



The Abdus Salam
International Centre for Theoretical Physics


United Nations
Educational, Scientific
and Cultural Organization


International Atomic
Energy Agency



SMR 1666 - 10

**SCHOOL ON QUANTUM PHASE TRANSITIONS
AND
NON-EQUILIBRIUM PHENOMENA IN COLD ATOMIC GASES**

11 - 22 July 2005

Feshbach resonances and Feshbach molecules

Presented by:

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Feshbach Resonances and Feshbach Molecules



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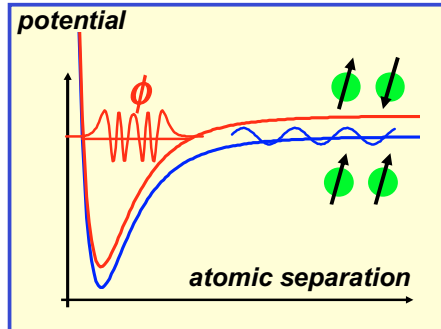


A Magnetic handle in ultracold gases



- **What is Feshbach resonance**
 - Single channel scattering
 - Two channel scattering
- **Li₂ molecules (open channel dominated)**
 - Atom-molecule thermalization
 - Collision properties of molecules
 - Precision molecular spectroscopy
 - Pairing gaps and collective modes (TOMORROW)
- **Cs₂ molecules (closed channel dominated)**
 - Highly rotation molecules (G- and L-wave)
 - Observation of Cs₄ quantum states
 - Molecule interferometry

Typical Feshbach resonances in cold atoms



Open channel
Singlet potential with a foreign bound state

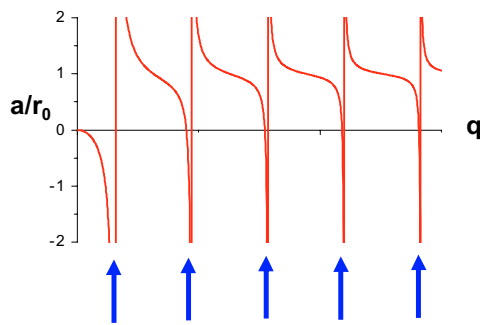
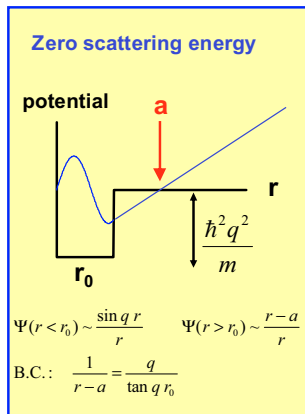
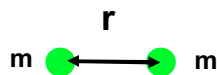
Closed channel
Triplet potential

Feshbach coupling
Hyperfine interaction
Feshbach tuning
Zeeman effect

Experimental observables

1. Inelastic and elastic cross sections
2. Mean-field shifts and many-body effects
3. Pairing of atoms

Resonant scattering (single channel)



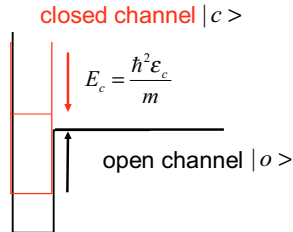
Formation of new bound state

Near the resonance, binding energy is

$$E_m = \frac{\hbar^2}{m(a_{bg} - r_0)^2}$$

Gribakin and Flambaum
PRA 48 563

Two-channel box model



$$H = \frac{\hbar^2}{m} \begin{pmatrix} -q_c^2 & \Omega \\ \Omega & -q_o^2 \end{pmatrix} \quad \text{for } r < r_0$$

$$= \begin{pmatrix} \infty & 0 \\ 0 & 0 \end{pmatrix} \quad \text{for } r > r_0$$

Exact Solution for E=0

$$\psi \sim \frac{\sin q_+ r}{r} |+\rangle + \frac{A \sin q_- r}{r} |-\rangle \quad \text{for } r < r_0$$

$$\sim \frac{r-a}{r} |o\rangle \quad \text{for } r > r_0$$

$$\frac{1}{r_0 - a} = \frac{q_+}{\tan q_+ r_0} \cos^2 \theta + \frac{q_-}{\tan q_- r_0} \sin^2 \theta$$

$|+\rangle, |-\rangle$: eigenstates for $r < r_0$

$$\tan 2\theta = \frac{2\Omega}{q_o^2 - q_c^2}$$

Approx.: $\Omega \ll q_c^2, q_o^2$ and $\epsilon_c \ll q_c / r_0$

$$\Rightarrow \frac{1}{a - r_0} = C + \frac{\gamma}{\epsilon_c}$$

Background properties: $C = \frac{1}{a_{bg} - r_0}$

Feshbach coupling: $\gamma = 2q_c^2 \theta^2 r_0^{-1}$

Magnetic dependence: $E_c + \mu B$

$$\Rightarrow a = a_{bg} \left(1 + \frac{\Delta B}{B - B_{res}} \right)$$

Feshbach bound state



Eigenstate for $E = -E_m$

$$\psi_m \sim \frac{\sin q_+ r}{r} |+\rangle + \frac{A_m \sin q_- r}{r} |-\rangle \quad \text{for } r < r_0$$

