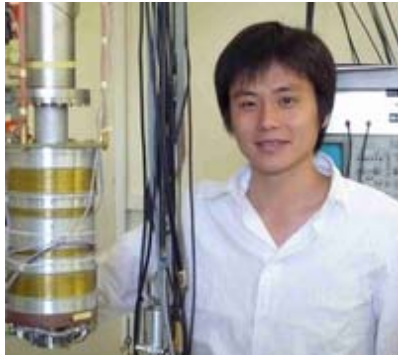


NMR study of the pressure induced Mott transition to superconductivity in the two phases of Cs_3C_{60}



Y. Ihara

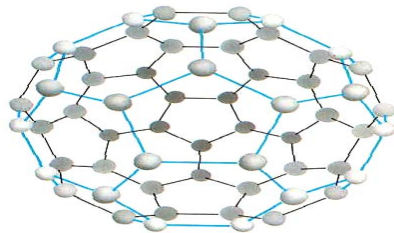


H. Alloul

P. Wzietek

Laboratoire de Physique des Solides, Université Paris XI, Orsay, France.

Daniele Pontiroli, Marcello Mazzani, Mauro Riccò.
CNISM and Dipartimento di Fisica, Università di Parma

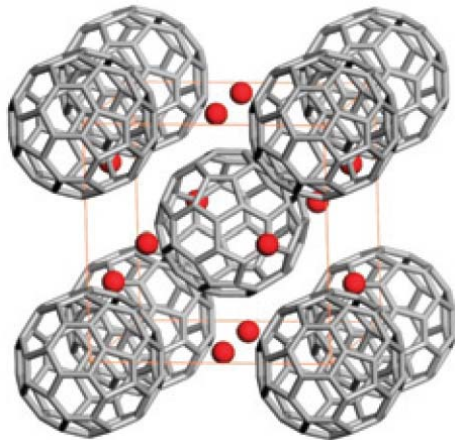


V. Brouet, H. Alloul *et al*
PRL, **82**, 2131 (1999); **86**, 4680 (2001);
PR B, **66**, 155122, 15123, 15124 (2002).

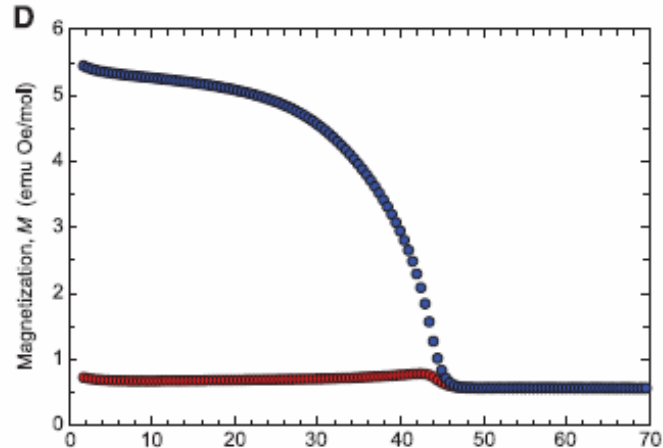


V. Brouet

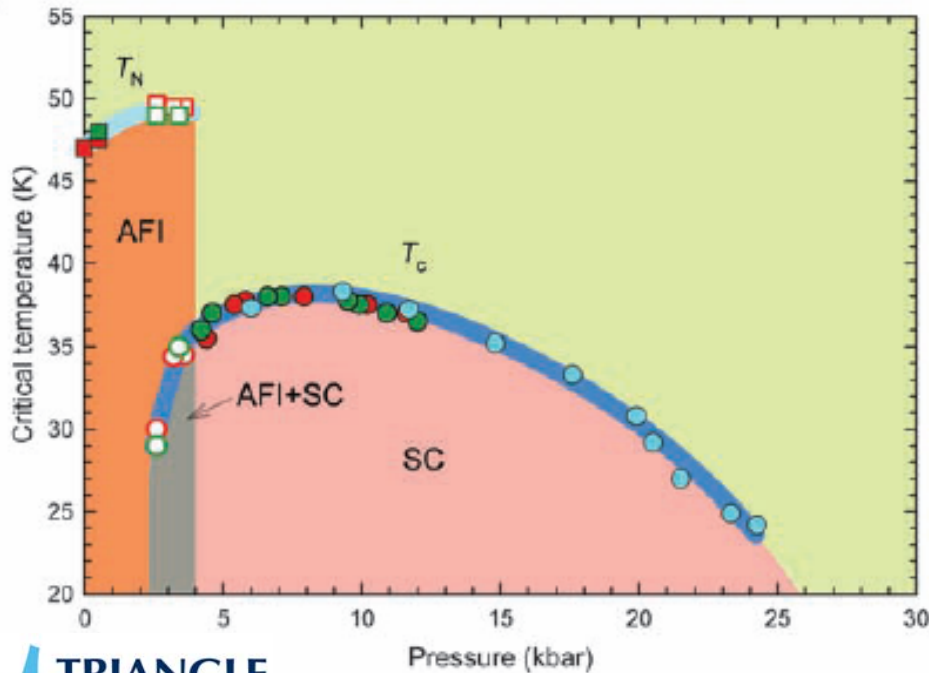
A15 Cs_3C_{60} : magnetic at ambient pressure



M



T



**SC induced by pressure
Proximity of magnetism
and SC.**

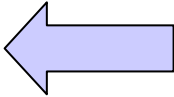
Palstra *et al.* Sol. Stat. Commun. **93** 327 (1995).

A Ganin *et al.* Nature Materials, April 2008

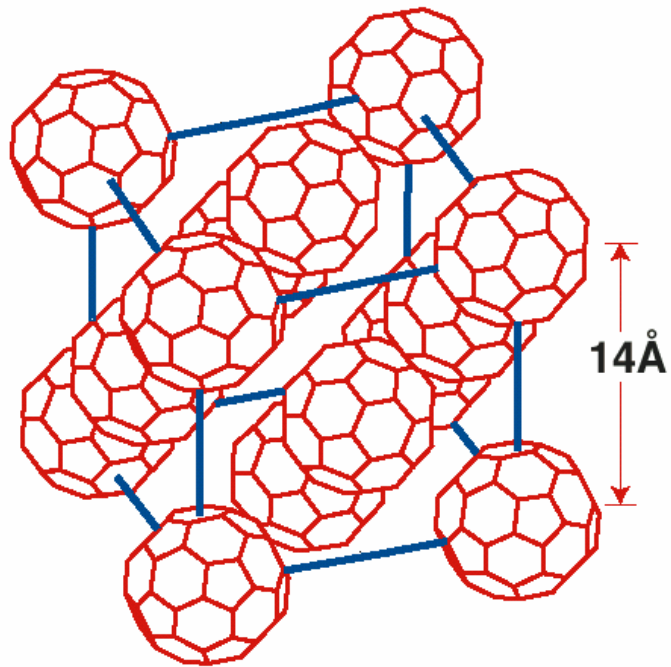
Takabayashi et al., Science 2009

NMR study of the pressure induced Mott transition to superconductivity in the two phases of Cs_3C_{60}

