000$ ions)



## Properties:

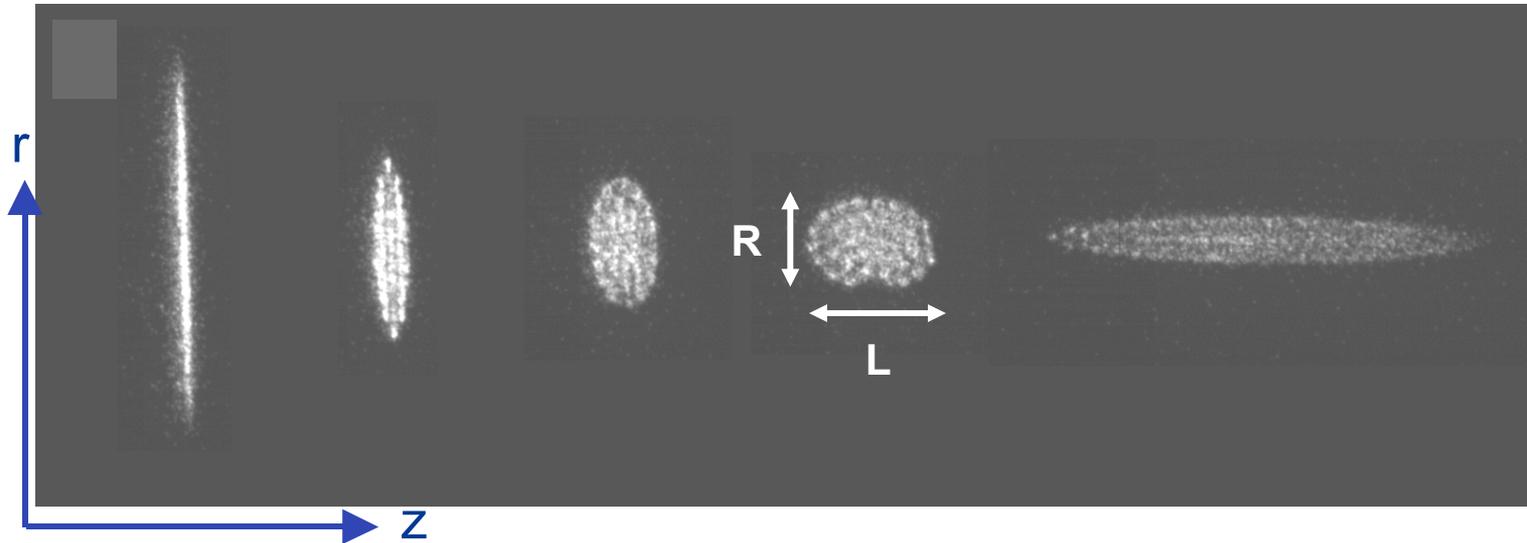
Shells with near-hexagonal structures

Uniform density of  $\sim 10^8 \text{ cm}^{-3}$

Temperature of  $\sim 10 \text{ mK}$

Trap life times of hours @  $P \sim 10^{-10} \text{ mBar}$

# Outer contours of crystals



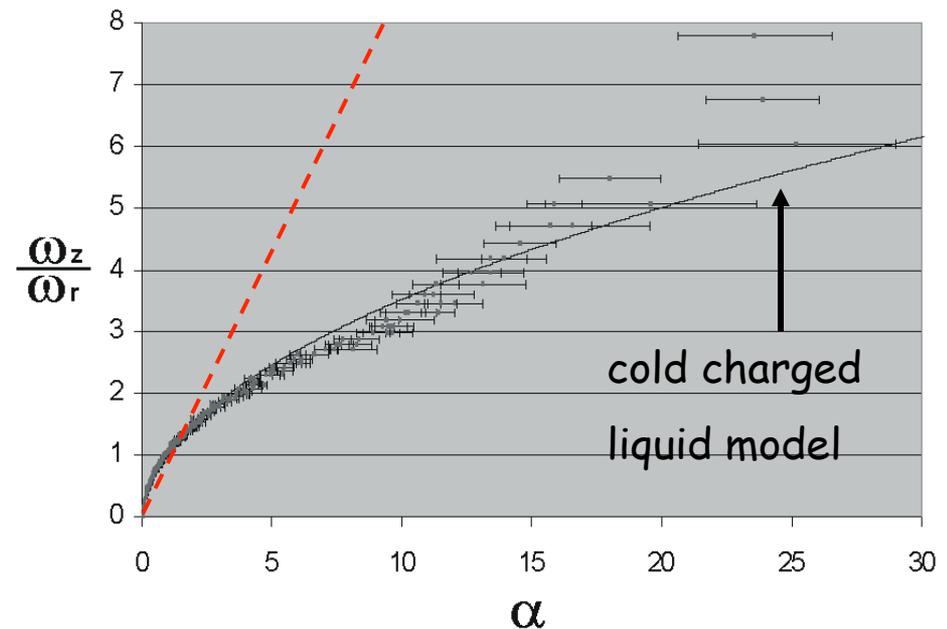
## Cold uniformly charged liquid model

In the oblate case,  $\alpha = R/L > 1$ ,

$$\frac{\omega_z^2}{\omega_r^2} = -2 \frac{\sin^{-1}(1 - \alpha^{-2})^{1/2} - \alpha(1 - \alpha^{-2})^{1/2}}{\sin^{-1}(1 - \alpha^{-2})^{1/2} - \alpha^{-1}(1 - \alpha^{-2})^{1/2}},$$

and in the prolate case,  $\alpha = R/L < 1$ ,

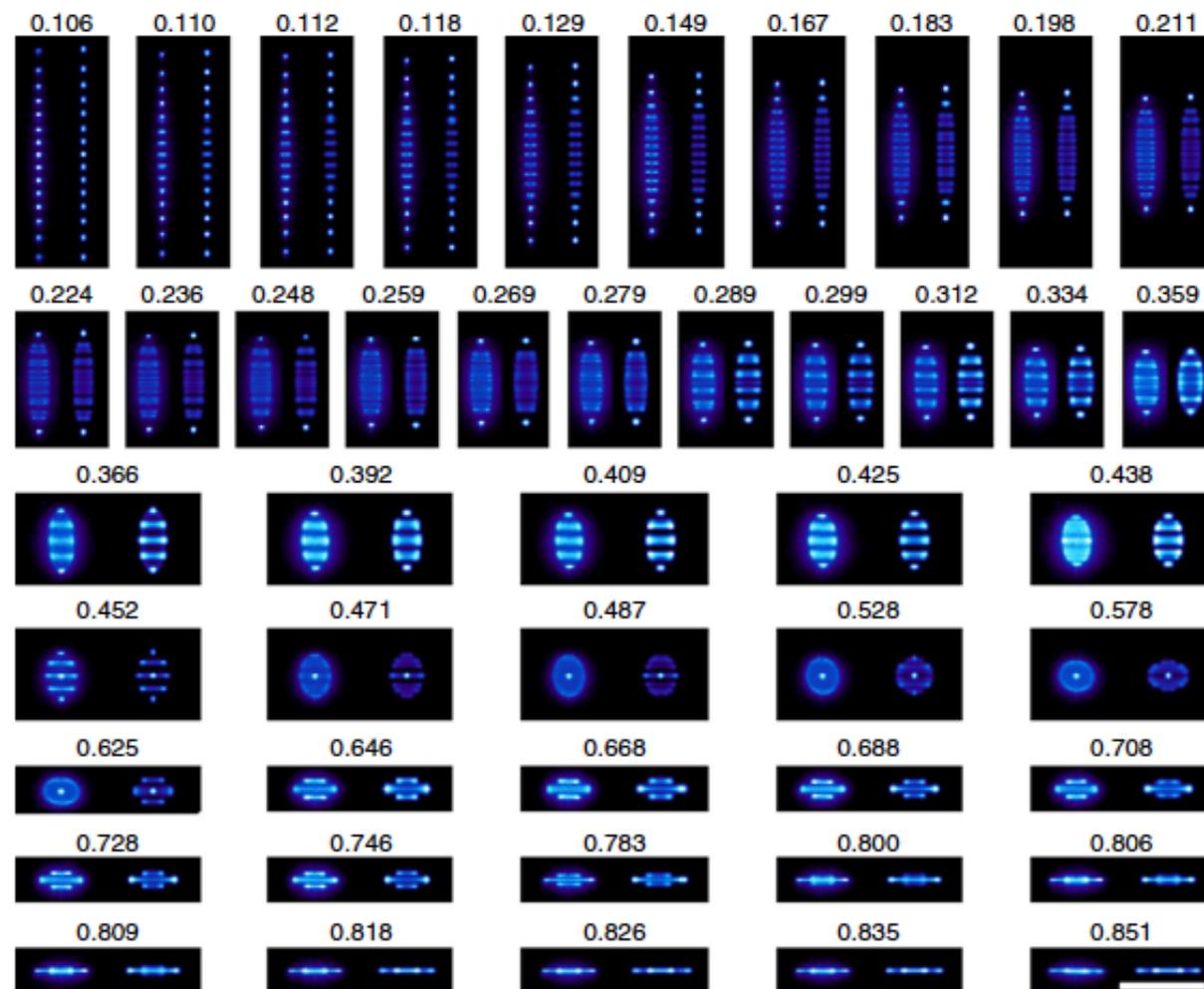
$$\frac{\omega_z^2}{\omega_r^2} = -2 \frac{\sinh^{-1}(\alpha^{-2} - 1)^{1/2} - \alpha(\alpha^{-2} - 1)^{1/2}}{\sinh^{-1}(\alpha^{-2} - 1)^{1/2} - \alpha^{-1}(\alpha^{-2} - 1)^{1/2}}.$$



L. Turner, Phys. Fluids **30**, 3196 (1987)

# Control of the conformations of ion Coulomb crystals in a Penning trap

Sandeep Mavadia<sup>1</sup>, Joseph F. Goodwin<sup>1</sup>, Graham Stutter<sup>1</sup>, Shailen Bharadia<sup>1</sup>, Daniel R. Crick<sup>1</sup>, Daniel M. Segal<sup>1</sup> & Richard C. Thompson<sup>1</sup>



## Structure and Madelung energy of spherical Coulomb crystals

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V. V. Avilov

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(Received 26 March 1991)

$$\mathbf{F}_{\text{conf}}^i = -K\mathbf{r}_i ,$$

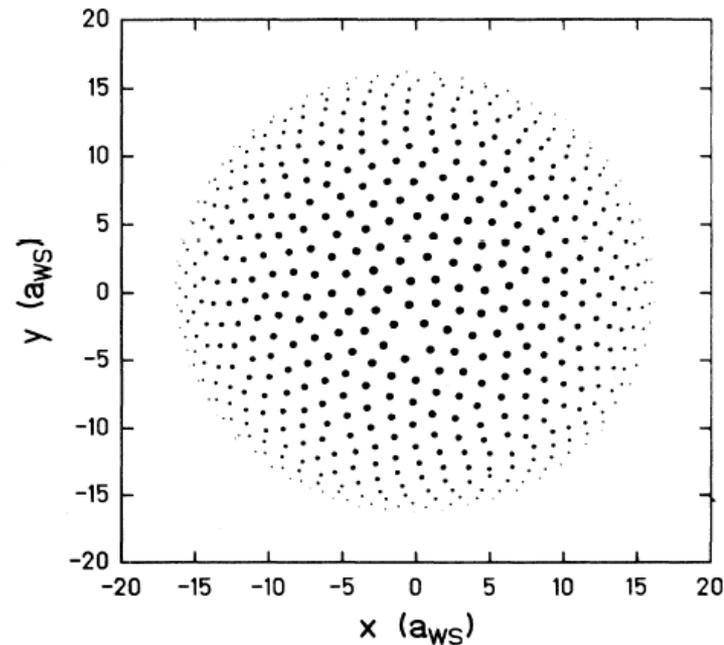


FIG. 1. The front hemisphere of the outer shell of the 5000-ion system. Note the hexagonal structure with approximately equilateral triangles and the point defect below the center.

TABLE I. Structures ( $N_M + N_{M-1} + \dots + N_1$  particles in subshells), rms radii, and energies of  $N$ -particle systems.

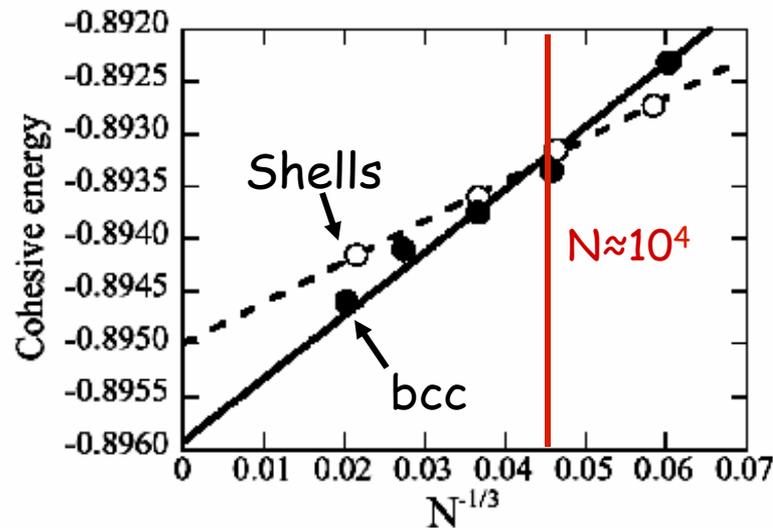
$N$	Structure	$R_{rms}$	$\epsilon$	$\Gamma$
2	2	0.6300	0.5953	$\infty^a$
3	3	0.8327	1.0400	$\infty^a$
4	4	0.9721	1.4174	$\infty^a$
5	5	1.0901	1.7820	$10^5$
6	6	1.1850	2.1065	$10^5$
7	7	1.2736	2.4321	$10^5$
8	8	1.3500	2.7331	$10^5$
9	9	1.4198	3.0238	$10^5$
10	10	1.4846	3.3058	$10^5$
11	11	1.5453	3.5822	$10^5$
12	12	1.6002	3.8407	$10^5$
13	12+1	1.6535	4.1009	$10^5$
26	24+2	2.1633	7.0197	$10^5$
27	24+3	2.1941	7.2206	$10^5$
28	24+4	2.2241	7.4200	$10^5$
29	26+3	2.2532	7.6162	$10^5$
60	48+12	2.9335	12.9085	$10^5$
61	48+12+1	2.9511	13.0611	$10^5$
900	356+247+154+92+38+12+1	7.4389	83.0047	$10^6$
923	361+253+160+93+43+12+1	7.5023	84.4277	$10^7$
1024	390+275+108+111+53+16+1	7.7696	90.5438	$10^6$
1200	448+309+212+132+70+25+4	8.1952	100.7410	$10^6$
1370	482+354+245+155+91+35+8	8.5683	110.1259	$10^6$
1415	491+365+244+163+94+44+13+1	8.6619	112.5436	$10^7$
1600	538+398+285+185+115+58+19+2	9.0270	122.2270	$10^6$
2000	633+470+357+246+163+85+40+6	9.7287	141.1975	$10^6$
2057	641+491+363+257+158+93+44+10	9.8209	144.6764	$10^8$
2837	805+634+480+356+246+173+143 <sup>b</sup>	10.9384	179.4717	$10^7$
2869	808+644+482+361+249+165+160 <sup>b</sup>	10.9796	180.8258	$10^8$
3816	986+793+618+481+938 <sup>b</sup>	12.0798	218.8847	$10^7$
3871	992+801+639+1439 <sup>b</sup>	12.1379	220.9917	$10^8$
5000	1183+980+802+2035 <sup>b</sup>	13.2229	262.2696	$10^9$