$$\sim \frac{e^{-\sqrt{\epsilon_m} a}}{r} |o\rangle \quad \text{for } r > r_0$$

Eigen energy $E_m = \frac{\hbar^2}{m} \epsilon_m$

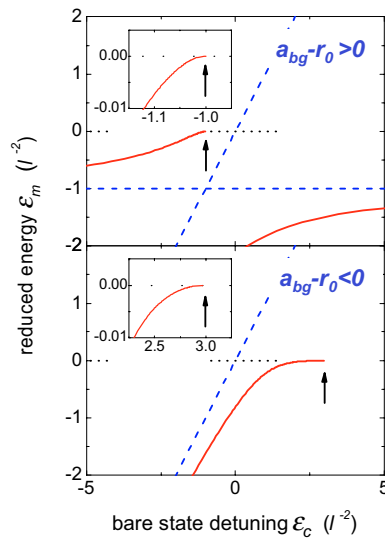
$$(\epsilon_m + \epsilon_c) \left(\sqrt{\epsilon_m} - \frac{1}{a_{bg} - r_0} \right) = \gamma$$

Mixing fraction

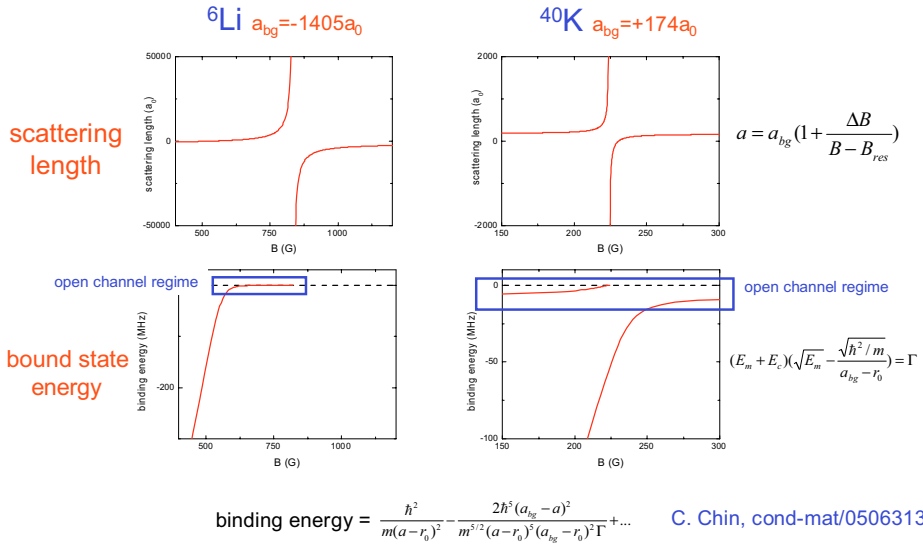
$$\psi_m = \sin \phi |c\rangle + \cos \phi |o\rangle$$

$$\sin^2 \phi = \int |\psi_c(r)|^2 dv = \frac{2\sqrt{\epsilon_m} \gamma}{(\epsilon_m + \epsilon_c)^2 + 2\sqrt{\epsilon_m} \gamma}$$

generic energy structure



Scattering length and bound state



Example: Li_2 and K_2 molecules

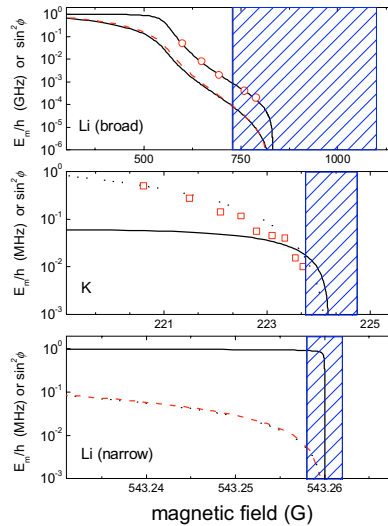


Excellent agreement with all calculations and measurements!!

• Small binding open channel

$$E_m = \frac{\hbar^2}{m(a_{bg} - r_0)^2}$$

Li (broad)
mixing frac. $< 10^{-3}$
K40
mixing frac. $< 10^{-1}$
Li (narrow)
mixing frac. ~ 1



Comparison:

Binding energy (Innsbruck)
Q scattering calc. (NIST)
Mixing fraction (Rice Univ.)