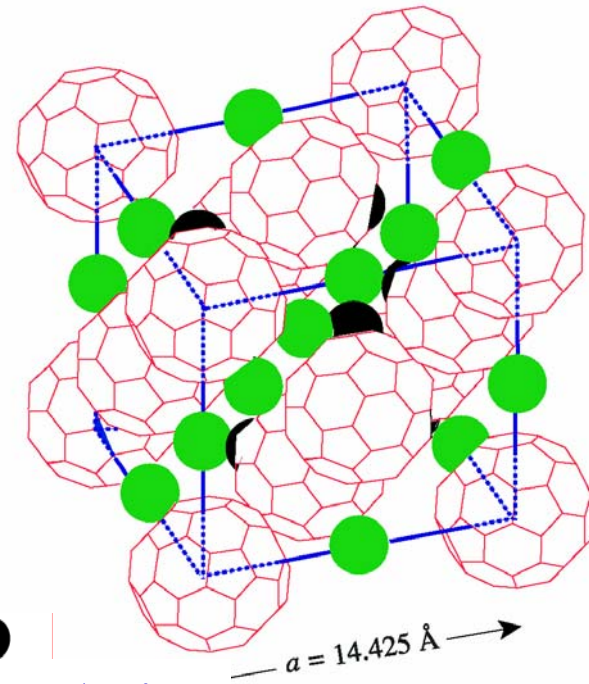
- **Introduction: A_3C_{60} and their superconductivity**
 - **LDA Electronic structure**
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 - **A_1C_{60} Mott insulator?**
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- **Expanded magnetic compounds and Cs_3C_{60} phases**
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sc and fcc phases of fullerides