"Magic numbers"

# 3D Long-range-order in Coulomb crystals

**Theory:** Wigner, Hansen, Pollock, Dubin, Hasse, Totsuji ..

Recent MD simulations:

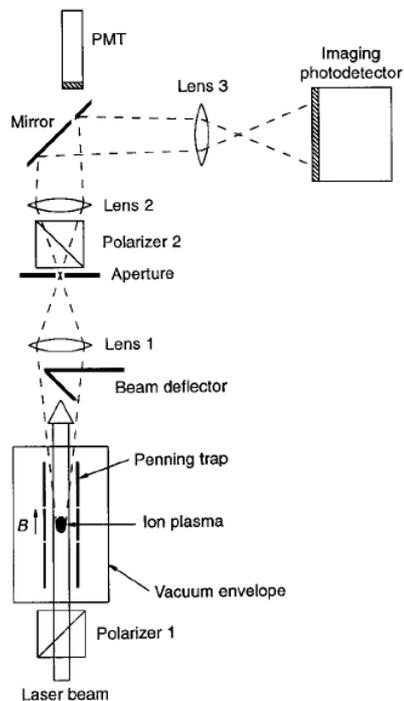


$$U_{\text{coh}} = \Delta U_{\text{sys}} / (Nq^2/a)$$

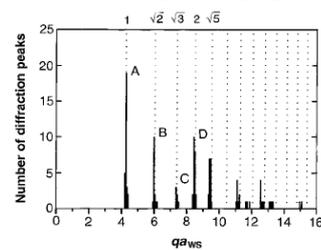
# Bragg Diffraction from Crystallized Ion Plasmas

W. M. Itano,\* J. J. Bollinger, J. N. Tan,† B. Jelenković,‡  
X.-P. Huang, D. J. Wineland

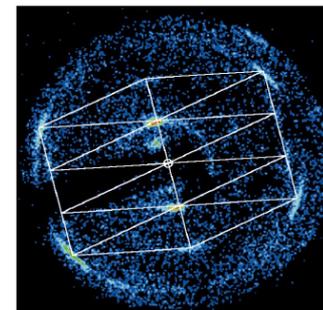
## Penning trap experiments with $\sim 10^5$ ${}^9\text{Be}^+$ ions



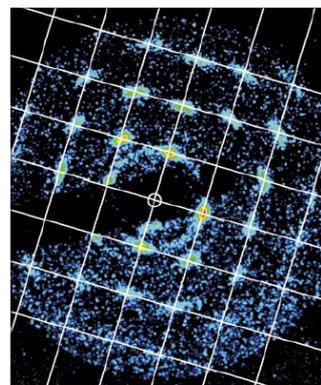
**Fig. 1.** Experimental setup. Laser light is directed through the ion plasma in the Penning trap. A diffraction pattern is created at a plane beyond lens 2, where rays that are parallel leaving the plasma are focused to a point. A mirror, placed near that plane, deflects the light to an imaging photodetector. An aperture placed inside a hole in the mirror allows diffracted light to be detected by a photomultiplier tube (PMT). The aperture is placed off the axis of the optical system, so the PMT generates a timing signal as the diffraction pattern rotates.



**Fig. 3.** Histogram representing the numbers of peaks (not intensities) observed as a function of  $qa_{ws}$ , where  $\mathbf{q} = \mathbf{k}_s - \mathbf{k}_i$  is the difference between the incident ( $\mathbf{k}_i$ ) and scattered ( $\mathbf{k}_s$ ) photon wave vectors. We analyzed 30 Bragg diffraction patterns from two approximately spherical plasmas having 270,000 and 470,000 ions. The dotted lines show the expected peak positions, normalized to the center of gravity of the peak at A ((110) Bragg reflections).



**Fig. 5.** A Bragg diffraction pattern with twofold symmetry. It matches the pattern expected for a bcc lattice oriented along a (115) direction. A diffraction spot is predicted at each intersection of the grid lines. The passive timing method and the MCP-RA detector were used. Here,  $\omega_i = 2\pi \times 149$  kHz,  $n_0 = 4.53 \times 10^9$  cm $^{-3}$ ,  $N = 4 \times 10^5$ ,  $\alpha = 1.20$ , and  $2r_0 = 1.12$  mm.



**Fig. 4.** Time-resolved Bragg diffraction pattern of the same plasma as in Fig. 2. Here and in Figs. 5 and 6 the small open circle marks the position of the undeflected laser beam. A bcc lattice, aligned along a (100) axis, would generate a spot at each intersection of the grid lines overlaid on the image. The grid spacing corresponds to an angular deviation of  $2.54 \times 10^{-2}$  rad. Here,  $\omega_i = 2\pi \times 125.6$  kHz,  $n_0 = 3.83 \times 10^9$  cm $^{-3}$ ,  $N = 5 \times 10^5$ ,  $\alpha = 0.98$ , and  $2r_0 = 1.36$  mm.

# Direct Observations of Structural Phase Transitions in Planar Crystallized Ion Plasmas

T. B. Mitchell,\* J. J. Bollinger, D. H. E. Dubin, X.-P. Huang, W. M. Itano, R. H. Baughman

Fig. 1. Schematic side view of the cylindrical Penning trap with its side- and top-view imaging optics. The insets show the variables used to characterize the intra- and inter-layer structure. The side-view inset also shows the central region of a lenticular ion plasma with three axial lattice planes. The rotation frequency of the ion plasma was controlled by applying properly phased sinusoidally time-varying electric potentials to the sixfold azimuthally segmented compensation electrodes.

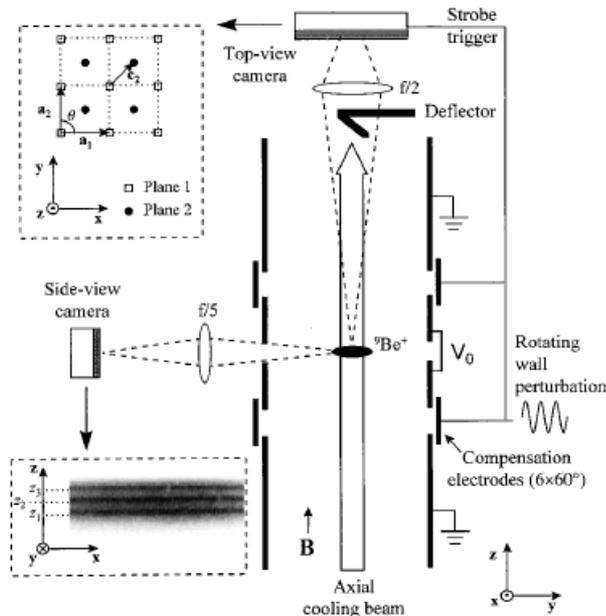


Table 1. Primitive and interlayer displacement vectors in the  $(x, y)$  plane for the observed phases. The primitive vector  $a_1$  defines the  $x$  direction, and  $|a_1| = |a_2| = a$ . Dashes, not applicable.

Phase	Symmetry	Stacking	Vectors			
			$a_1$	$a_2$	$c_2$	$c_3$
I	Hexagonal	Single plane	$(a, 0)$	$(a \cos 60^\circ, a \sin 60^\circ)$	—	—
III	Square	Staggered	$(a, 0)$	$(0, a)$	$(a_1 + a_2)/2$	$(0, 0)$
IV	Rhombic	Staggered	$(a, 0)$	$(a \cos \theta, a \sin \theta)$	$(a_1 + a_2)/2$	$(0, 0)$
V	Hexagonal	hcp-like	$(a, 0)$	$(a \cos 60^\circ, a \sin 60^\circ)$	$(a_1 + a_2)/3$	$(0, 0)$
$V_{fcc}$	Hexagonal	fcc-like	$(a, 0)$	$(a \cos 60^\circ, a \sin 60^\circ)$	$(a_1 + a_2)/3$	$2(a_1 + a_2)/3$

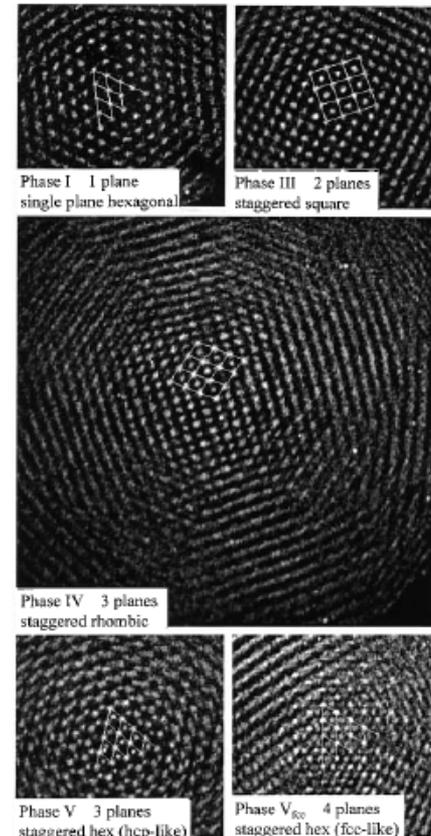


Fig. 2. Top-view  $(x, y)$  images of the five structural phases observed in the experiment, with lines showing a fit of the central ions to the indicated structure.

## The Structure of the Cylindrically Confined Coulomb Lattice

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Received February 27, 1990

For an infinitely long cylinder of radius  $R_0$  with a charge density  $\sigma$  per unit length the repulsive Coulomb force for particles with charge  $q$  is

$$F_{\text{int}} = 2q\sigma \begin{cases} r/R_0^2, & r < R_0 \\ 1/r, & r > R_0 \end{cases} \quad (1.1)$$

and with the external confining (focusing) force

$$\mathbf{F}_{\text{conf}} = -qK\mathbf{r} \quad (1.2)$$

the equilibrium condition is that

$$F_{\text{conf}} + F_{\text{int}} = 2\sigma q/R_0^2 - qK \equiv 0 \quad (1.3)$$

yielding

$$R_0 = \sqrt{2\sigma/K}. \quad (1.4)$$

tions of infinite one-component plasmas [7, 8]. Outside the cylinder, for  $r > R_0$  there is, of course, an increasing radial restoring force

$$F_{\text{tot}} = -qK(r - R_0^2/r). \quad (1.5)$$

The average density on the inside of the cylinder being uniform, the average volume per particle is  $\pi R_0^2 q/\sigma = 2\pi q/K$ , and setting this equal to a volume of  $4\pi a^3/3$  yields a Wigner-Seitz radius,

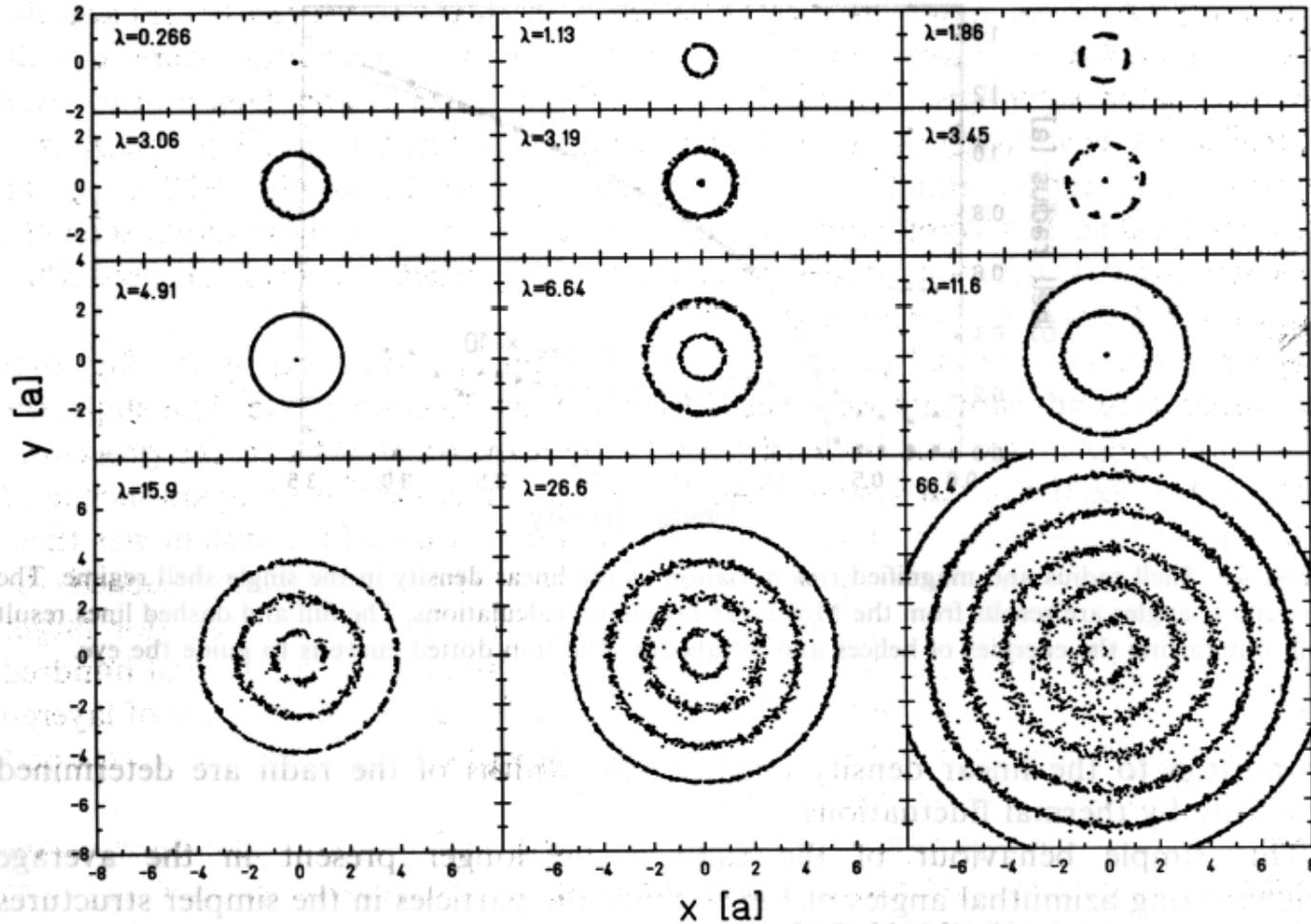
$$a = \left(\frac{3q}{2K}\right)^{1/3}. \quad (1.6)$$