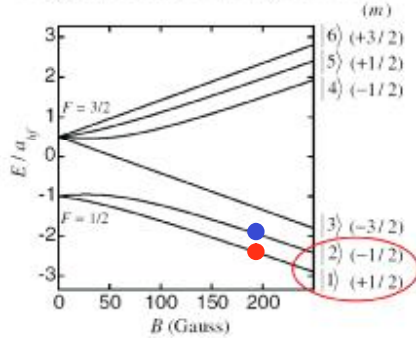
Binding energy (JILA)
Q scattering calc. (NIST)

Position & width (Rice Univ.)
Binding energy (NIST)

⁶Li experiment

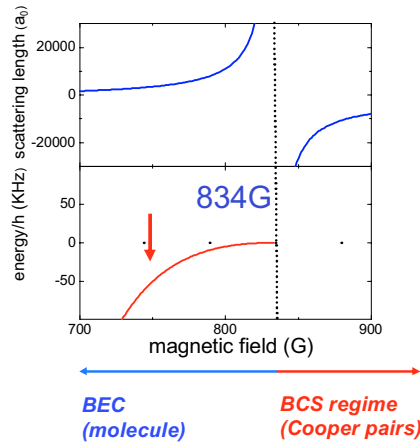


⁶Li ground state in a magnetic field



At high fields,
 $|2\rangle = |2s, m_s=-1/2, m_l=0\rangle$
 $|1\rangle = |2s, m_s=-1/2, m_l=-1\rangle$

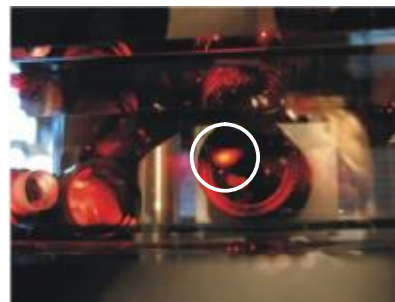
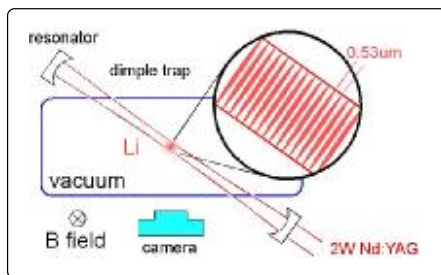
scattering between $|1\rangle$ and $|2\rangle$



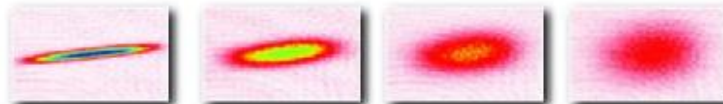
Li Experimental Setup



- Resonator set up at Brewster's angle. *Opt. Lett* 26, 1837 (2001)
- Efficient loading into the optical trap: $\sim 10^7$ thermal atoms



Absorption images



Time Evolution

Atoms ↔ molecules: thermalization

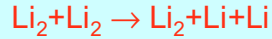


B = 690 G:
 mol. bind. energy $E_m = k_B 8\mu\text{K}$
 \gg therm. energy $k_B T = k_B 2.5\mu\text{K}$

Atom → molecule: recombination



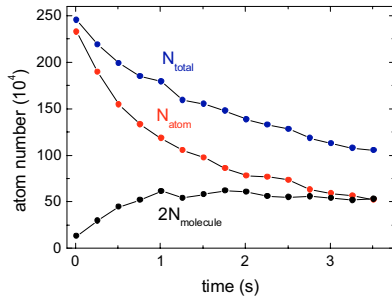
Molecule → Atom: coll. dissociation



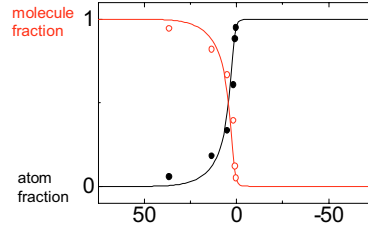
Thermal equilibrium

$$N_M / N_A = \phi_A e^{E_m/kT}$$

ϕ : phase space density



Exp: S. Jochim et al. PRL 91, 240402 (2003)



Theo: C. Chin and R. Grimm, PRA (2004)

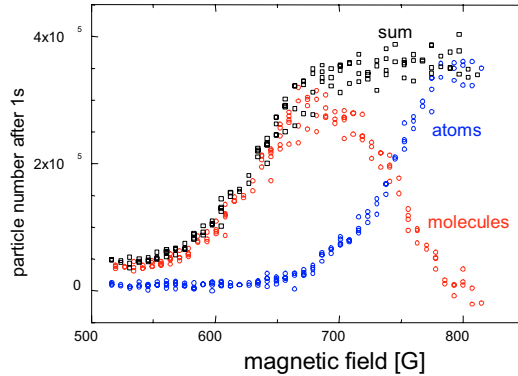
Collision properties of molecules



Small a
 Large binding energy

molecule decay atom-molecule conversion

Large a
 Small binding energy



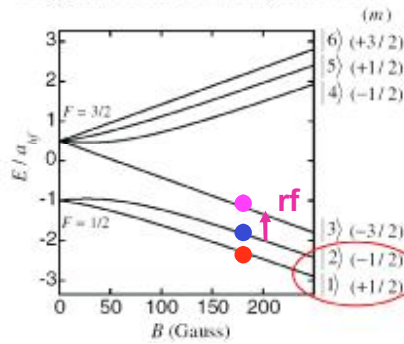
Experiment:
 Long lifetimes near the resonance!!
 > 40s at 764G
 < 100ms below 500G