Solid C₆₀



● | tetrahedral A site

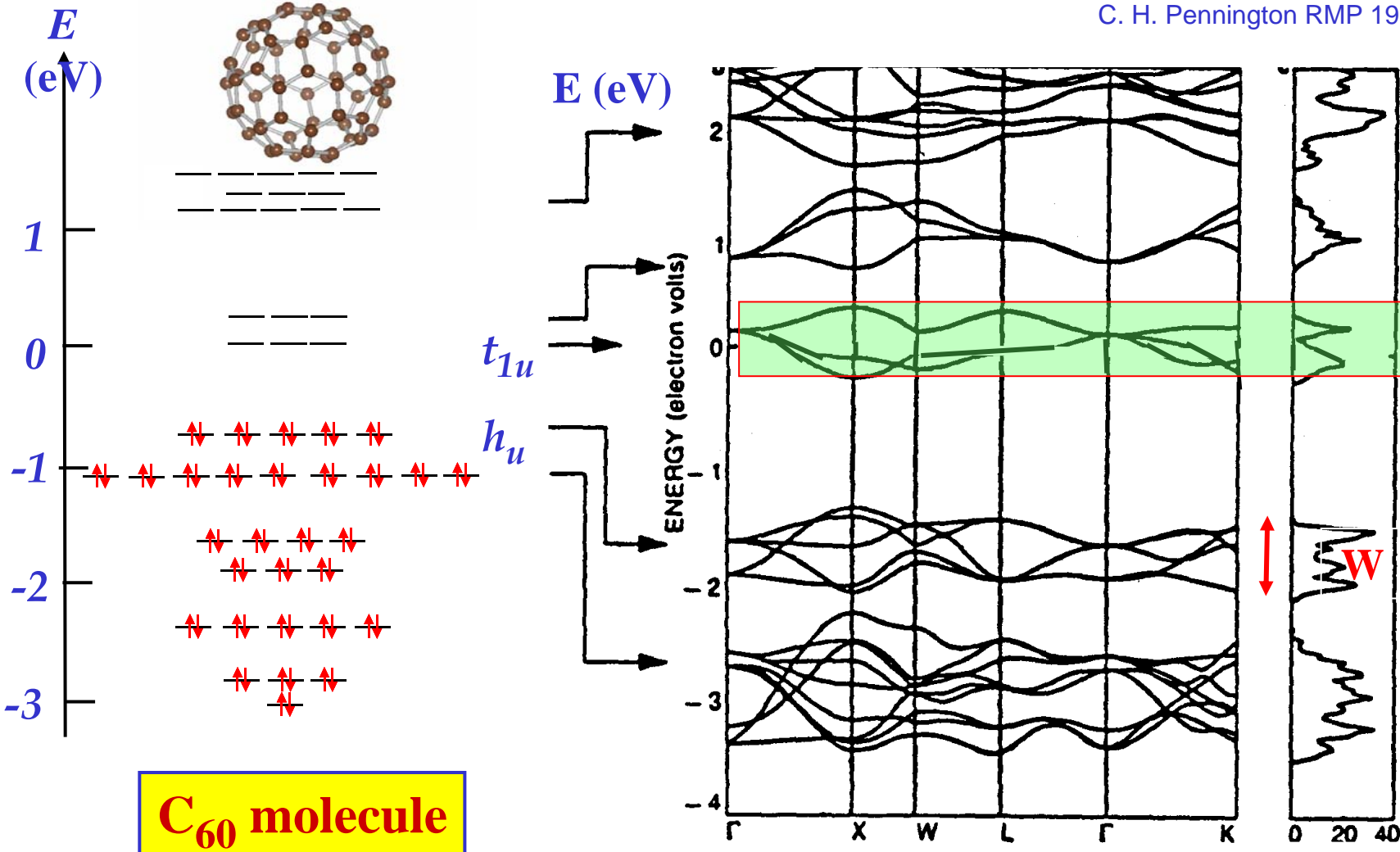
● | octahedral A site



High T_c
superconductors

Electronic structure of doped fullerides

C. H. Pennington RMP 1995.

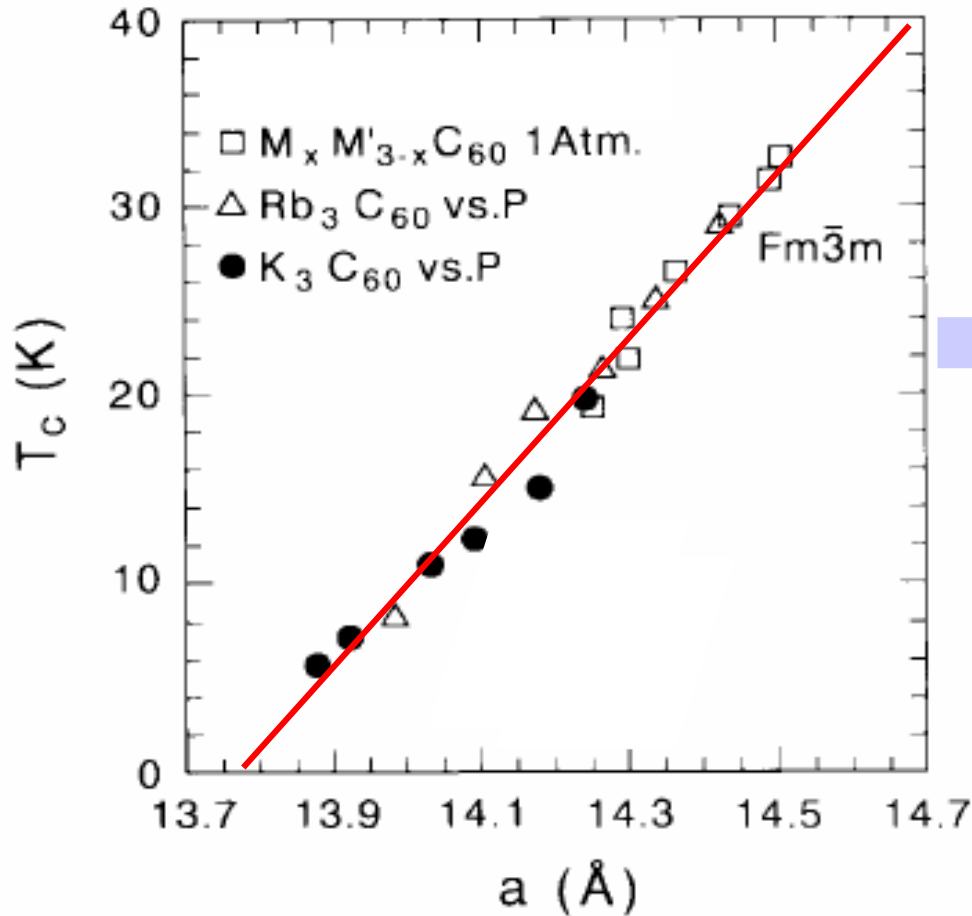


C₆₀ molecule

C₆₀ solid

W = 0.5 eV

Superconductivity in A_3C_{60}



$$k_B T_c = 1.14 \hbar \omega_D \exp \left[- \frac{1}{V_0 n(E_F)} \right]$$

T_c depends only on $n(E_F)$

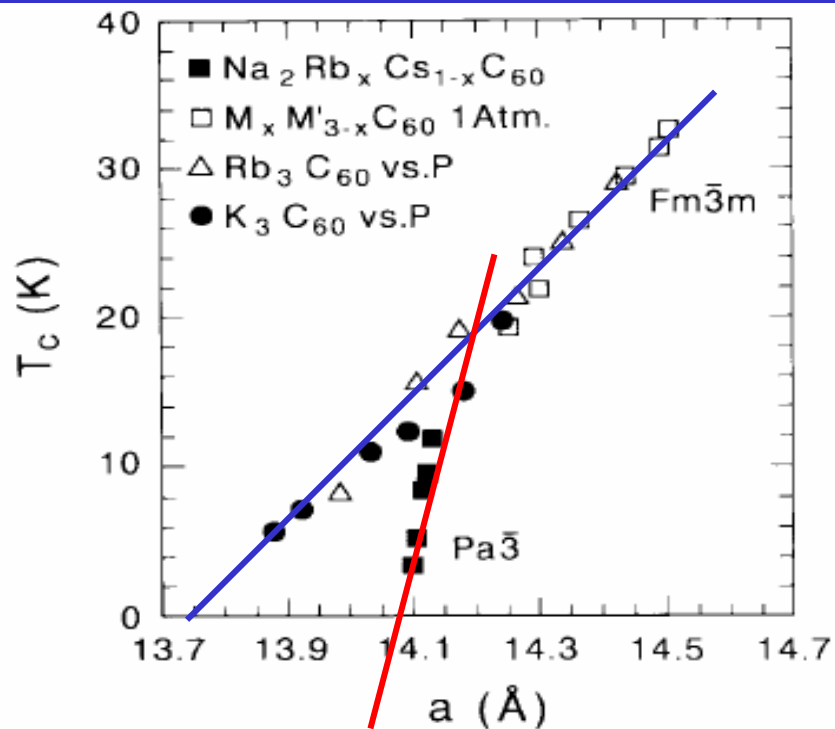
$\hbar \omega_D$ and V_0

are molecular properties

**Only C_{60} phonons
are involved**

The scaling between T_c and the lattice parameter is assumed to support a BCS-like mechanism

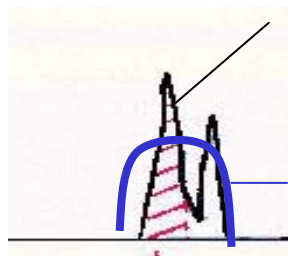
Difference between $\text{Na}_2\text{AC}_{60}$ and the other A_3C_{60}



Yildirim *et al.* MRS proceedings **359** 273 (1995).

$\text{Na}_2\text{AC}_{60}$ sc Pa3 structure
other A_3C_{60} are fcc

- Different slope for sc phases ?
- Different variation of $n(E_F)$?
- Alkali plays a role after all ?



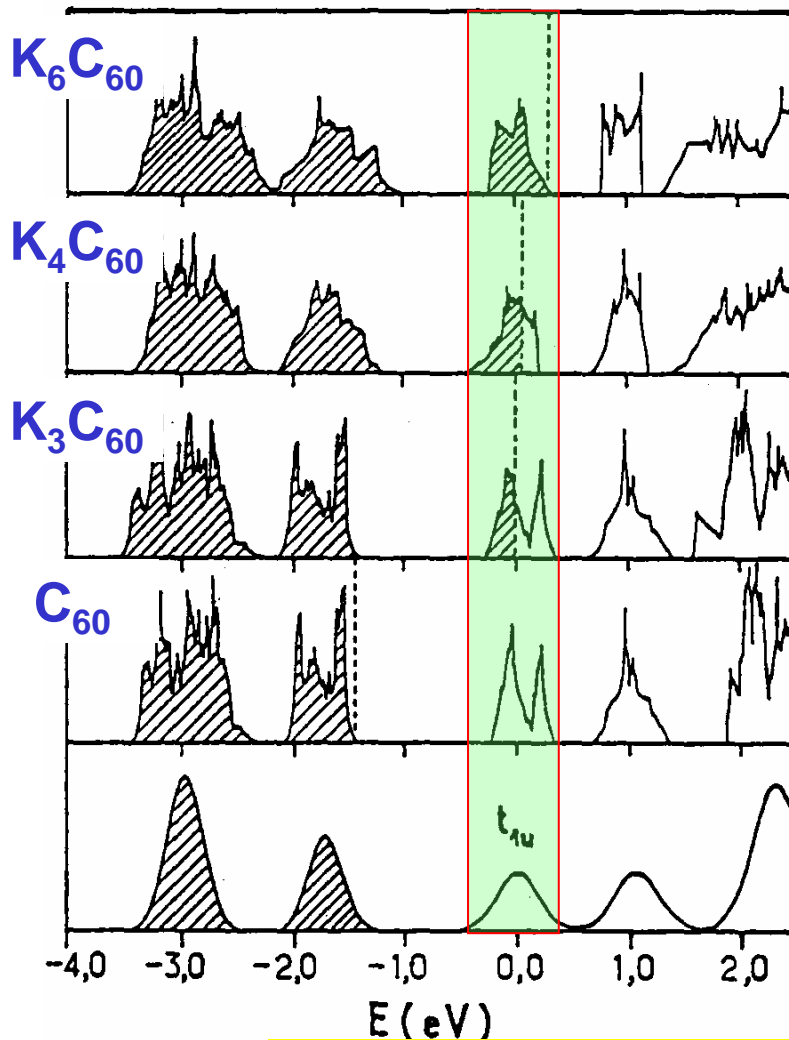
ordered

disordered

What happens then for other n values ?

The anomalous properties of A_nC_{60} compounds

t_{1u} band



Electronic doping of triply degenerate t_{1u} orbitals

All compounds with $n < 6$ should be metallic.

But...

Crystalline symmetry depends on doping level.

$x = 4$: Insulator (bct)

K_4C_{60} , G. Zimmer EPL 1994, V. Brouet PRB 2002.

$x = 3$: Superconductor (fcc, cubic)

$x = 2$: Insulator (cubic)

Na_2C_{60} , F. Rachdi PRB 1997, V. Brouet PRL 2001, PRB 2002.

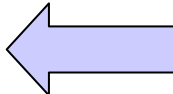
$x = 1$: Metal-Insulator transition

polymer- Cs_1C_{60} , V. Brouet PRL 1996, B. Simovic Synth. Met. 1999.

cubic- Cs_1C_{60} , V. Brouet PRL 1999.