Note that the average distance between neighbouring ions in large systems is almost twice this value. In units of  $a$  we define the dimensionless linear particle density

$$\lambda \equiv \sigma a/q \quad (1.7)$$

# The structure of cylindrically confined infinite crystals

## Projections to the radial plane



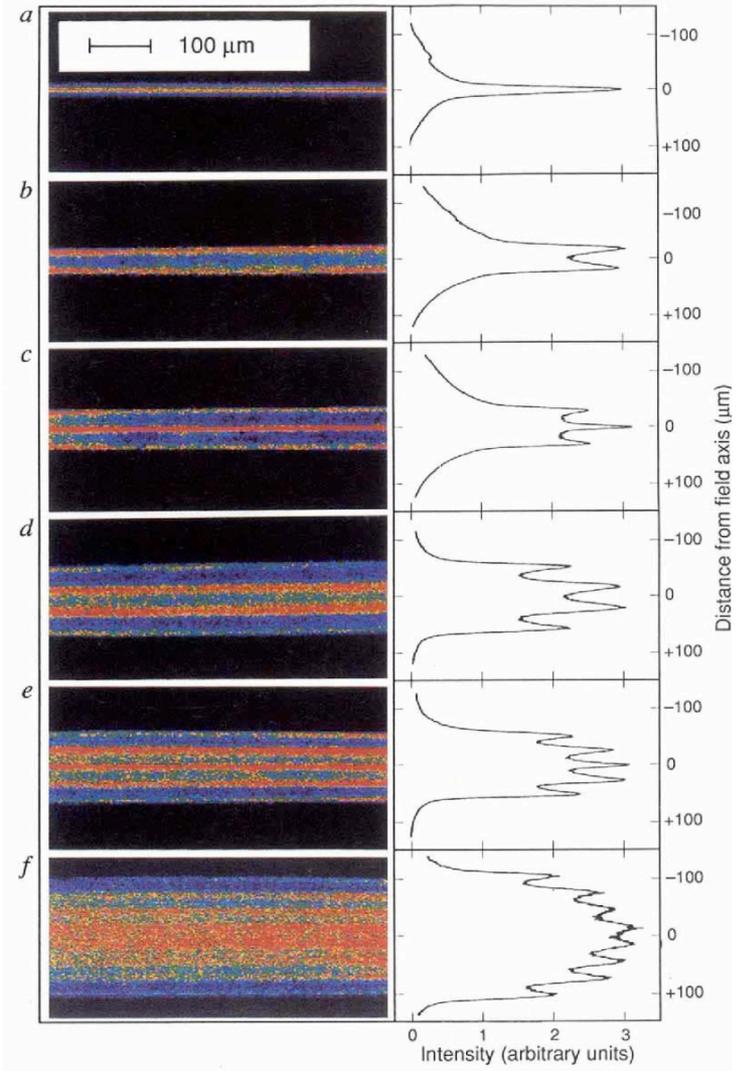
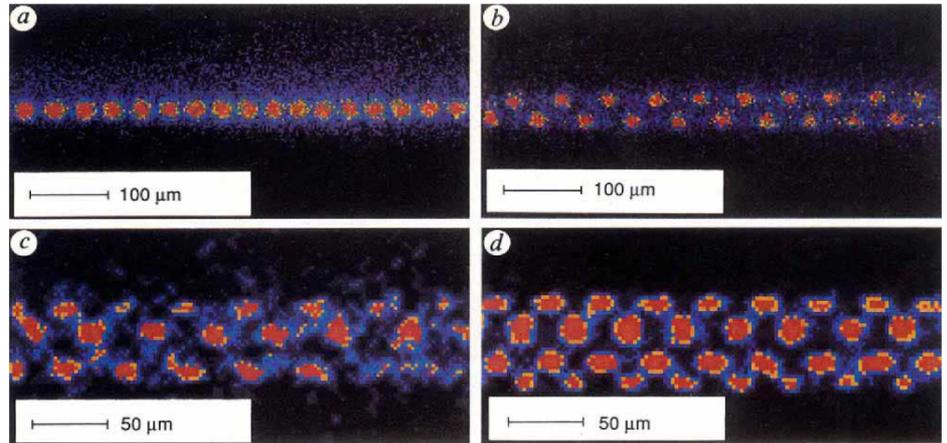
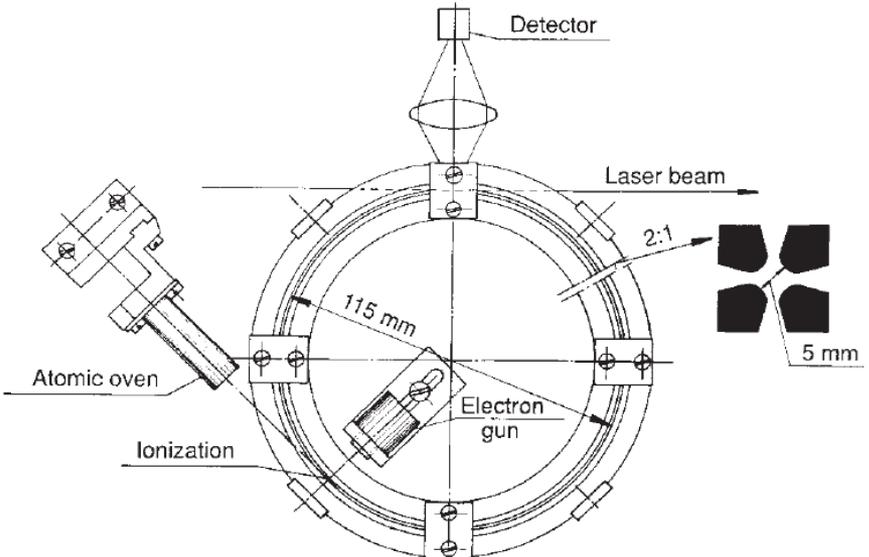
Dimensionless linear charge density:  $\lambda \equiv \sigma a / q$

# The structure of cylindrically confined infinite crystals

## Structures for increasing linear densities

Total linear density	Structure
$0 < \lambda < 0.709$	String
$0.709 < \lambda < 0.964$	Zigzag
$0.964 < \lambda < 1.25$	1st Helix solution
$1.25 < \lambda < 1.39$	Tetrahedron
$1.39 < \lambda < 1.70$	1st Helix solution
$1.70 < \lambda < 2.19$	Tetrahedron
$2.19 < \lambda < 2.52$	2nd Helix solution
$2.52 < \lambda < 2.96$	Helix
$2.96 < \lambda < 3.10$	3rd Helix solution
$3.10 < \lambda < 5.7$	Shell + string
$5.7 < \lambda < 9.5$	2 Shells
$9.5 < \lambda < 13$	2 Shells + string
$\lambda = 19.9$	3 Shells + string
$\lambda = 26.6$	4 Shells
$\lambda = 66.4$	6 Shells

# Observation of multi-shell structures in a quadrupole storage ring



**G. Birkel, S. Kassner & H. Walther**

NATURE · VOL 357 · 28 MAY 1992



FIG. 1(color). (a) CCD picture integrated over 2 sec of a very prolate crystal of about 3500 ions. (b) Molecular dynamic simulation with 3500 ions in a crystal with roughly the same aspect ratio as the one in (a).

## Large Ion Crystals in a Linear Paul Trap

M. Drewsen, C. Brodersen, L. Hornekær, and J. S. Hangst

*Institute of Physics and Astronomy, University of Aarhus, DK-8000 Aarhus, Denmark*

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*and University of Chicago, Chicago, Illinois 60637*

(Received 5 December 1997)

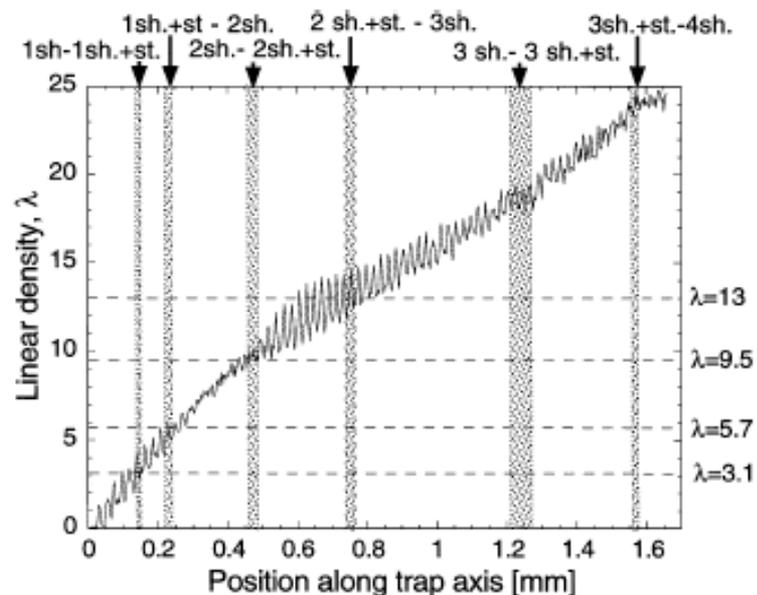


FIG. 2. The dimensionless linear density  $\lambda$  as a function of position along the crystal is shown in Fig. 1a (solid line). The shaded vertical areas indicate the uncertainties in the measured position of structural transitions, while the dashed horizontal lines correspond to linear densities where these transitions should happen according to MD simulation for infinite long crystals.

# Two-species ion Coulomb crystals

## Effective oscillation freq.'s:

$$\omega_r = 1/2 \beta \Omega, \quad \beta = (1/2 q^2 + a)^{1/2}$$

$$\omega_z = (-1/2 a)^{1/2} \Omega$$

$$q = \frac{4QU_{RF}}{m\Omega^2 r_0^2} \quad a = -\frac{\alpha QU_{end}}{m\Omega^2 r_0^2}$$

## Atomic density:

$$n_i = (\epsilon_0 U_{rf}^2) / (M_i r_0^4 \Omega^2)$$

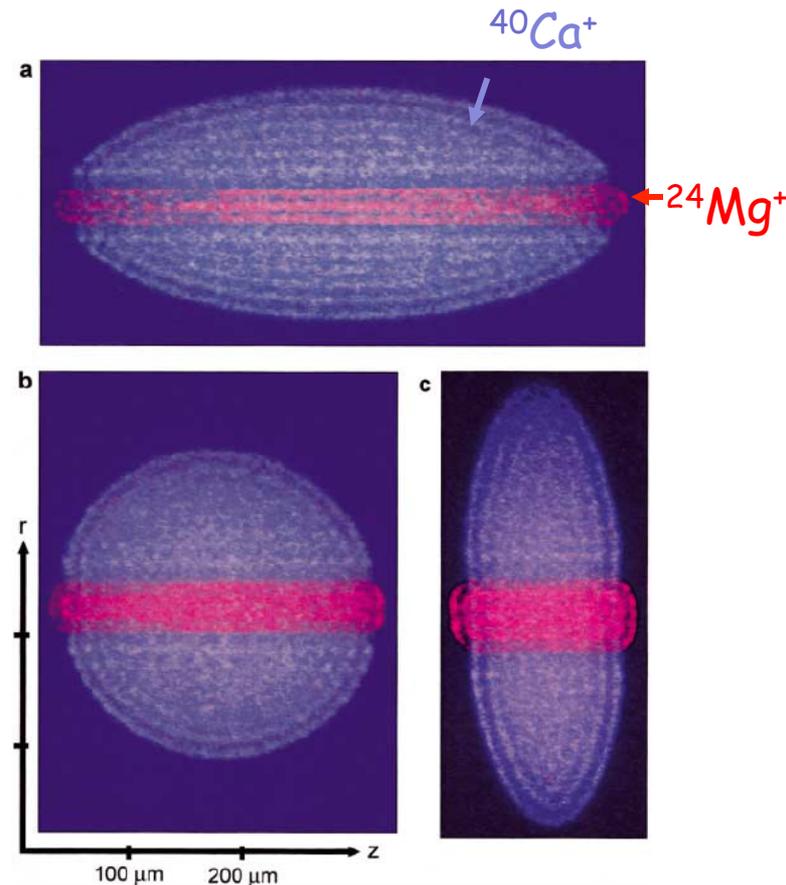
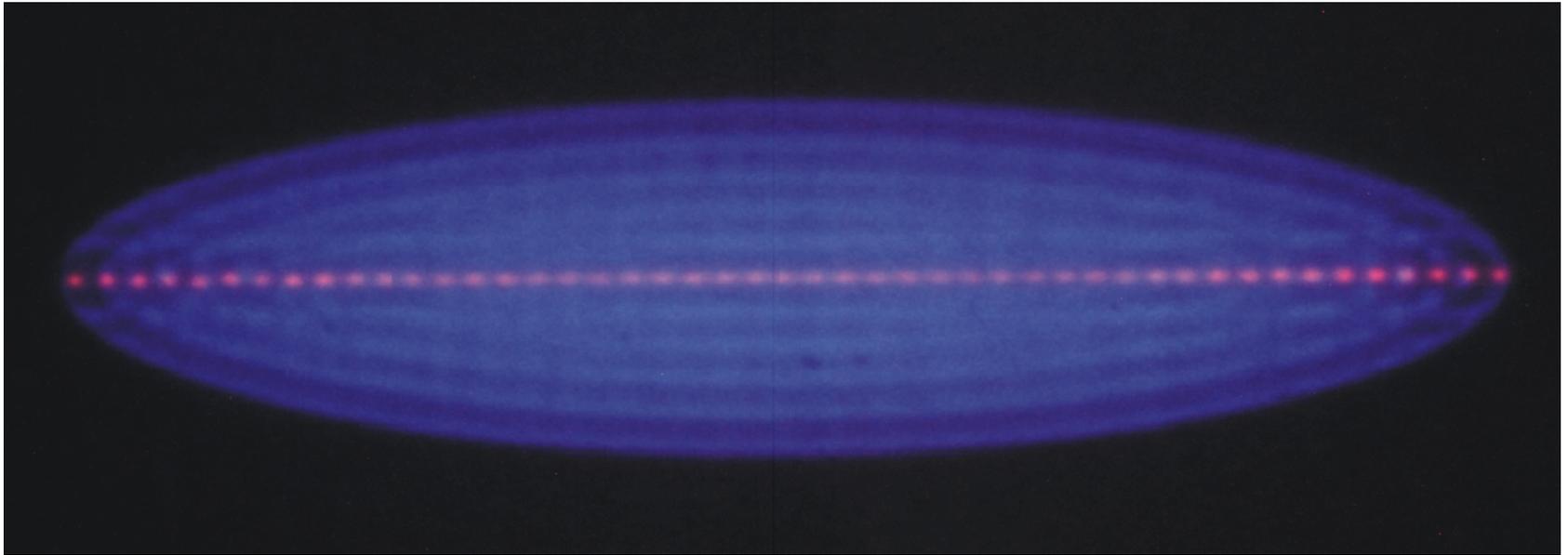


FIG. 2 (color). A  $^{40}\text{Ca}^+ / ^{24}\text{Mg}^+$  bicrystal at three different end cap voltages. The crystal is symmetric under rotations around the trap axis,  $z$ , and contains approximately 300  $^{24}\text{Mg}^+$  ions (red) and 3000  $^{40}\text{Ca}^+$  ions (blue). The ratios of the axial and effective radial trapping frequencies for  $^{40}\text{Ca}^+$  and  $^{24}\text{Mg}^+$  ions in the three cases shown are (a)  $\omega_{z,\text{Mg}^+} / \omega_{r,\text{Mg}^+} = 0.4$  and  $\omega_{z,\text{Ca}^+} / \omega_{r,\text{Ca}^+} = 0.6$ ; (b)  $\omega_{z,\text{Mg}^+} / \omega_{r,\text{Mg}^+} = 0.7$  and  $\omega_{z,\text{Ca}^+} / \omega_{r,\text{Ca}^+} = 1.0$ ; (c)  $\omega_{z,\text{Mg}^+} / \omega_{r,\text{Mg}^+} = 1.1$  and  $\omega_{z,\text{Ca}^+} / \omega_{r,\text{Ca}^+} = 1.8$ .

$\sim 3000 \text{ } ^{40}\text{Ca}^+ + 47 \text{ } ^{24}\text{Mg}^+$



Equidistantly spaced ions in the string!

Similar to an infinite long single component crystal!