Theory:
 decay rate: $C r_e (r_e/a)^{2.55}$
 Mol. scattering length $a_m = 0.6a$
 Petrov et al '04

Molecular binding energy: RF spectroscopy



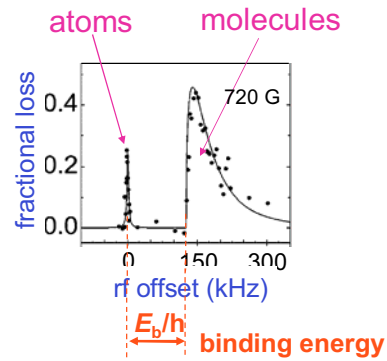
^8Li ground state in a magnetic field



State 3 is initially unpopulated

$$f_{\text{atom}} = E_{23}$$

$$f_{\text{molecule}} = E_{23} + E_b + 2E_K$$

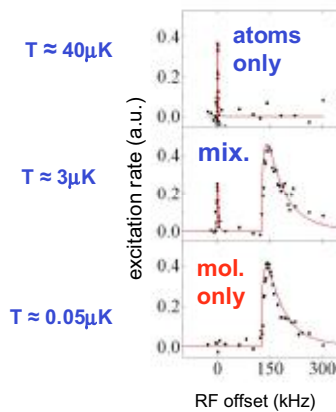


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Applications of RF spectroscopy



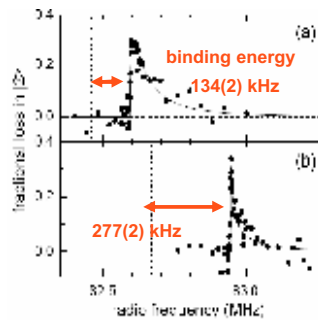
1. Distinguish atoms and weakly bound molecules



Also by Jin group at JILA

2. Precision molecular spectroscopy

Cooperation: A. Simoni, E. Tiesinga and P. Julienne, NIST



Experiment:
M. Bartenstein et al,
PRL **94**, 103201
(2005)

Lineshape theory
C. Chin and P. Julienne
PRA **71**, 012713
(2005)

3. Observation of Fermionic pairing gap

Outside the scope of this talk.

Please see
C. Chin et al.,
Science **04**

Bound-free and bound-bound transitions



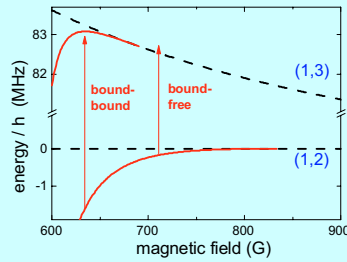
Bound-free transitions

$$\Gamma_f = \frac{h\Omega}{2} \left| \int \psi_K^*(r) \phi_m(r) dr \right|^2$$

$$= \frac{4}{\hbar} \left(1 - \frac{a'}{a}\right)^2 \frac{K^{1/2} E_b^{1/2} E_b^*}{(K + E_b)^2 (K + E_b^*)}$$

Bound-bound transitions

$$\Gamma_b = \frac{h\Omega}{2} \left| \int \phi_m^*(r) \phi_m(r) dr \right|^2 = \left(\frac{a-a'}{a+a'}\right)^2$$



Precision determination of scattering properties without collisions!!

EXPERIMENT:

Bound-free transitions

720.13(4)G RF:82.593(2)MHz

694.83(4)G RF:82.944(2)MHz

Bound-bound transitions

676.09(3)G RF:83.2966(5)MHz

661.44(2)G RF:83.6645(3)MHz



THEORY (Simoni, Tiesinga, Julienne)

$a_s = 45.167(8)a_0$, $a_1 = -2140(18)a_0$

Feshbach resonance positions

$|1\rangle + |2\rangle$: 834.1 ± 1.5 G

$|1\rangle + |3\rangle$: 690.4 ± 0.5 G

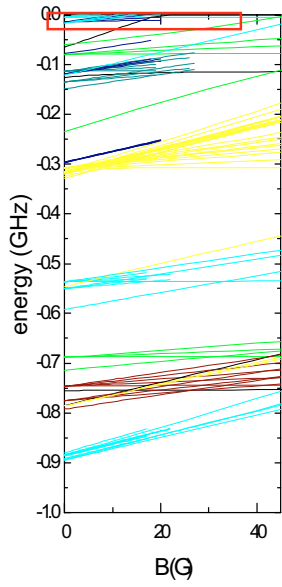
$|2\rangle + |3\rangle$: 811.2 ± 1.0 G

Brief summary on Li₂ molecule

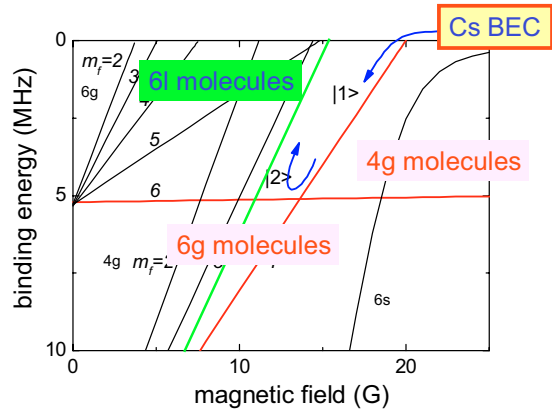


- **Molecule formation by recombination and thermal equilibrium**
 - Efficient atom-molecule conversion when $kT \ll E_m$
- **Collisional stability with positive scattering length**
 - Evaporation to BEC and BEC-BCS crossover (tomorrow)
- **Precision molecular RF spectroscopy**
 - Precise determination of the molecular energy structure
 - Pairing gap (tomorrow)
- **Collective modes and Fermionic hydrodynamics (tomorrow)**