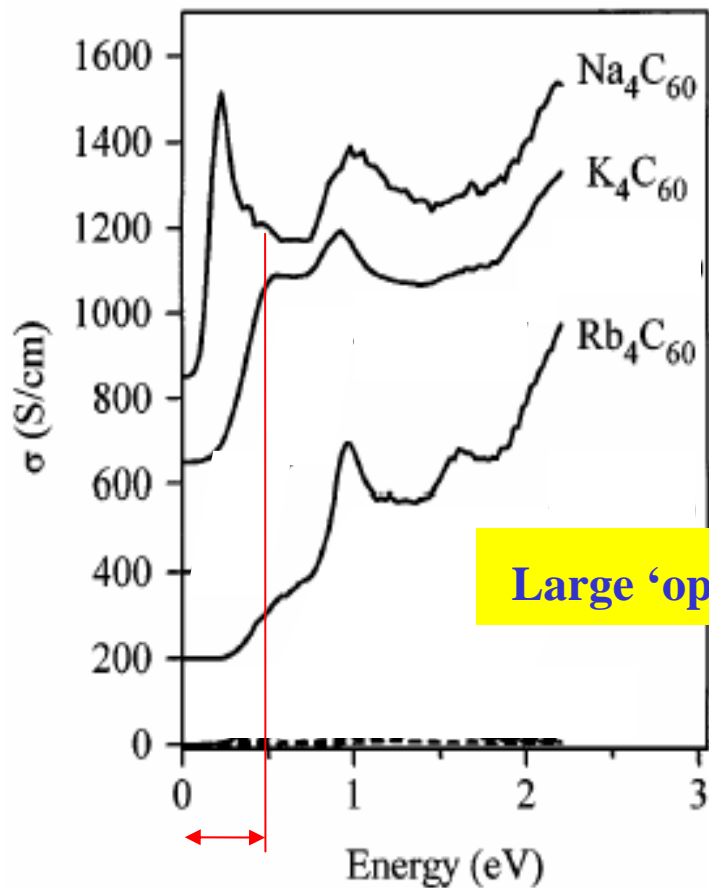
Strong correlations and Jahn-Teller distortions (JTD)

NMR study of the pressure induced Mott transition to superconductivity in the two phases of Cs_3C_{60}

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Electronic properties of A_4C_{60} (bct) and Na_2C_{60} (cubic)

Optical conductivity σ extracted by EELS

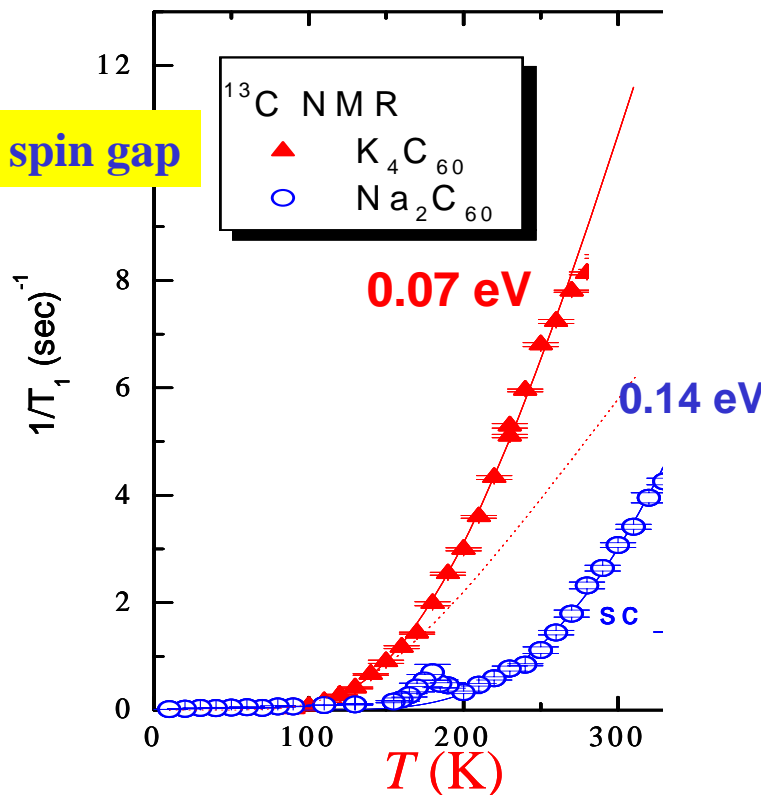


0.5 eV

M. Knupfer *et al*, PRL **79**, 2714 (1997).

NMR $1/T_1$

Small spin gap



K_4C_{60} : G. Zimmer *et al* EPL (1994),
V. Brouet *et al*, PRB 66 155122
(2002). Na_2C_{60} , F. Rachdi *et al*, PRB 1997,
V. Brouet *et al* PRL 2001. PRB 2002.

Insulating non magnetic ground state

Stabilization of the singlets

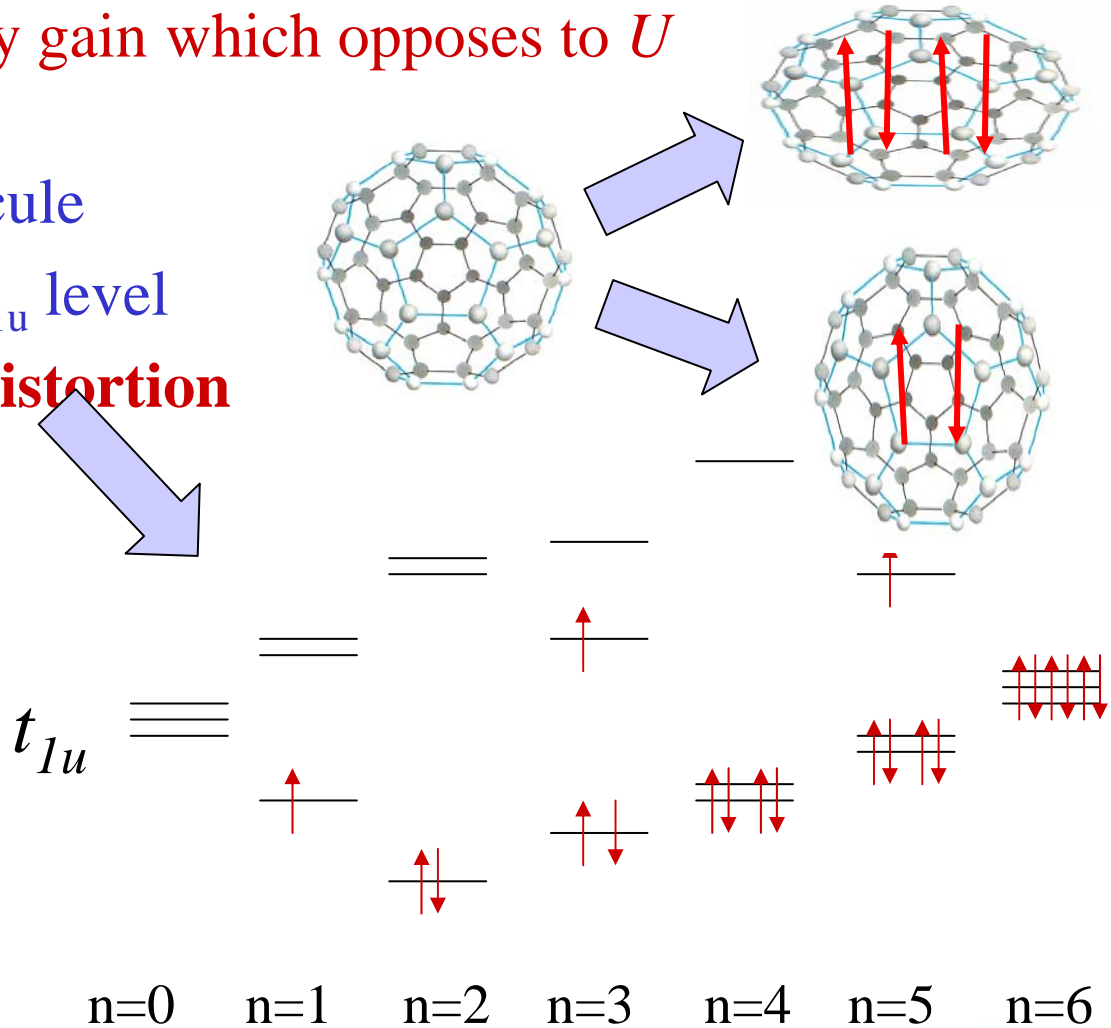
Why singlets? Two electrons on a ball costs an energy U
 So there is an energy gain which opposes to U