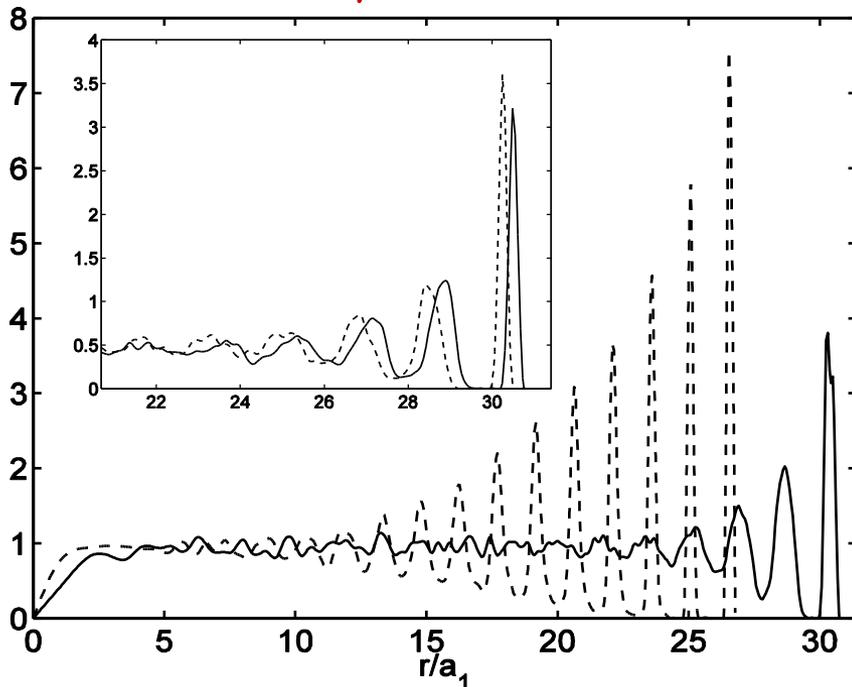
# Coulomb crystals of two species with the same charge-to-mass ratio

Simulations: 50:50 systems with  $q_2=2q_1$ ,  $m_2=2m_1$

Full mixing and double shells

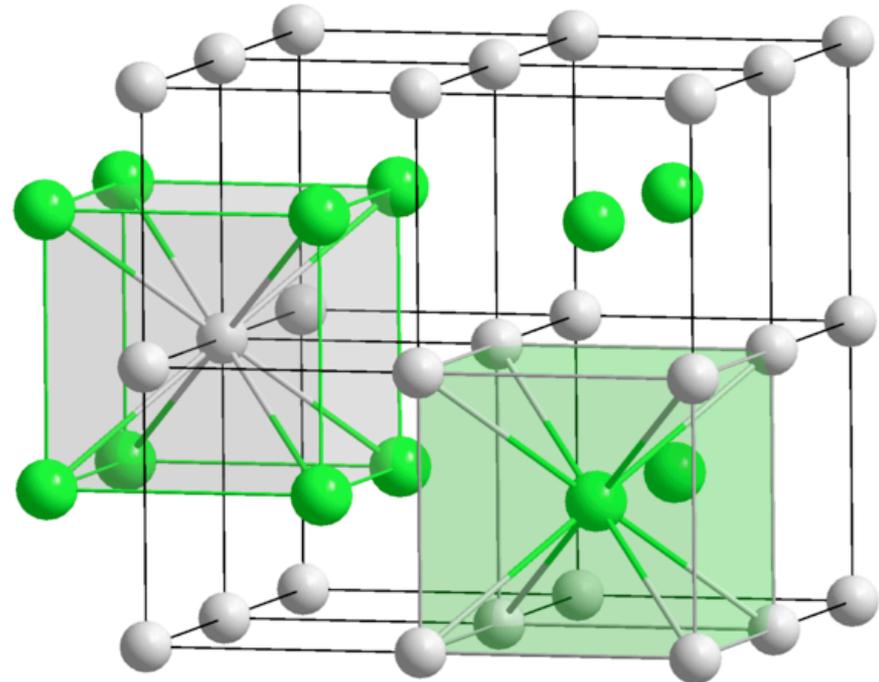
Radial density distributions

20,288 ions



$$n_i = (\epsilon_0 U_{\text{rf}}^2) / (M_i r_0^4 \Omega^2)$$

Ground state vs.  $N^{-1/3}$

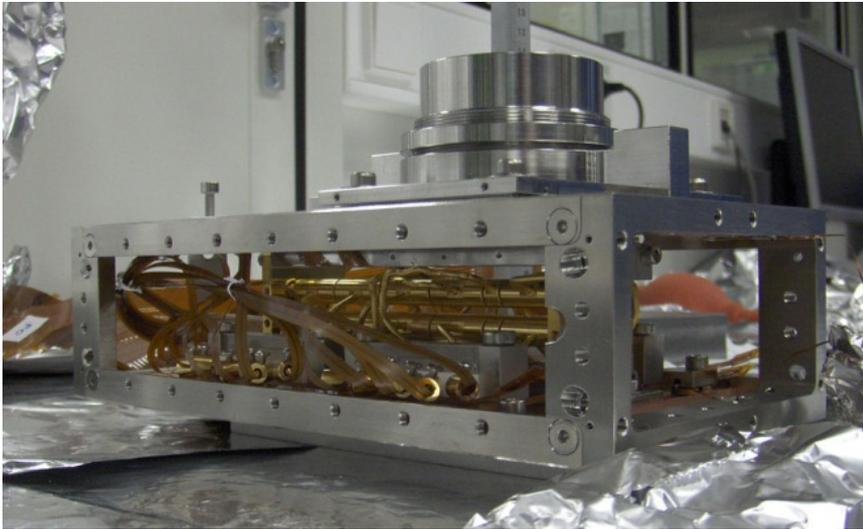


# Coulomb crystals of two species with the same charge-to-mass ratio

Potential Exp.: 50:50 systems of  $^{24}\text{Mg}^+$  and  $^{48}\text{Ca}^{2+}$

More extreme cases, e.g.:

$^9\text{Be}^+$  -  $^{27}\text{Al}^{3+}$ ,  $^{36}\text{Ar}^{4+}$ ,  $^{45}\text{Sc}^{5+}$ ,  $^{54}\text{Fe}^{6+}$ , ...,  $^{180}\text{Hf}^{20+}$ ,  $^{189}\text{Os}^{21+}$ , ...



... the sensitivity of the mixing on the charge-to-mass ratio...

...one per thousand is enough to observe at least partial segregation of the two species.

Ref: PRL 91 165001 (2003)

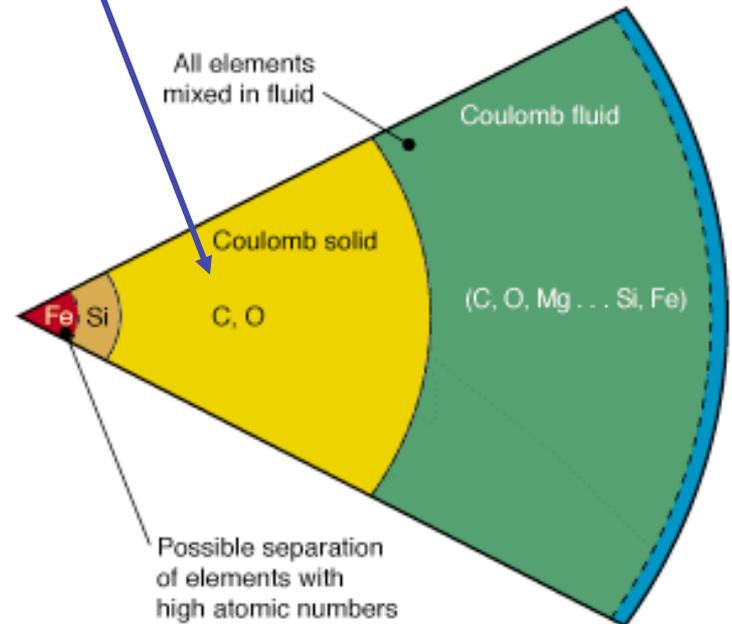
**This can be tested experimentally!**

# White Dwarfs (revisited)

The “Sun” 5 Billion years from now!

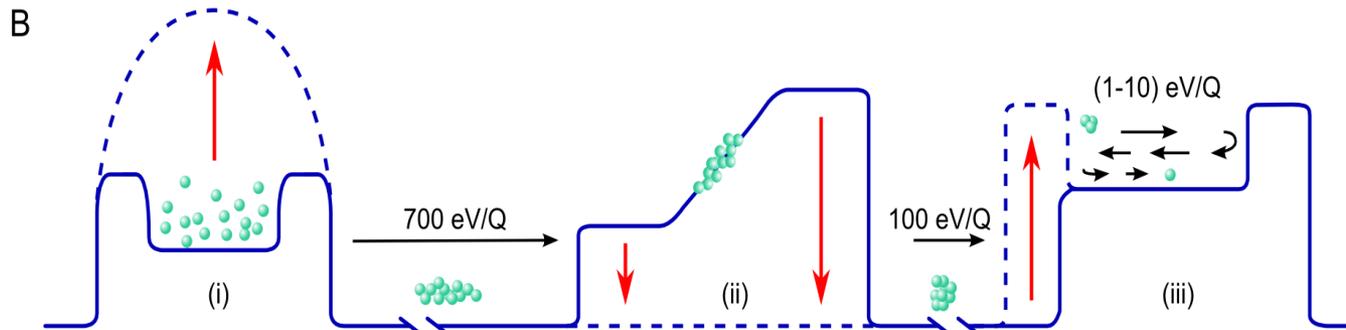
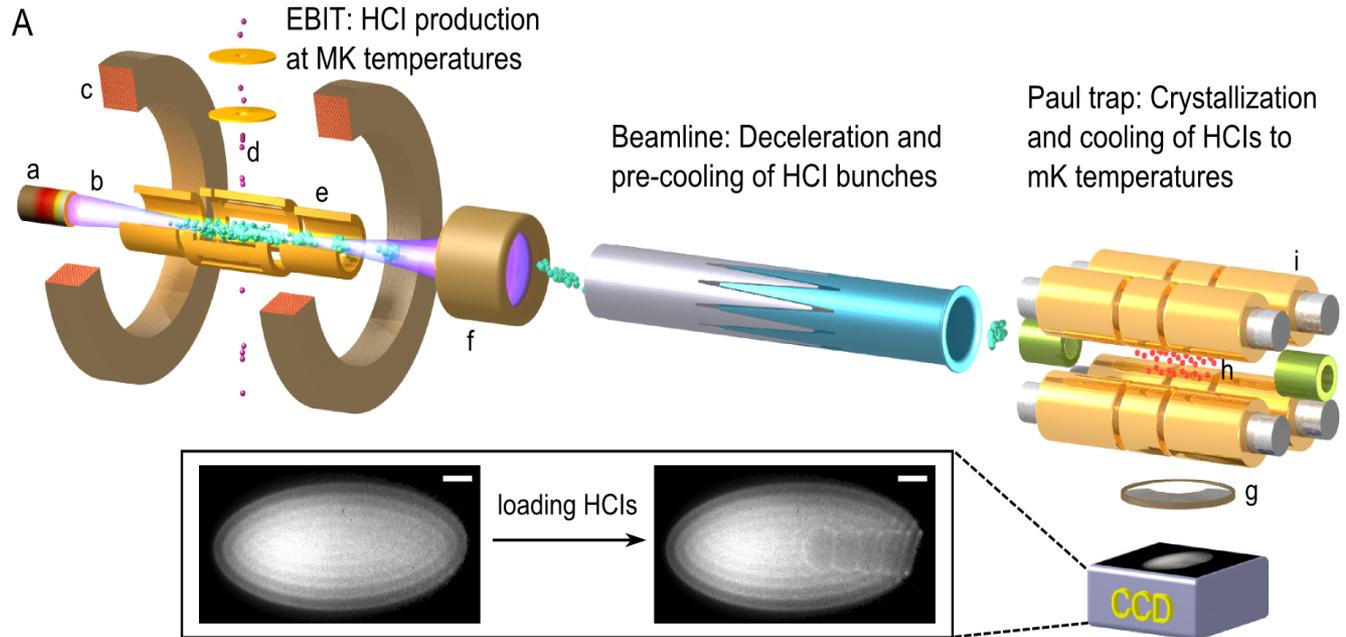


$^{12}\text{C}^{6+}$  and  $^{16}\text{O}^{8+}$

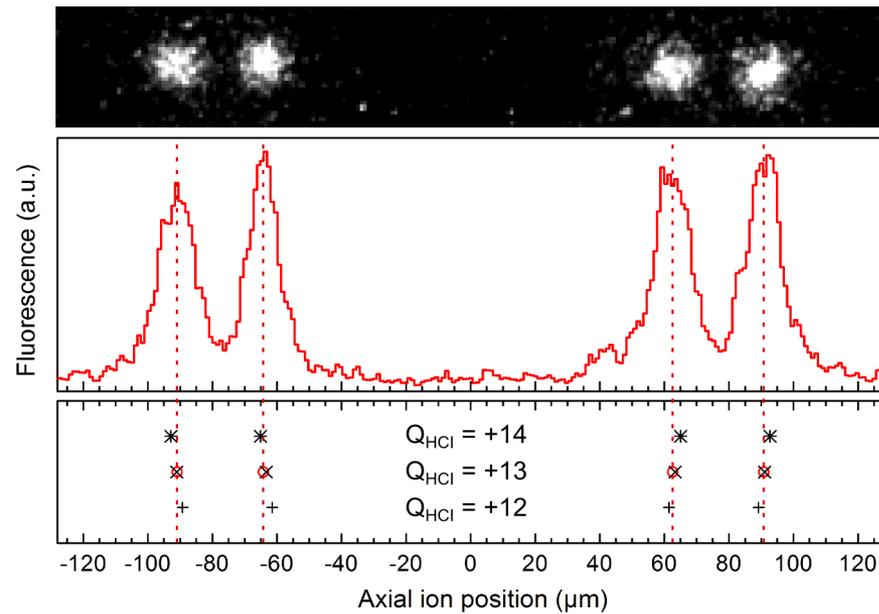


In collaboration with Piet Schmidt, PTB,  
and J. José R. Crespo López Urrutia, MPIK (Project leader)

$^9\text{Be}^+$  and  $^{40}\text{Ar}^{13+}$



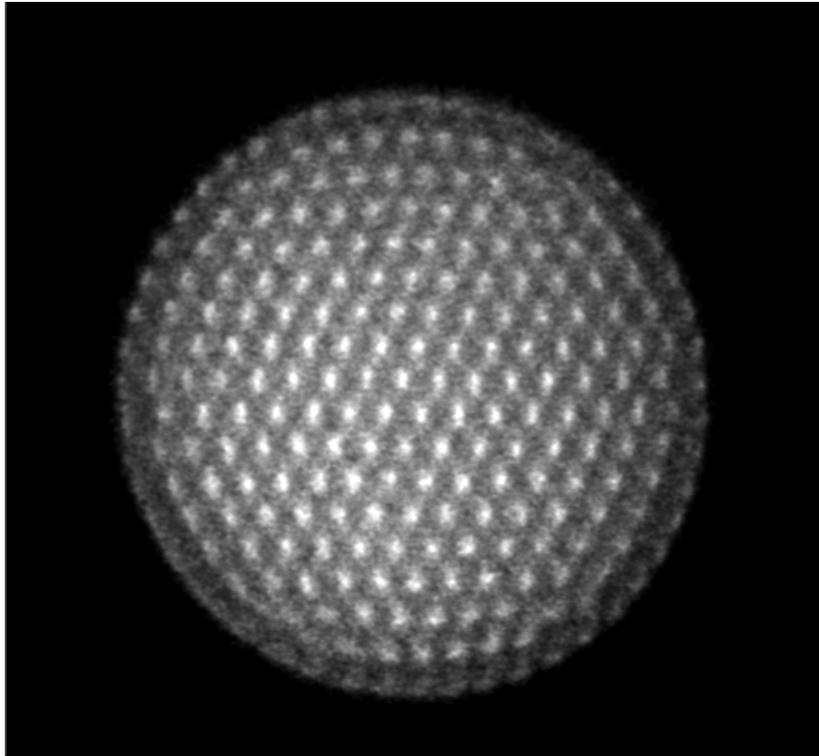
# Four ${}^9\text{Be}^+$ and one ${}^{40}\text{Ar}^{13+}$ ions