Experiments on Cs₂ molecules



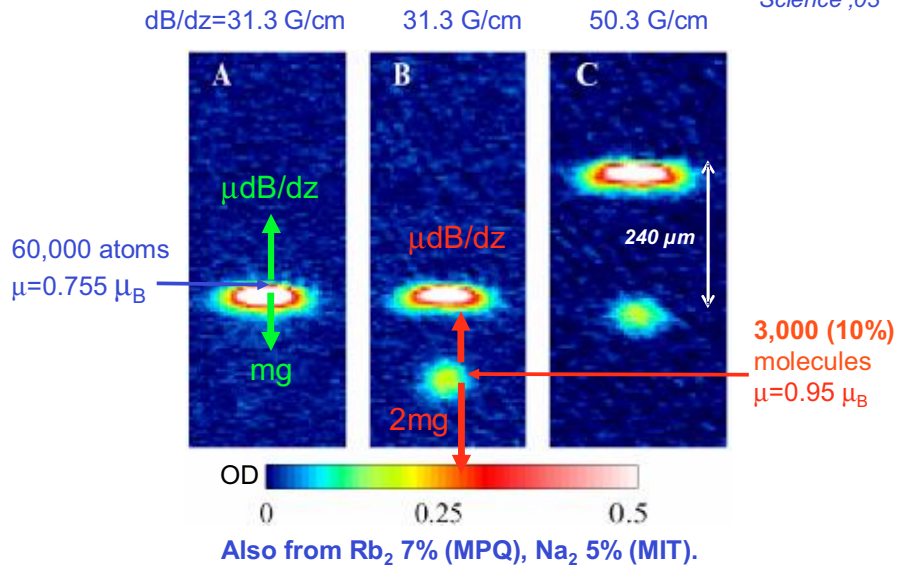
G-wave molecules: orbital angular momentum $L = 4 \hbar$
 G-wave resonance: coupling which changes L by $4\hbar$
 due to high order spin-orbit interaction



Stern-Gerlach separation (5/2003)



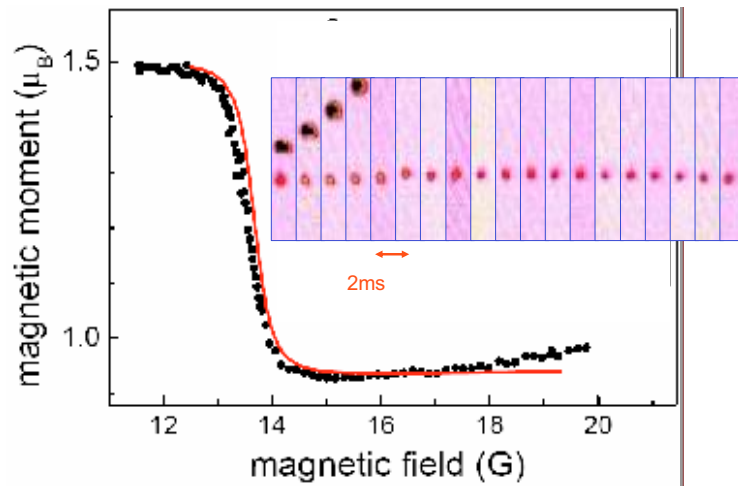
Science, 03



mapping out the avoided crossing

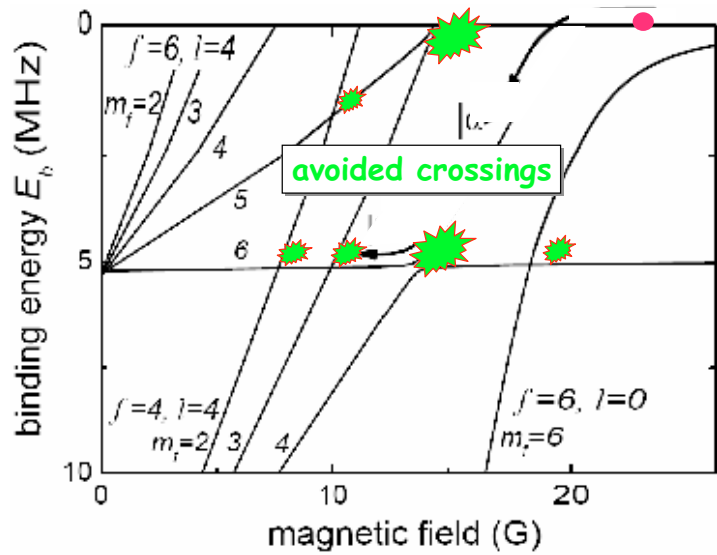


levitation condition \Rightarrow magnetic moment (precision 1%)