For a charged molecule

The degeneracy of the t_{1u} level is lifted by a **Jahn Teller distortion**

The Jahn-Teller splitting of the t_{1u} level depends on the C_{60} charge

Manini, Tosatti PRB94

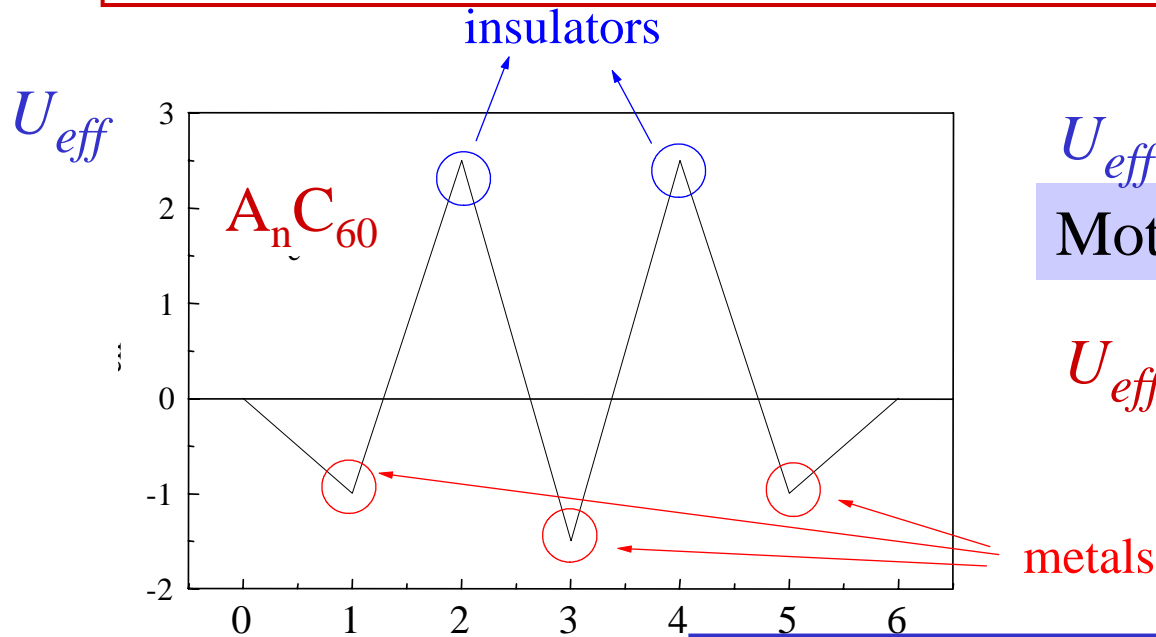


The larger gain per electron is for $n=2$ or 4

Strong electronic correlations induce the insulating behaviour

In all A_nC_{60} the coulomb repulsion U is large

The Jahn-Teller distortion adds a contribution U_{eff} to U
Yields a $S=0$ ground state rather than $S=1$ for $n = 2$ or 4



U_{eff} adds to U for even n

Mott Jahn-Teller insulators

U_{eff} reduces U for odd n

metals

**Effective electronic interactions
mediated by Jahn-Teller distortions**

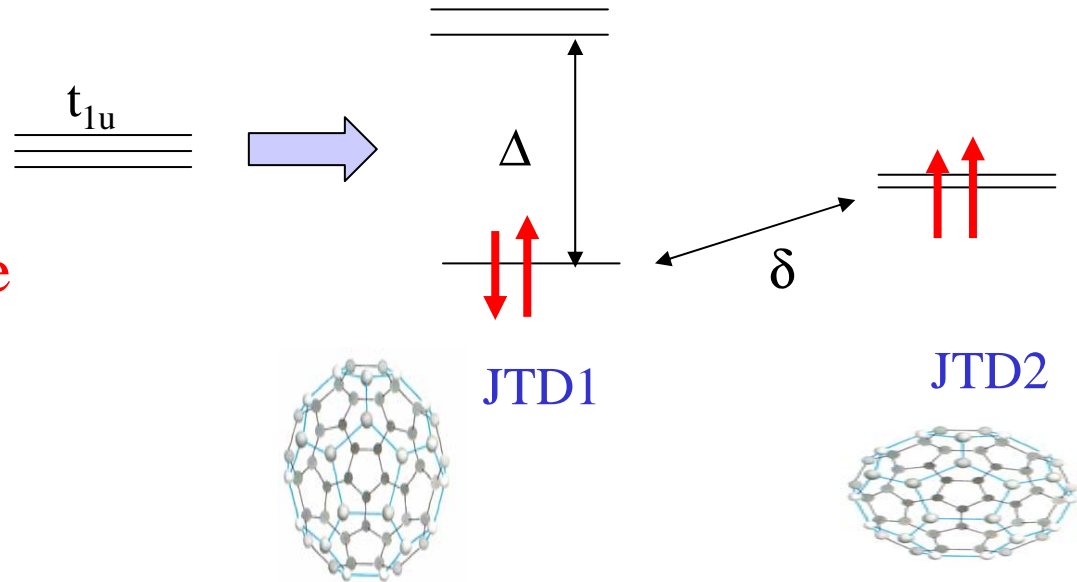
n
M. Héritier, W. Victoroff
O. Gunnarson
M. Fabrizio, E. Tosatti

Insulating states of A_4C_{60} and Na_2C_{60}

(Fabrizio, Tosatti PRB97)