### **III. Metastable long-range-ordered Coulomb clusters**

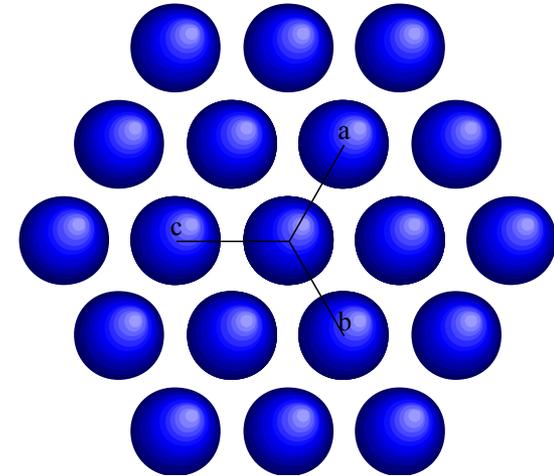
# Observed structures in a linear Paul trap

Number of  $^{40}\text{Ca}^+$  ions  $\sim 2700$



$\sim 300 \mu\text{m}$

Cubic structures viewed along one of the  $\langle 111 \rangle$  directions:



The structures are the same, BUT densities different!

$$(n_{\text{bcc}} = 2n_{\text{sc}} = 4n_{\text{fcc}})$$

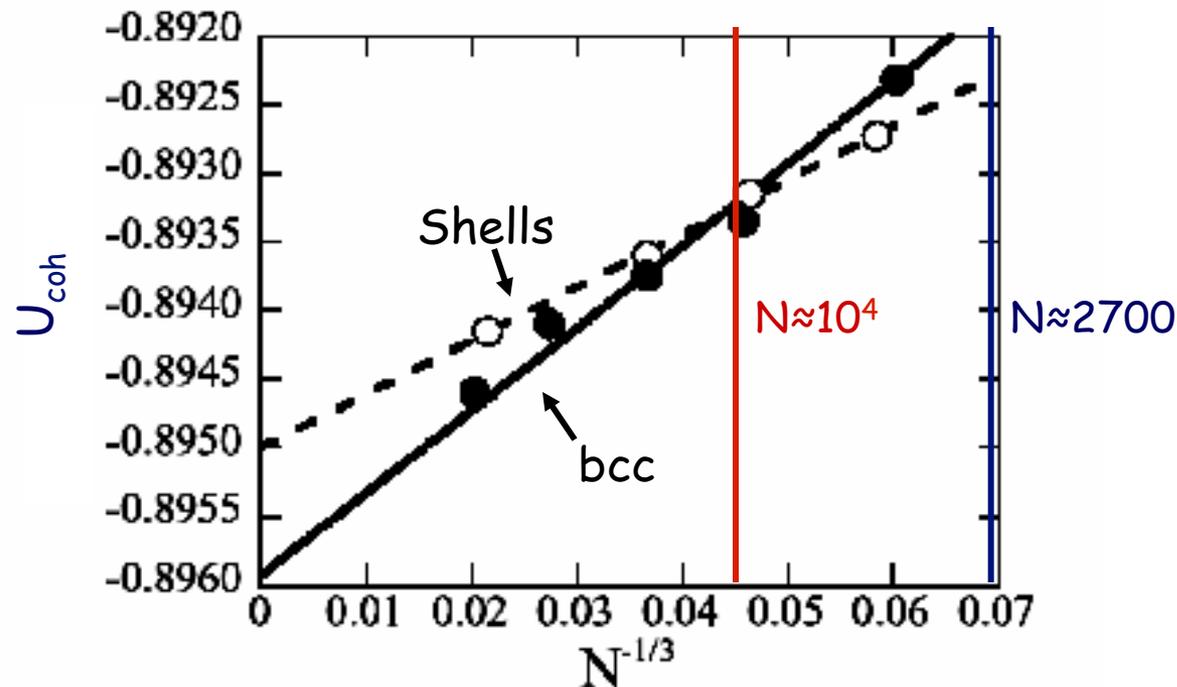
Expected ion density:  $n_{\text{ions}} = 2.4 \pm 2 \times 10^8 \text{ cm}^{-3}$

Ion density assuming bcc structure:  $n_{\text{bcc}} = 2.3 \pm 2 \times 10^8 \text{ cm}^{-3}$

# 3D Long-range-order (theory)

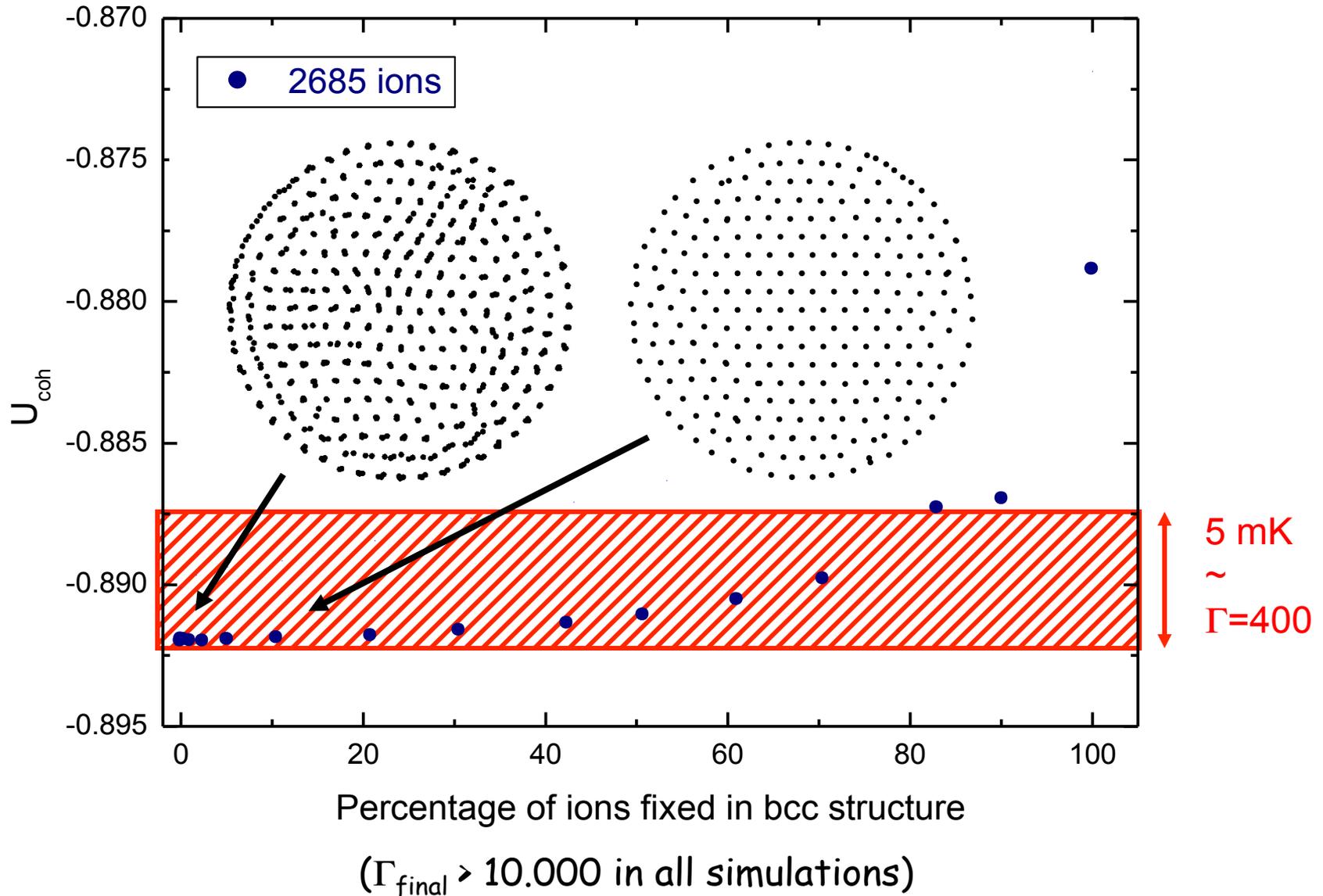
Theory: Wigner, Hansen, Pollock, Dubin, Hasse, Totsuji, ...

Recent MD simulations:



Cohesive energy:  $U_{\text{coh}} = \Delta U_{\text{sys}} / (Nq^2/a)$

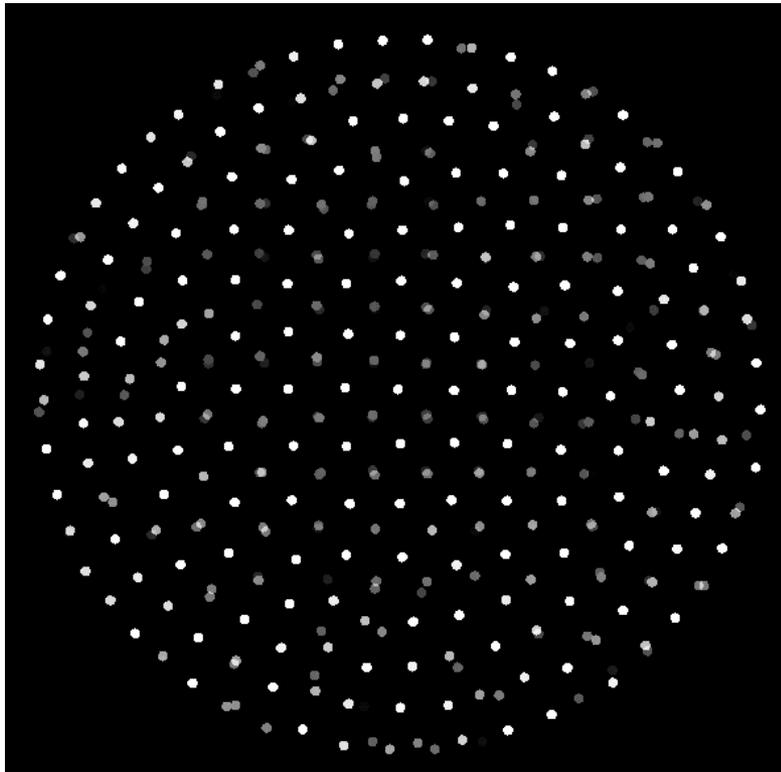
# MD simulations of spherical clusters with fixed central bcc-structures



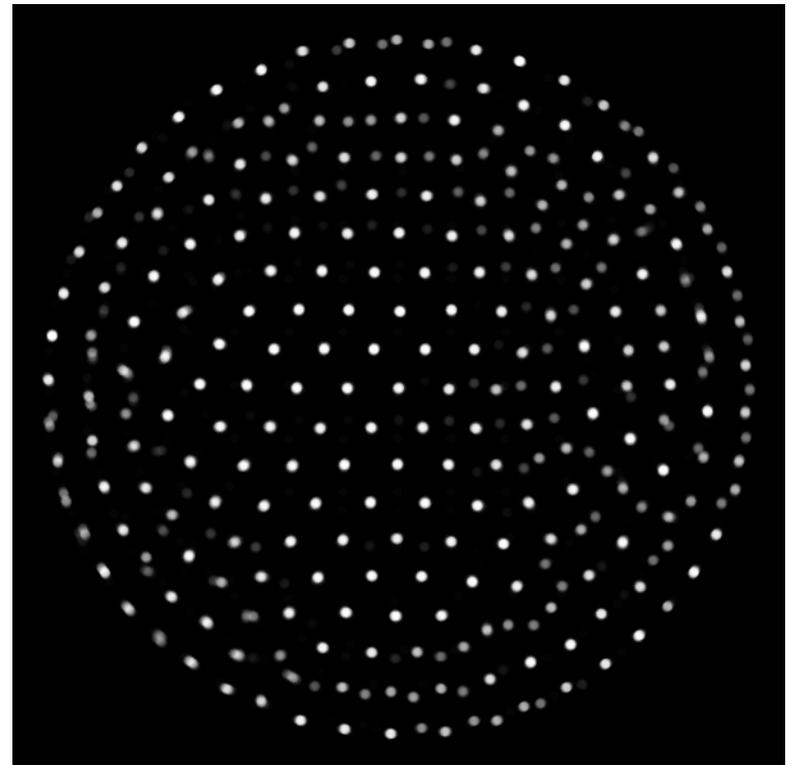
# MD simulation

2685 ions,  $T=1$  mK ( $\Gamma \sim 2000$ )

Video: 10 ms real time



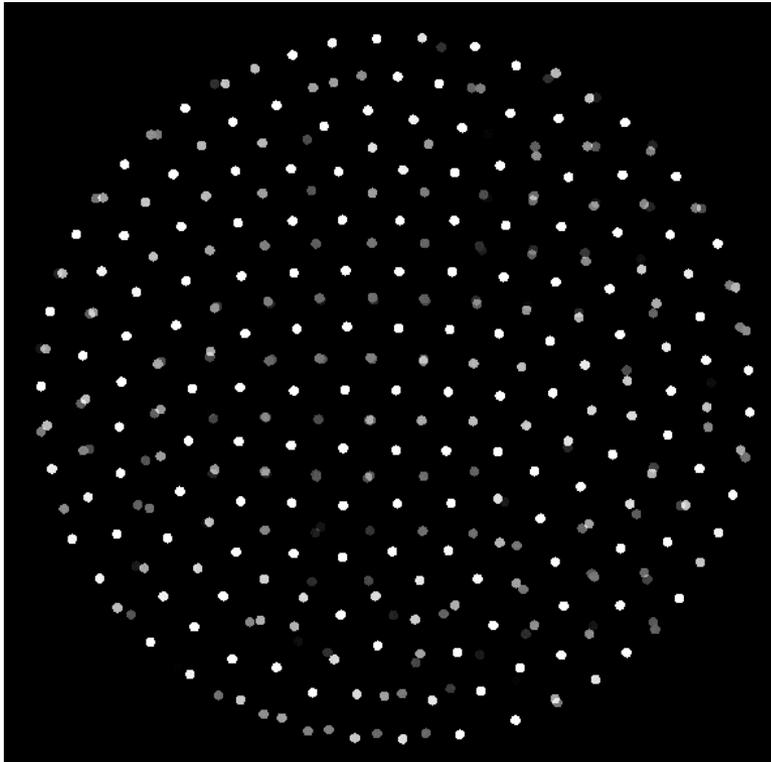
Averaged ion distribution (10 ms)