theory by E. Tiesinga et al., NIST

Convert 4g molecules to 6g molecules

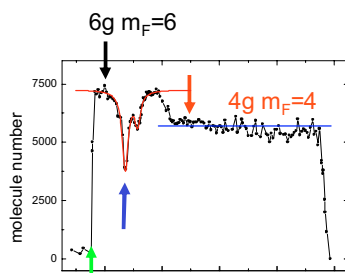


data from NIST group

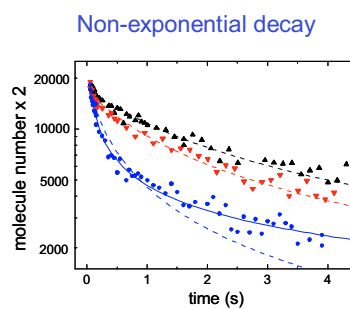
Feshbach resonances of Cs₄ molecules



- Leave the molecules in the trap for 300ms.



What happened here?

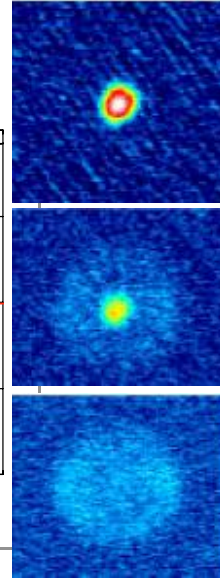
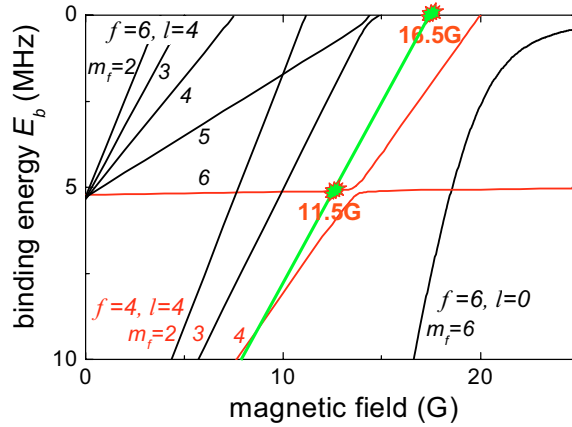


Binary decay constant:		density dependence resonance structure ⇒ molecular Feshbach resonances induced by Cs ₄ states
6g m _F =6	3×10 ⁻¹¹ cm ³ /s	
4g m _F =4	5×10 ⁻¹¹ cm ³ /s	
on resonance (unitarity limit)	2×10 ⁻¹⁰ cm ³ /s 4×10 ⁻¹⁰ cm ³ /s	

puzzling observations



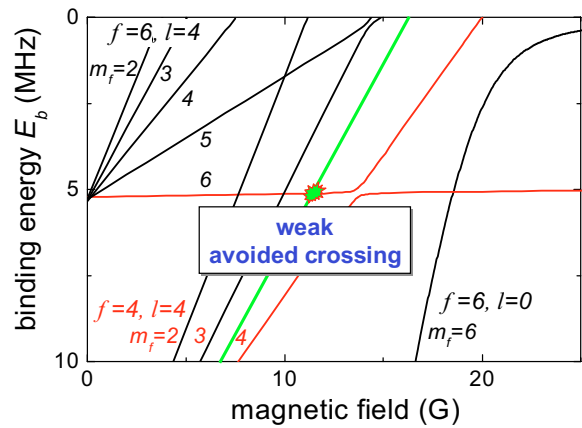
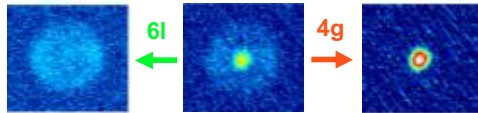
The mysterious state dissociates at 16.5G!!



interesting experimental tools



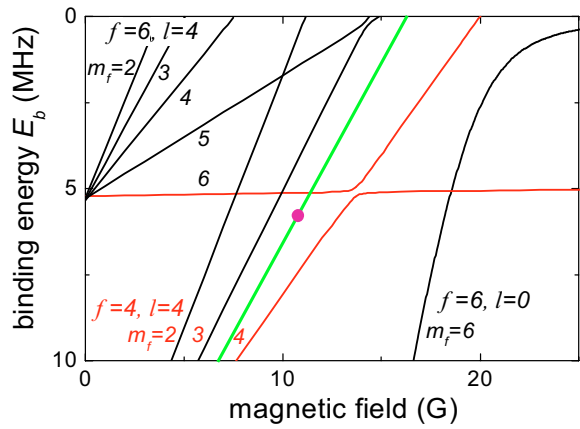
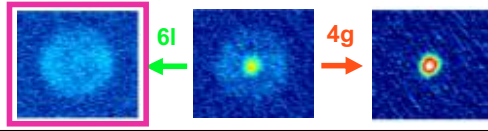
detection method for dissociation channel