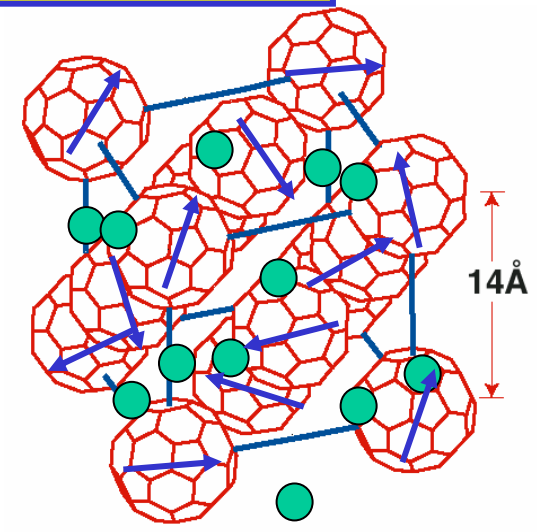
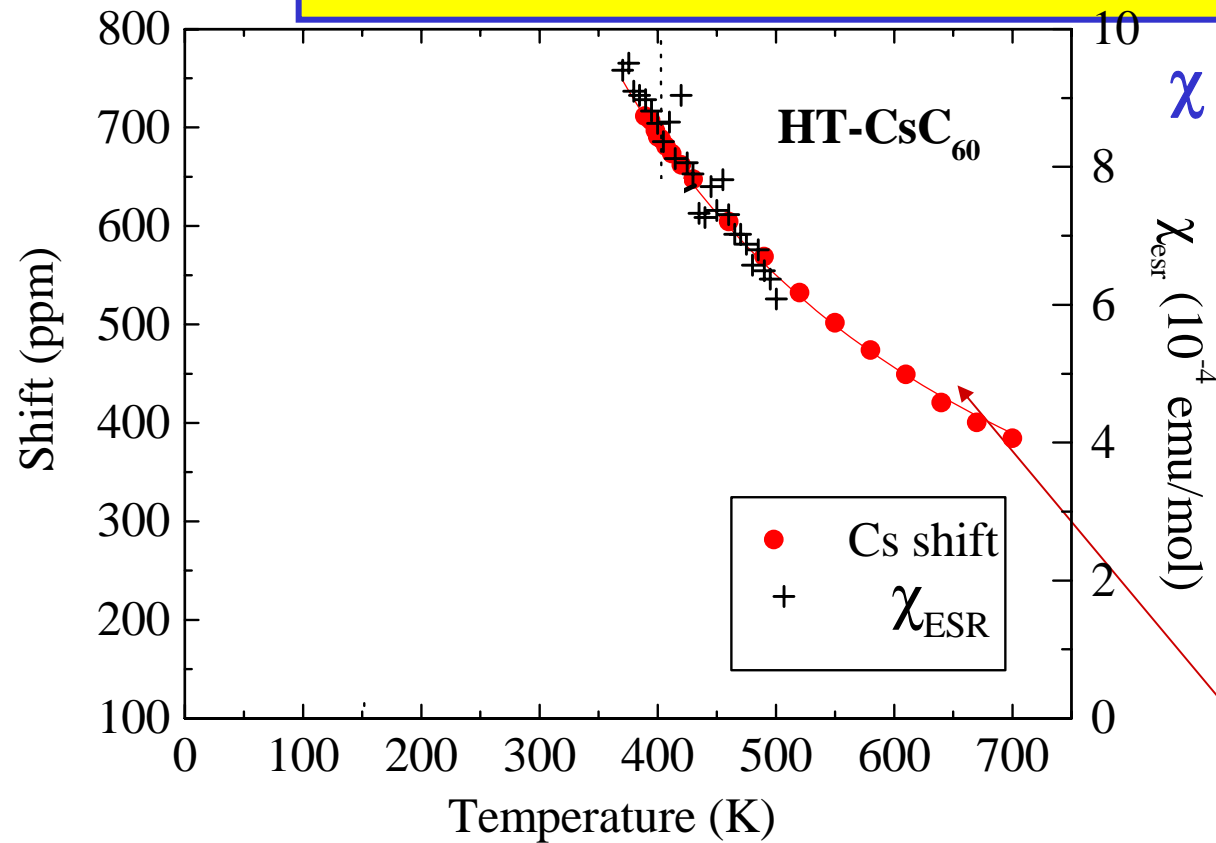
Experimental situation

- Non magnetic ground state
- 2 different gaps



- **Hund's rule is not obeyed** in the fundamental JTD1 state
- **Excited state is triplet** and corresponds to a different JTD2
- NMR detects the **Singlet Triplet excitation δ**

The special case of the cubic phase of CsC₆₀



$$\chi_P = \frac{g^2 \mu_B S(S+1)}{3k_B(T+\theta)}$$

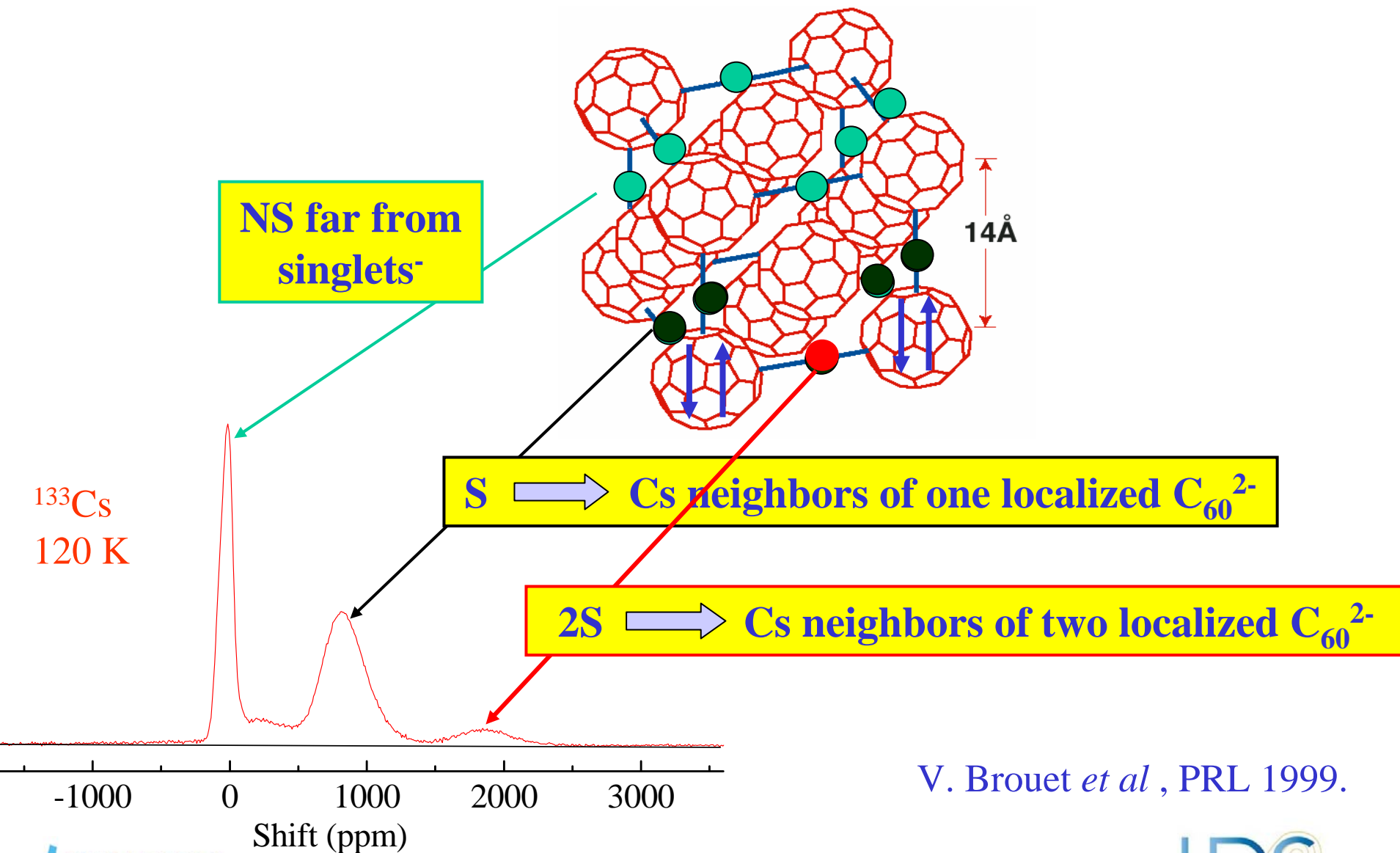
with $S=1/2$ and $\theta < 5\text{K}$

**Strong correlations
for A₃C₆₀ as well ?**

Both NMR and ESR indicate a
paramagnetic Curie law for fcc AC₆₀

So CsC₆₀ is a Mott insulator

Evidence for charge segregation: singlet states



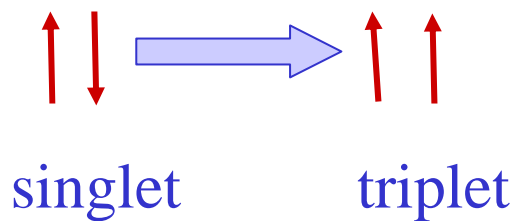
V. Brouet *et al*, PRL 1999.