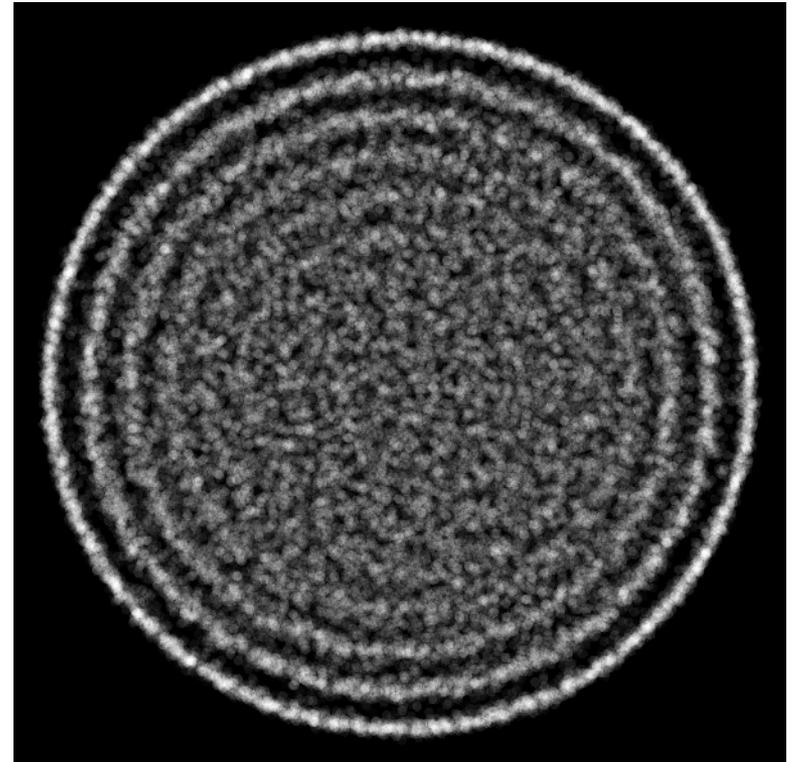
# MD simulation

2685 ions,  $T=20$  mK ( $\Gamma\sim 100$ )

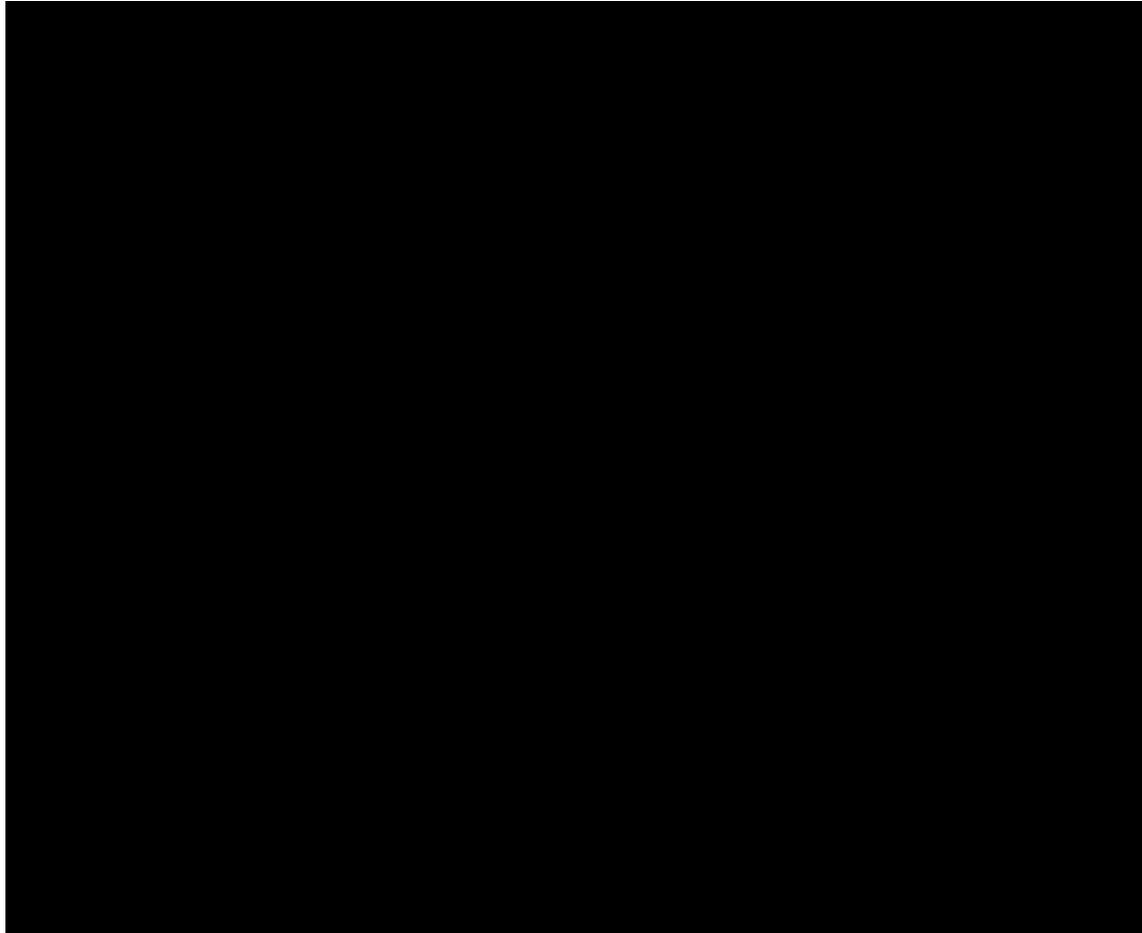
Video: 10 ms real time



Averaged ion distribution (10 ms)

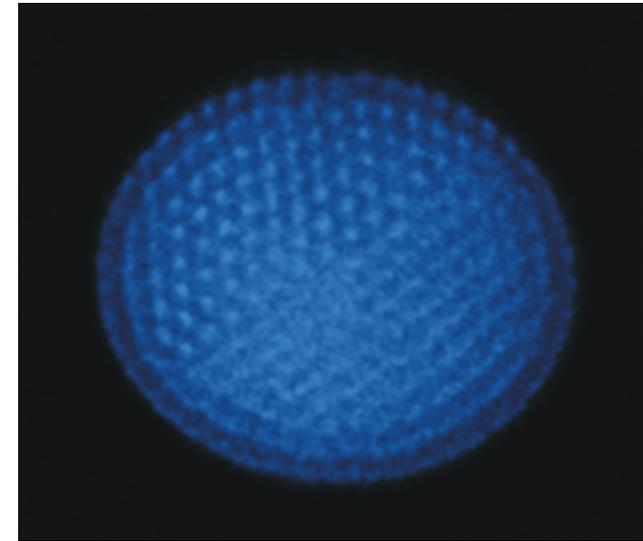
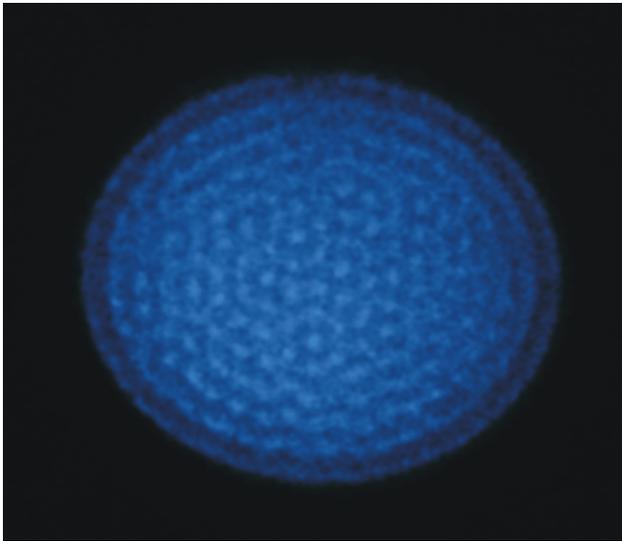
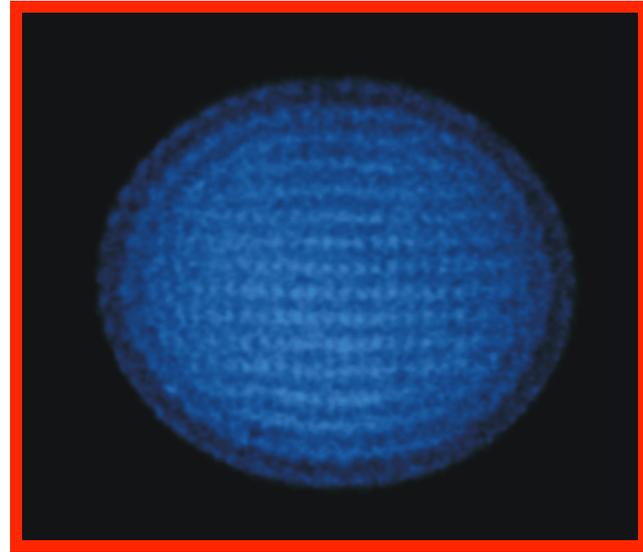
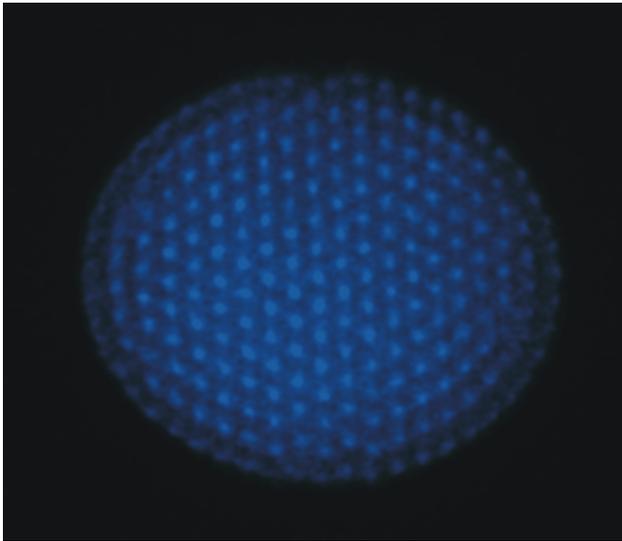


# Real-time video



Number of ions: ~2500

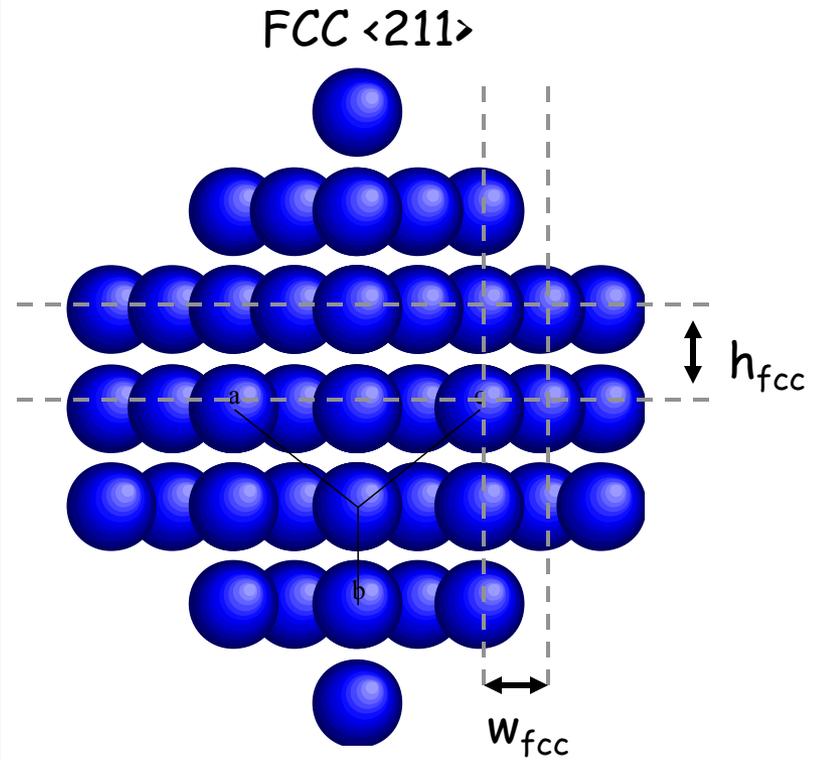
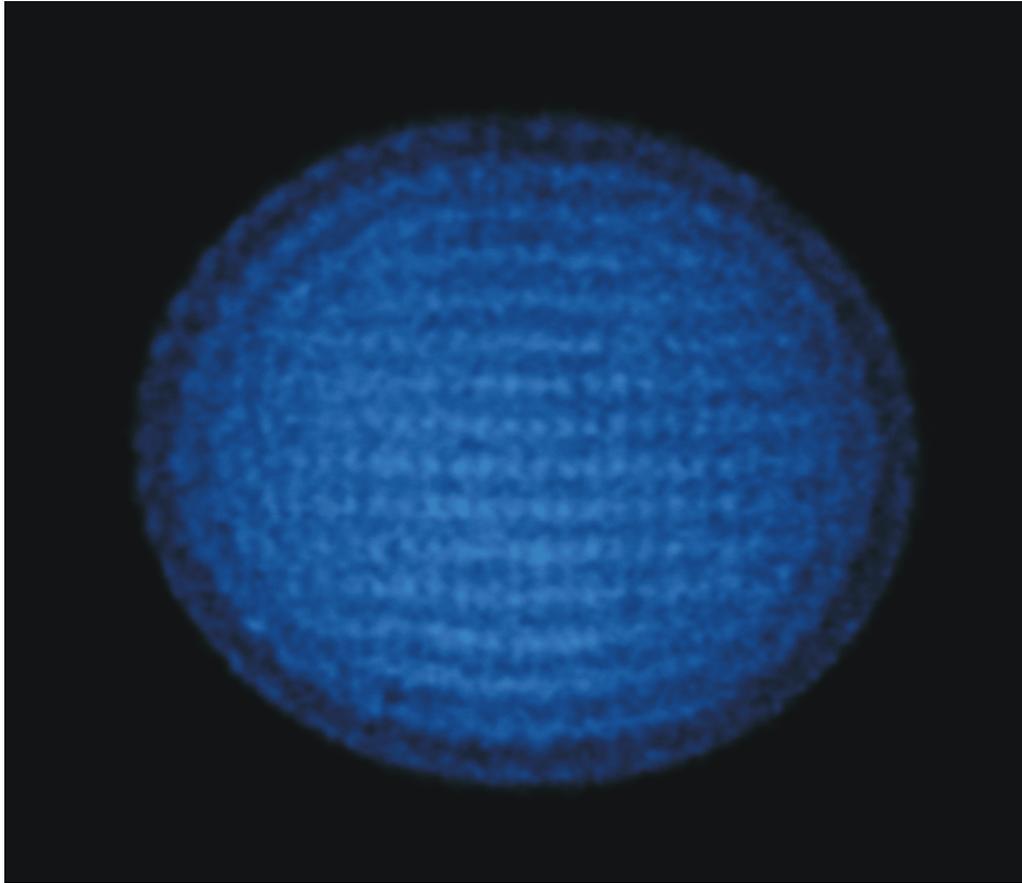
# "Snap-shots"