first experiment



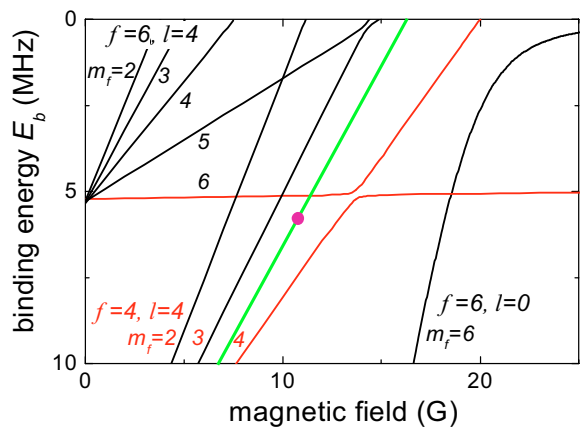
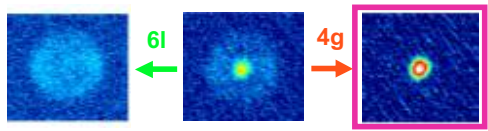
detection method for dissociation channel



first experiment



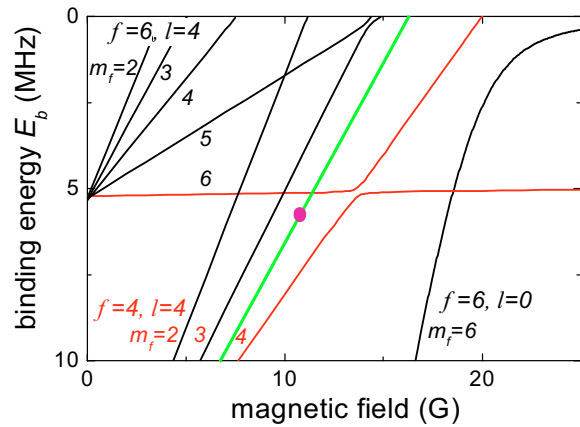
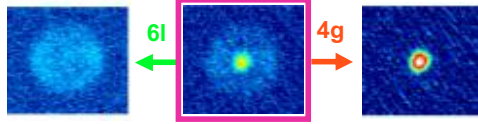
detection method for dissociation channel



first experiment



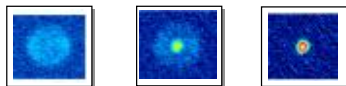
detection method for
dissociation channel