Comparison of Cs NMR in QC- CsC_{60} and HT- CsC_{60}

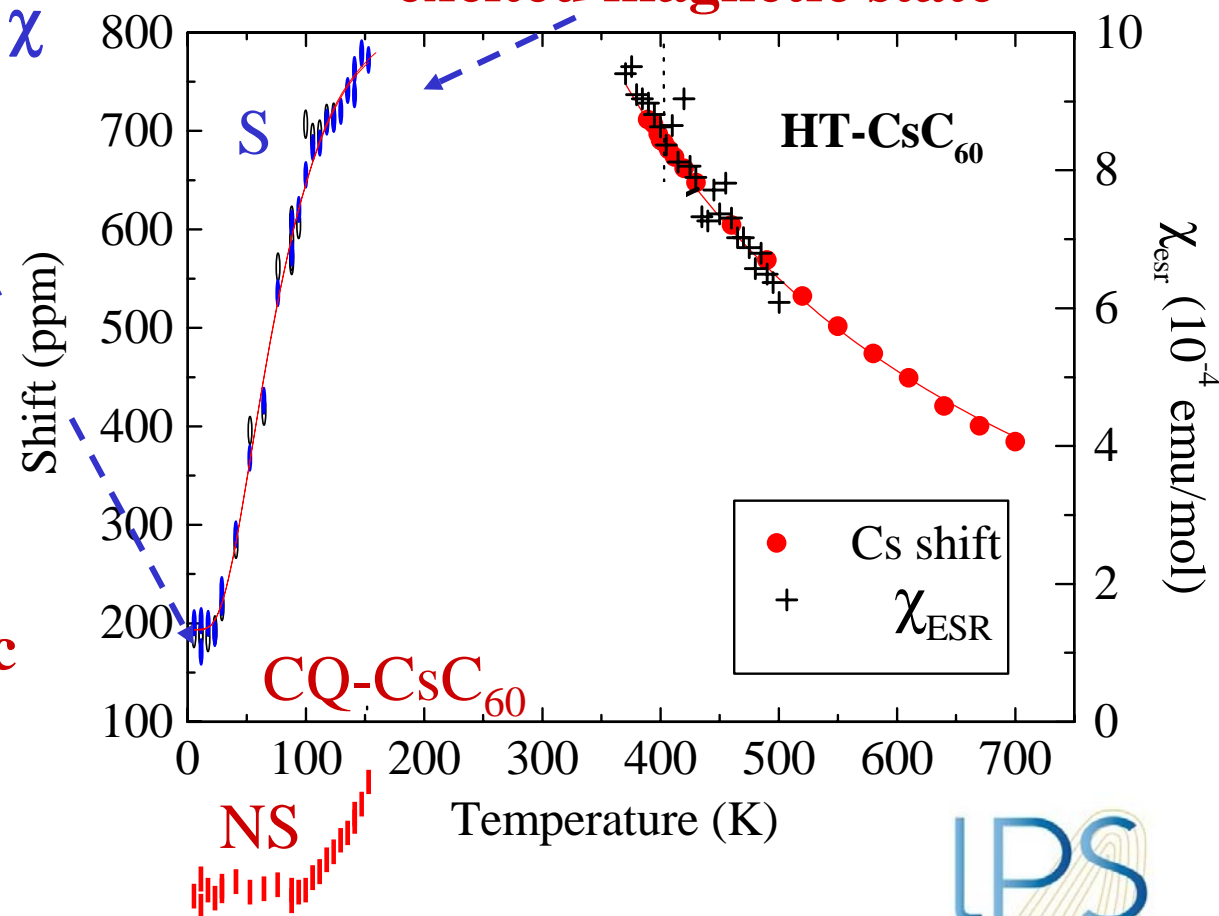
Very different from HT cubic phase

S (and 2S) sense a transition from a **local non-magnetic ground state**

..... towards an **excited magnetic state**



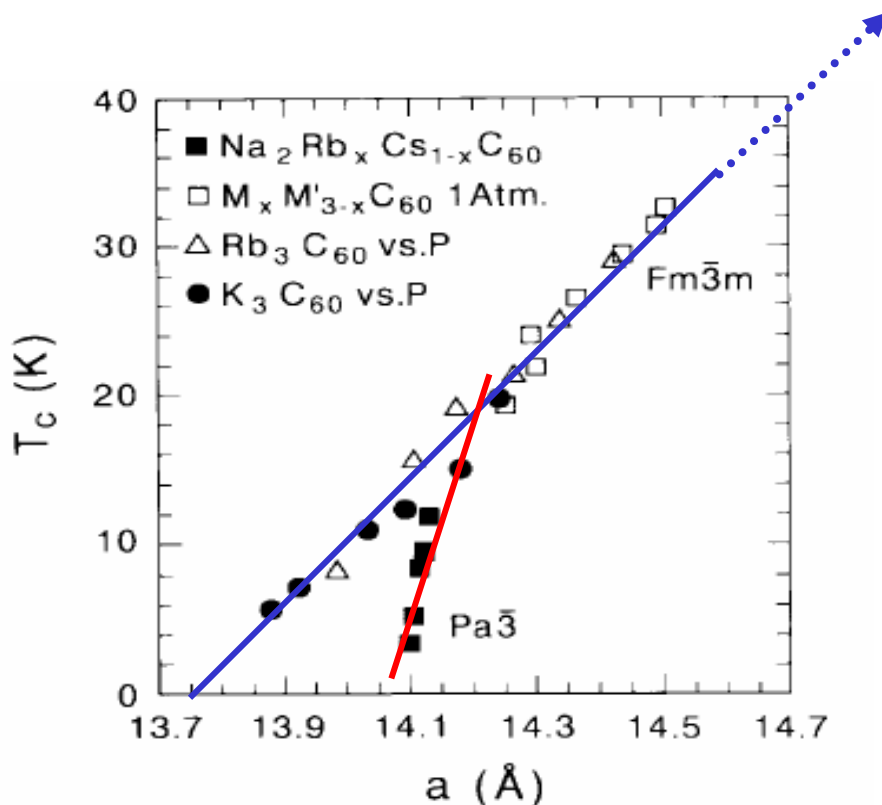
NS has a non magnetic **T** dependence



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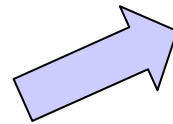
Superconductivity in A_3C_{60}



Higher T_c ?

Mott transition?

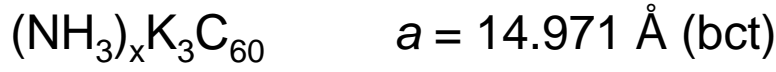
Alkali Ionic radius



- Li
- Na
- K
- Rb
- Cs
- Fr



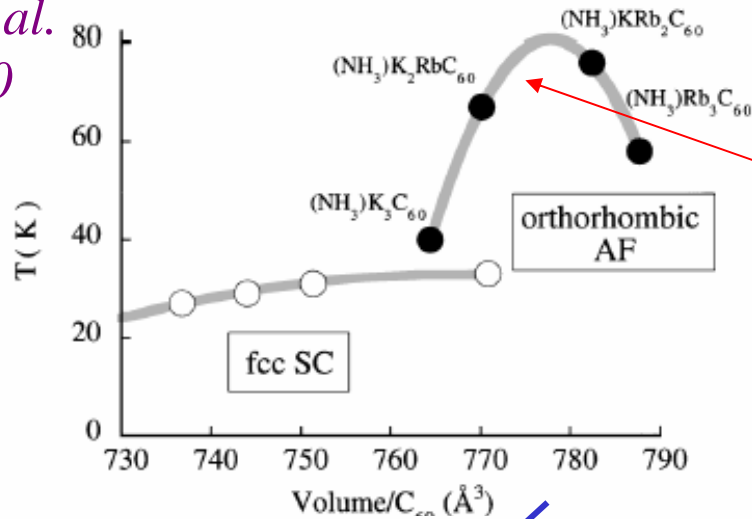
Expanded Mott state in K_3C_{60} expanded by NH_3 molecules



Rosseinsky Nature **364** 425 (1993).