1 frame = 100 ms  $\approx 10^4 \tau_{osc}$

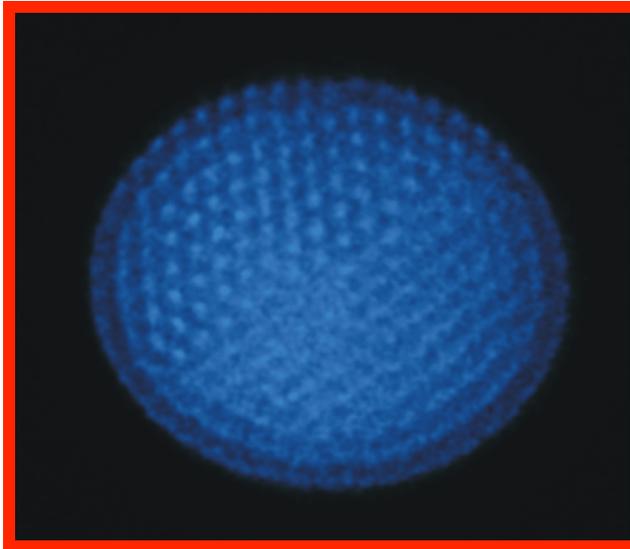
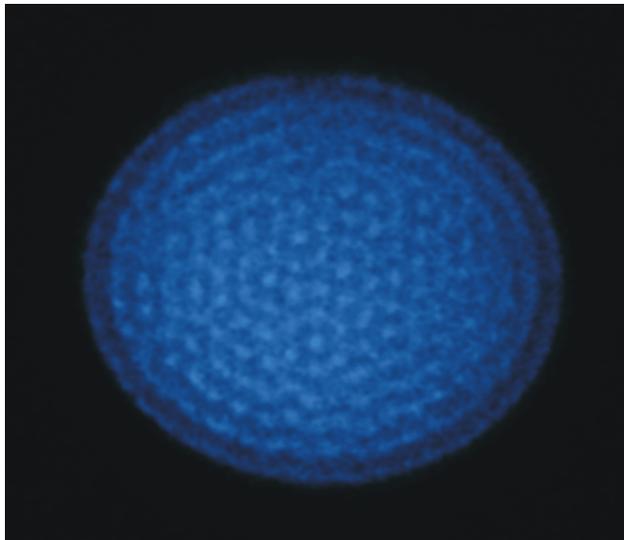
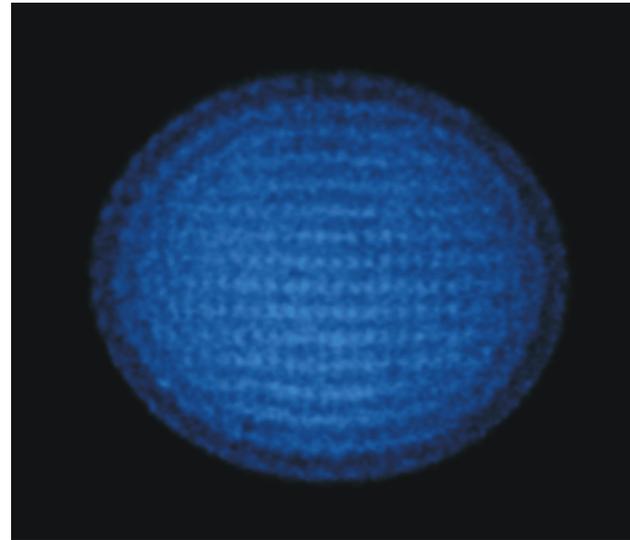
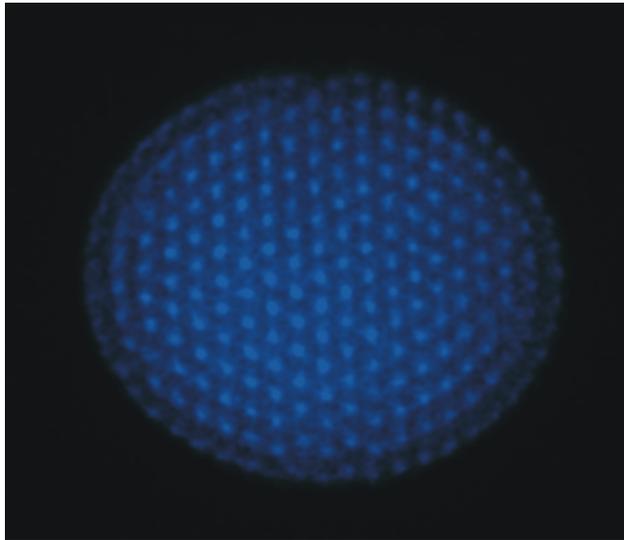
arXiv:1202.2544 (2012)

# FCC structure



$$h_{fcc}/w_{fcc} = 2^{3/2}/3^{1/2} = 1.63$$

# "Snap-shots"

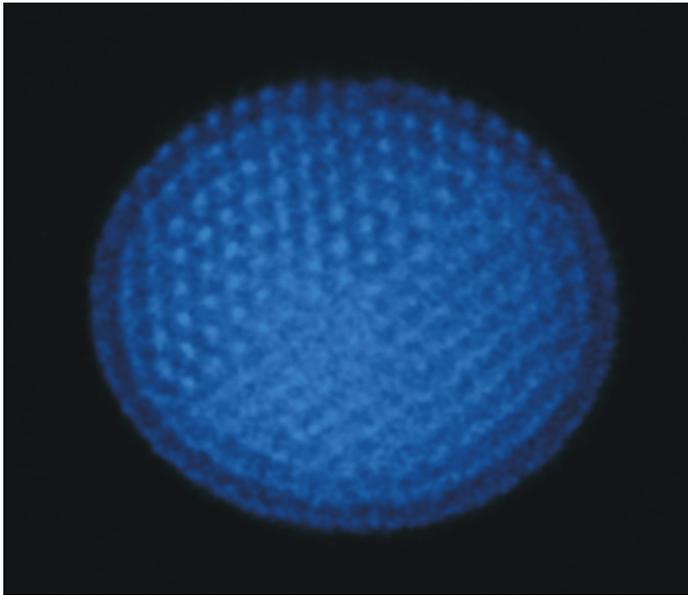


1 frame = 100 ms  $\approx 10^4 \tau_{osc}$

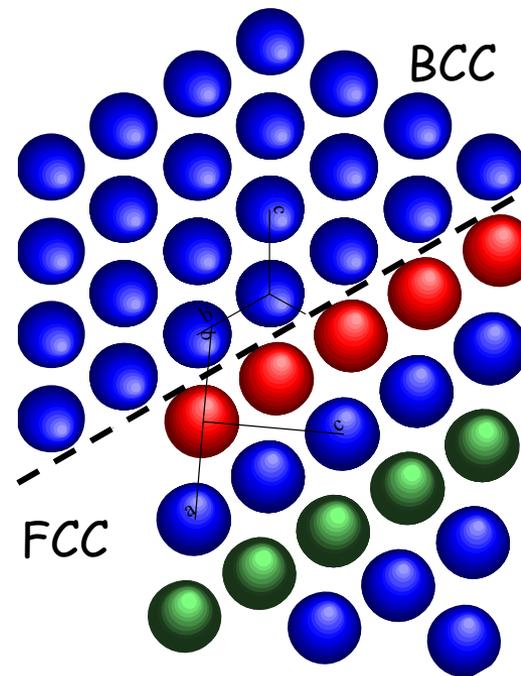
arXiv:1202.2544 (2012)

# BCC $\langle 110 \rangle$ - FCC $\langle 111 \rangle$ interface

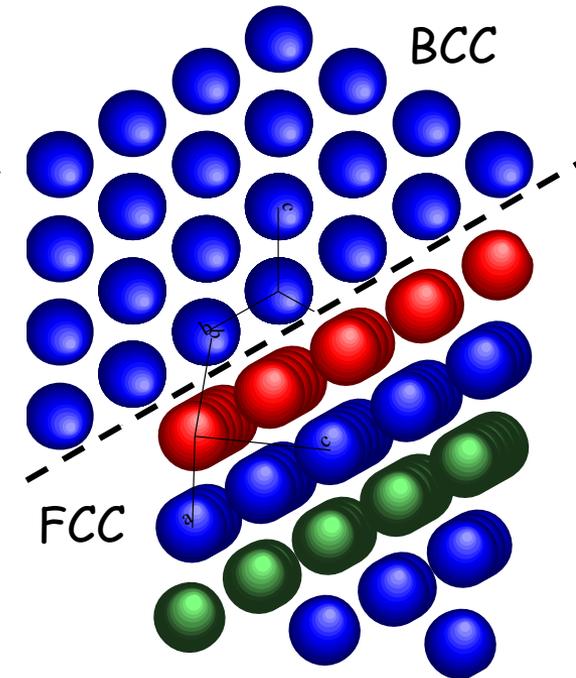
Interface viewed along the  $\langle 111 \rangle$  direction of the bcc structure



Kurdjumov-Sachs



Nischiyama-Wassermann



# Fcc <111> - bcc <110> interface also observed in Penning traps?

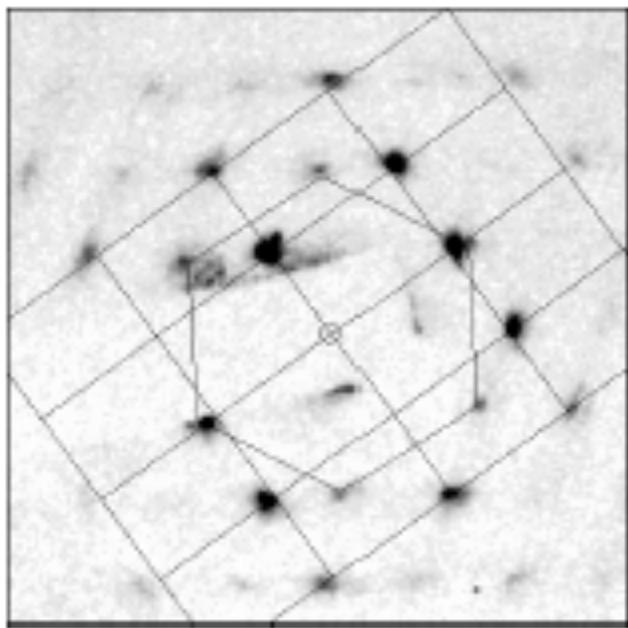


Fig. 6. Time-resolved Bragg diffraction pattern showing a superposition of twofold and sixfold symmetric patterns. The rectangular grid connects the points for which diffraction spots are predicted for a bcc lattice oriented along a <110> direction. An fcc lattice oriented along a <111> direction would generate diffraction spots at the vertices of the hexagon. The orientation of the hexagon has been adjusted to fit the data, and it differs by about 3° from that of the rectangular grid. The active timing method and the CCD camera were used. Here,  $\omega_c = 2\pi \times 70$  kHz,  $n_0 = 2.15 \times 10^{18}$  cm<sup>-3</sup>,  $N = 5 \times 10^5$ , and  $2r_0 = 2.27$  mm.

Expected orientation:

Kurdjumov-Sachs:

5.3°

Nischiyama-Wassermann

0°

Penning trap experiment

3°

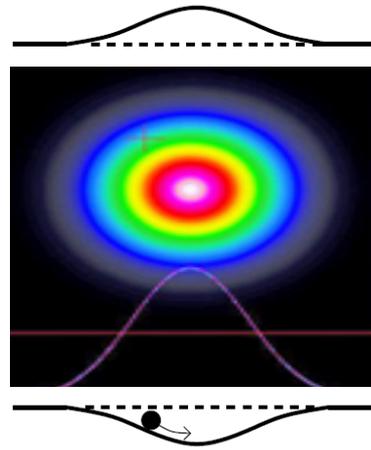
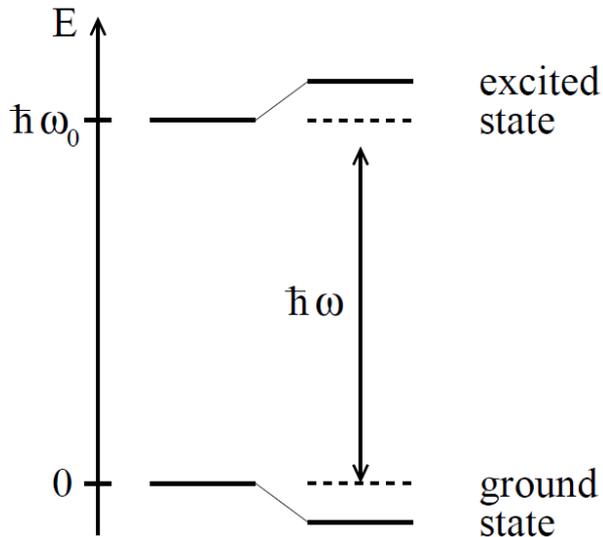
## IV. Control of crystal structures by light induced forces

# Introduction to optically induced potentials

Generally, we can polarize atoms/ions:

$$\tilde{p} = \alpha \tilde{E} \quad \Rightarrow \quad U_{\text{dip}} = -\frac{1}{2} \langle \mathbf{p} \mathbf{E} \rangle = -\frac{1}{2\epsilon_0 c} \text{Re}(\alpha) I$$

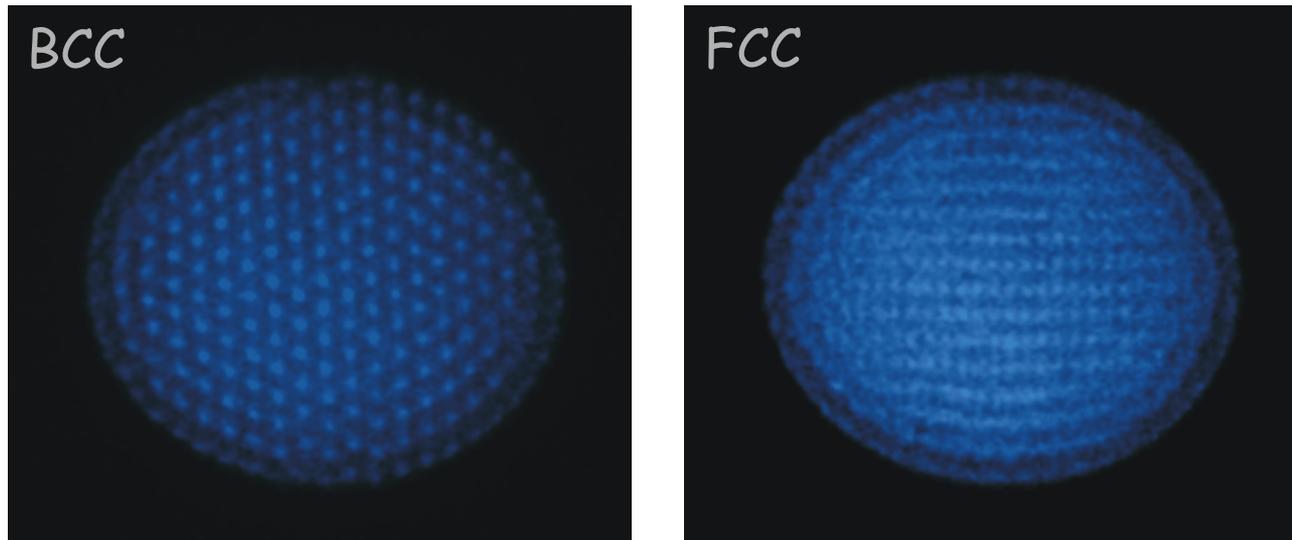
Two level atom/ion:



$$U_{\text{dip}}(\mathbf{r}) = \frac{3\pi c^2}{2\omega_0^3} \frac{\Gamma}{\Delta} I(\mathbf{r})$$
$$\Delta \equiv \omega - \omega_0$$