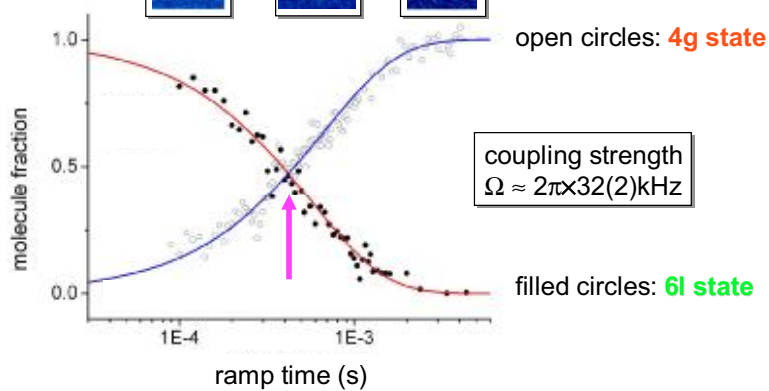
Landau-Zener → coupling strength



fast

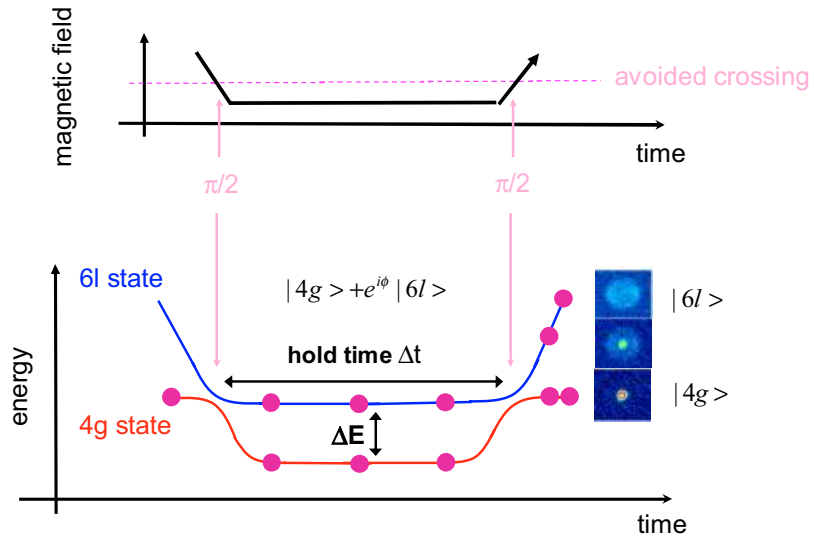


slow

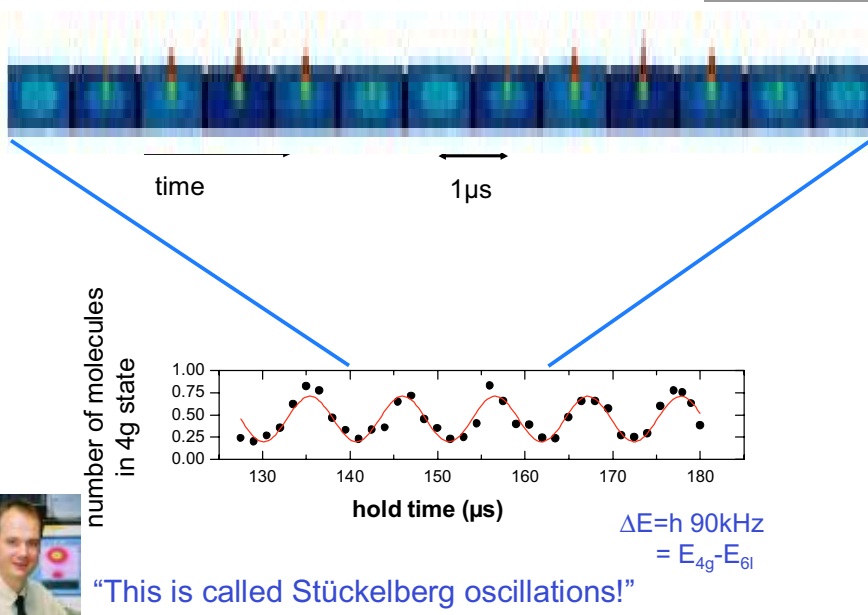


molecular beam splitter: coherence ?

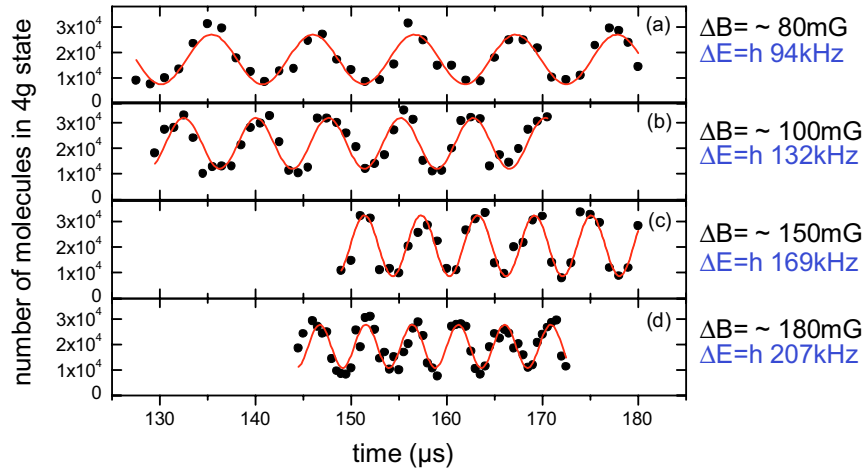
interferometer



After some struggle... It worked!!



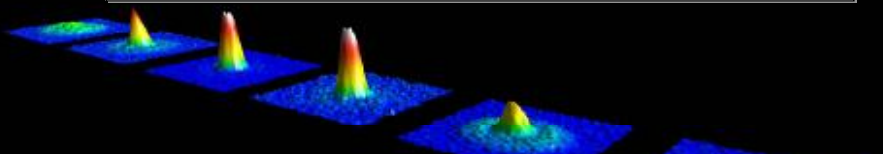
Stückelberg oscillations



molecule interferometer for future precision experiments !

Conclusion

Efficient creation of S-, G- and L-wave molecules
Precision determination of $\Psi(\mathbf{r})$, E_b , μ , α of molecules
Tetramer states and molecule interferometry
Molecules in many-body regime.



Cold molecules roadmap

Formation \rightarrow single molecule \rightarrow two- and many-body \rightarrow applications



What is this mysterious state?



Date: Mon, 10 May 2004 14:41:24 -0400

From: Cheng Chin [<mailto:Cheng.Chin@uibk.ac.at>]

Sent: Wednesday, November 10, 2004 16:09

To: Paul S. Julienne; Eite Tiesinga

Cc: Rudolf Grimm; Charles W. Clark; Christoph Naegerl

Subject: s p d f g h i k l m n o q r t u v w x y z

Hi folks,

I found the reference for the spectroscopy notation in
Phys. Rev. 33, 900.906 (1929), "Report on Notation for Atomic Spectra"
by H. N. Russell, A. G. Shenstone, and Louis A. Turner

It is **s p d f g h i k l m n o q r t u v w x y z**.
There is no j as Charles said.

.....

So we have **ultracold L-wave molecules**. ☺