# Exploitation of optically induced potentials

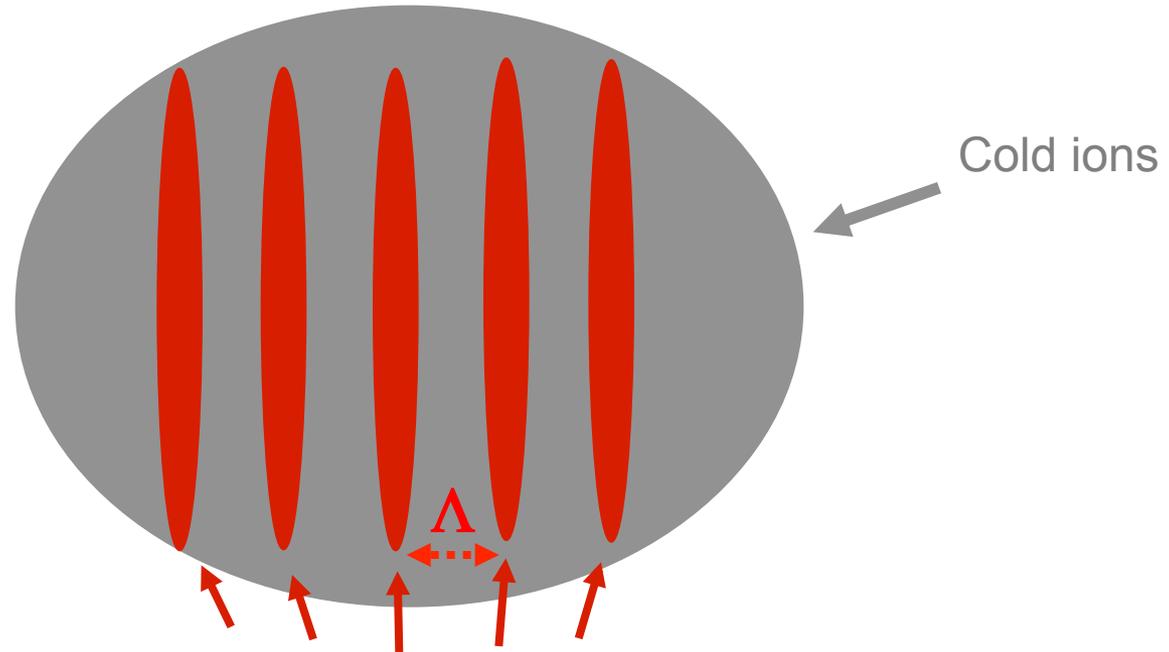
## A) Control of structural phases of Coulomb crystals



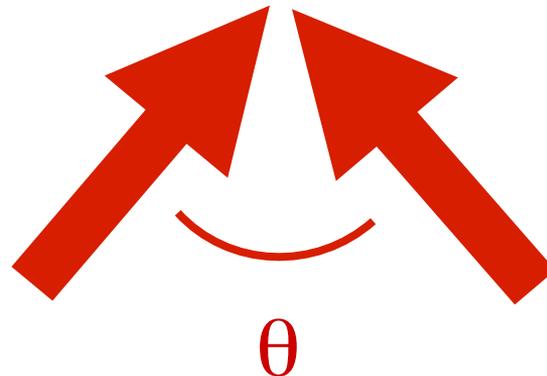
### Idea:

Try to clamp specific lattice planes by standing wave optical fields.

# Standing wave field induced potential

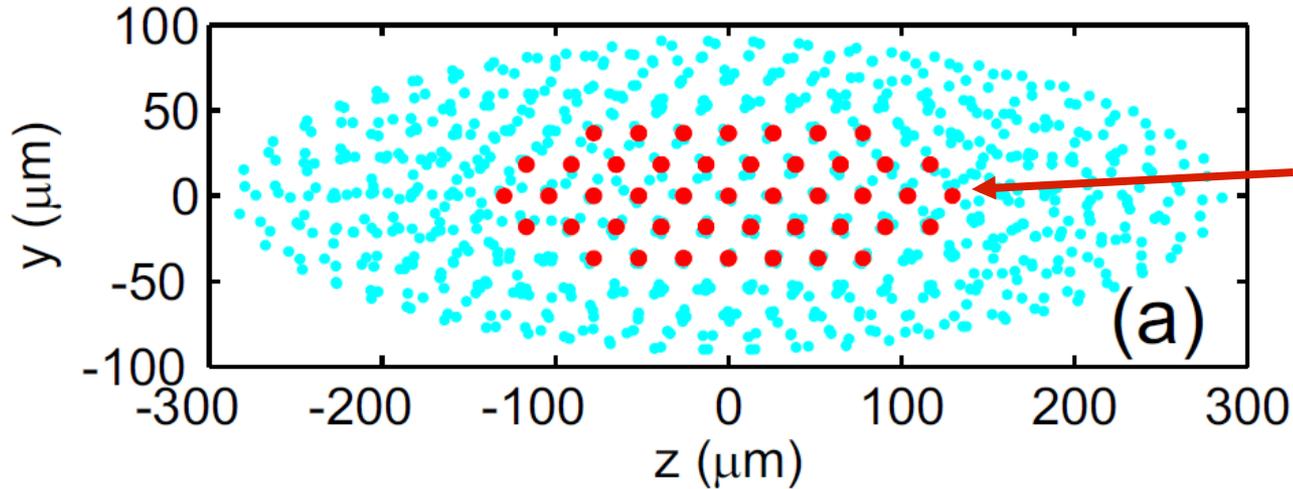


Interference pattern from two crossing light beams



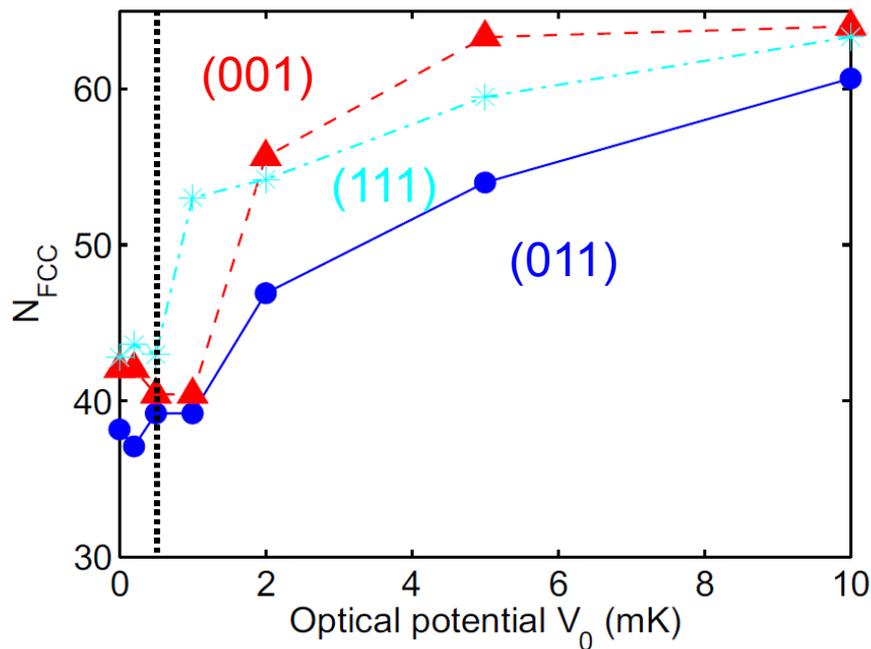
Standing wave period  $\Lambda$  can be controlled by the angle  $\theta$

# MD simulation results with 1000 ions



64 ions initially fixed in a fcc structure

Next, release the initially fixed ions, and see if they stay:

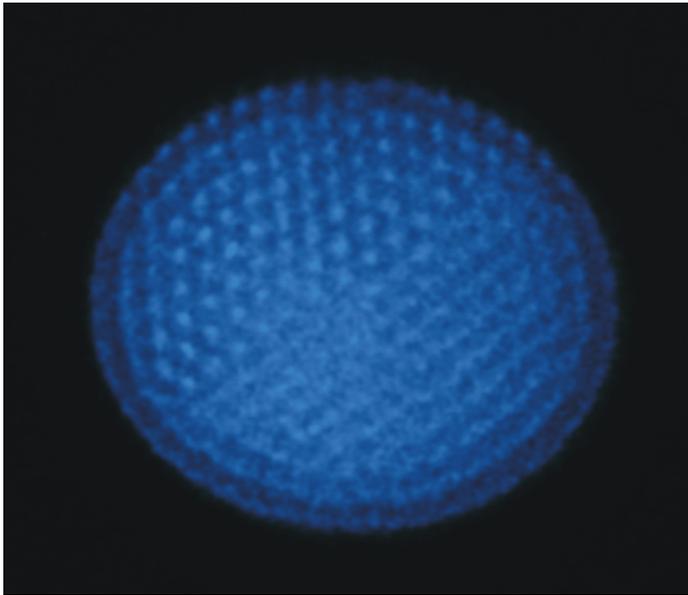


P. Horak, A. Dantan, and M. Drewsen,  
Phys. Rev. A. **86**, 043435 (2012)

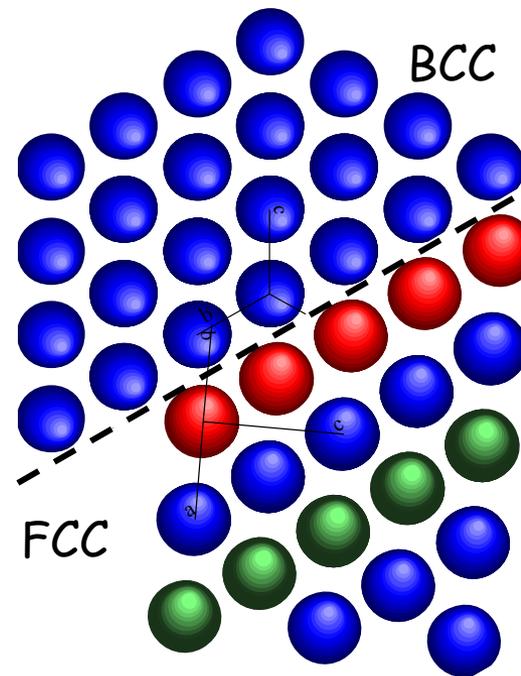
**Can we induce a structural phase transition?**

# BCC $\langle 110 \rangle$ - FCC $\langle 111 \rangle$ interface

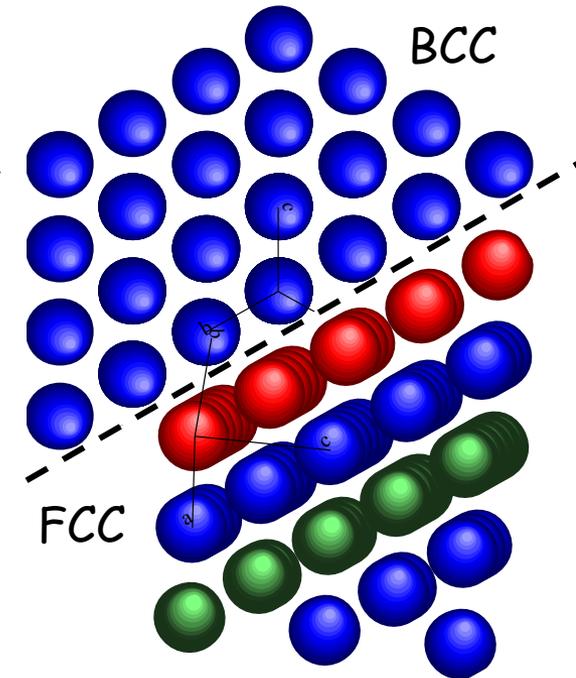
Interface viewed along the  $\langle 111 \rangle$  direction of the bcc structure



Kurdjumov-Sachs



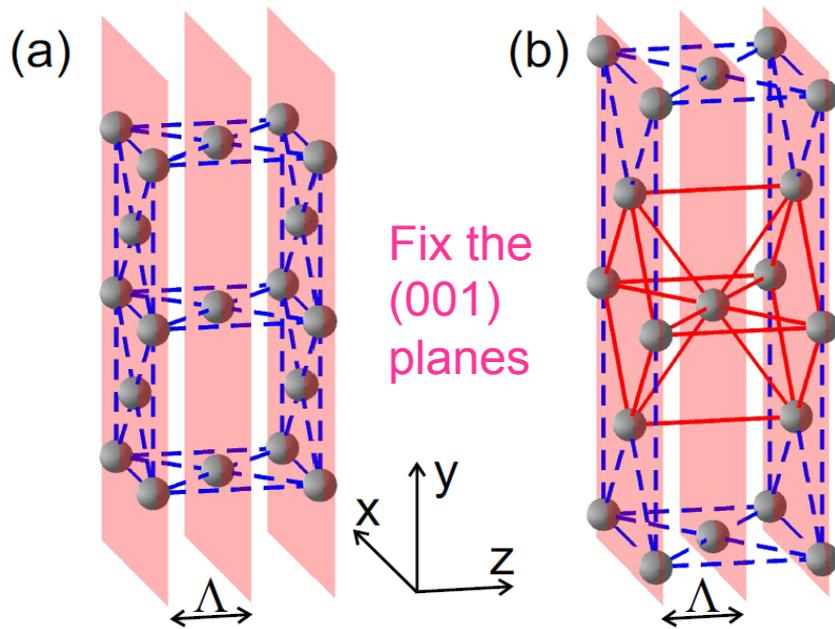
Nischiyama-Wassermann



# Can we induce a structural phase transition?

The Bain path:

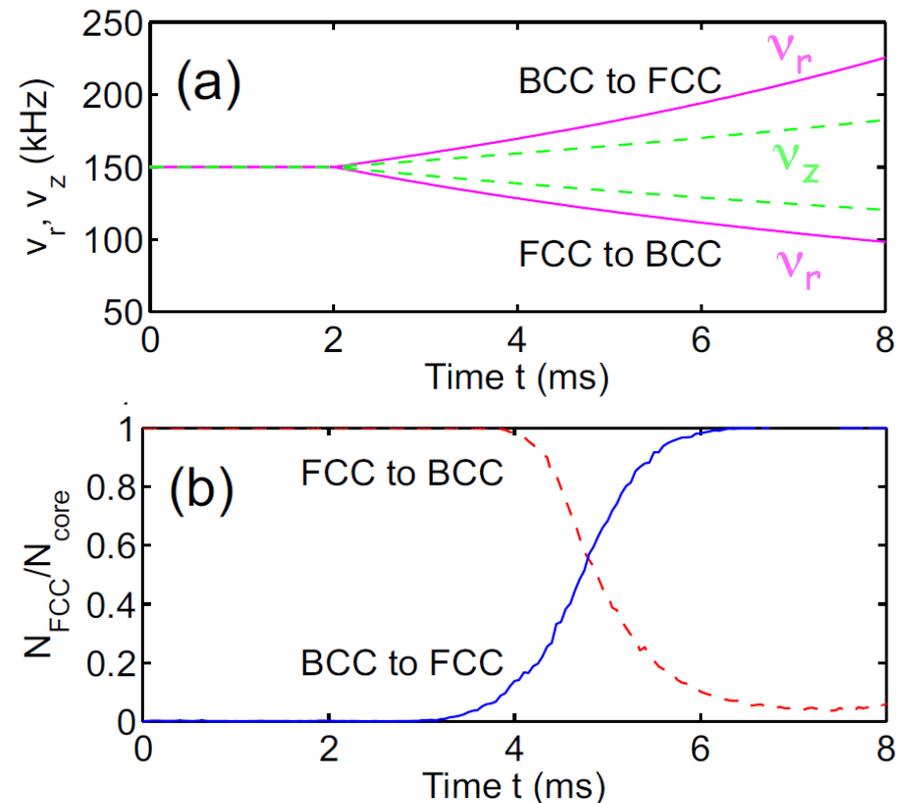
fcc  $\longleftrightarrow$  bcc



Keeps the number of particles in planes constant, but changes the density by a factor of 2 !

Can this be achieved experimentally?

Seems so:



P. Horak, A. Dantan, and M. Drewsen, Phys. Rev. A. **86**, 043435 (2012)