Metal-insulator transition in expanded A_3C_{60}

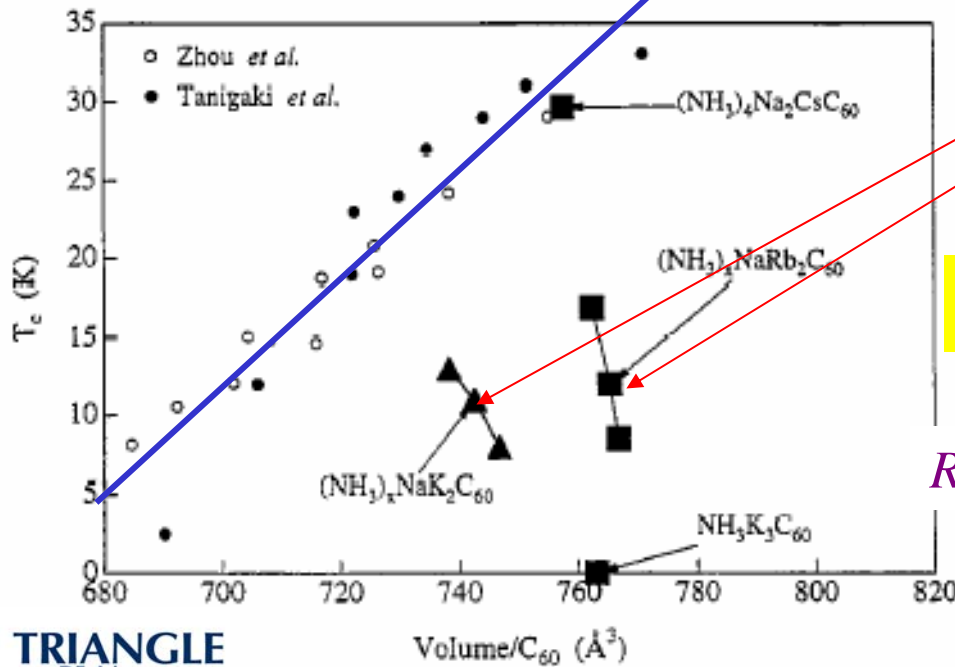
Iwasa et al.
PRL2000



**For large lattice spacings:
the system becomes AF.**

But:

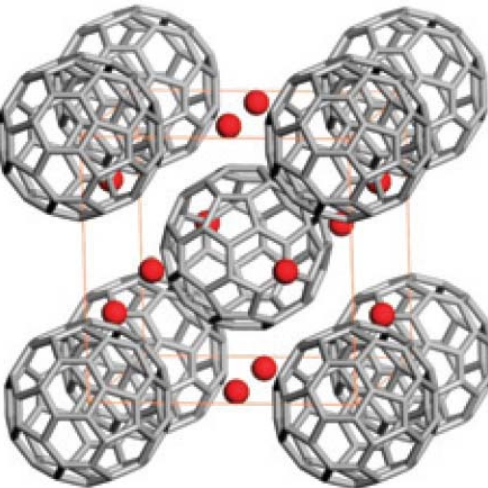
**structural distortions also
induce magnetism**



Lowering of the degeneracy?

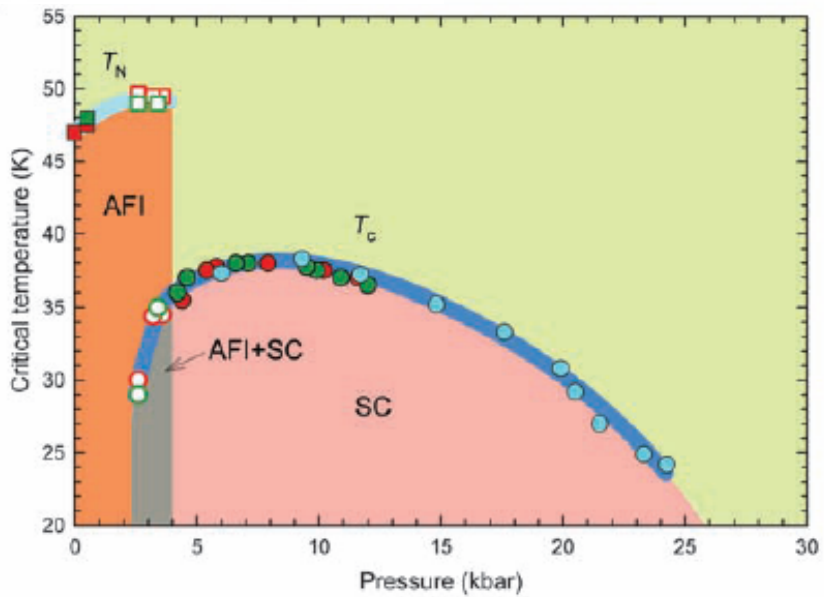
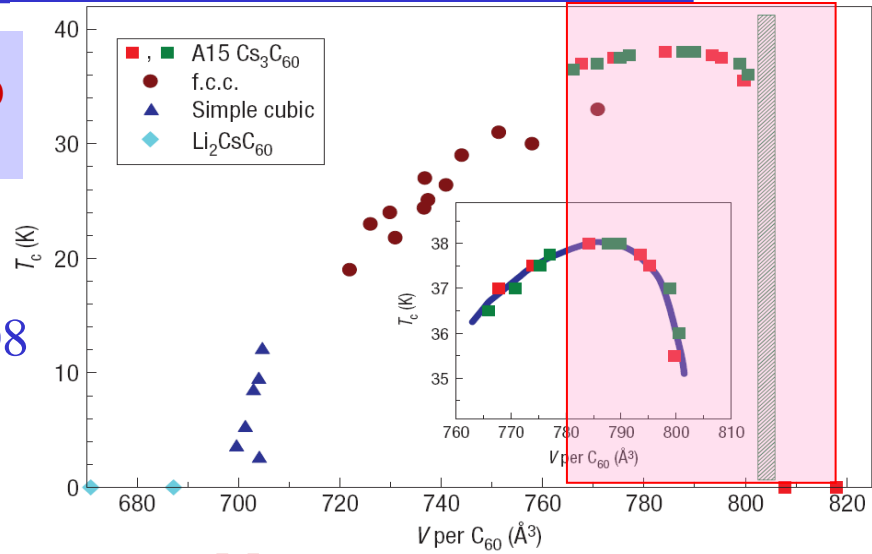
Rosseinsky, Maniwa, Iwasa, Prassides

A15 Cs_3C_{60} pressure induced superconductivity

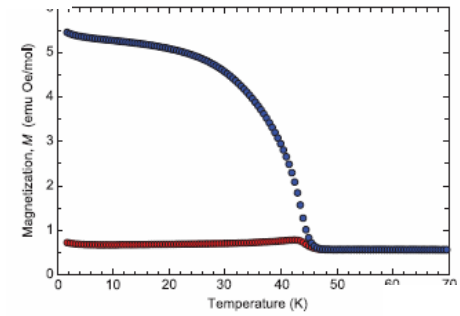


New phase of Cs_3C_{60} with A15 structure

A. Ganin et al
Nature Materials, 2008



M

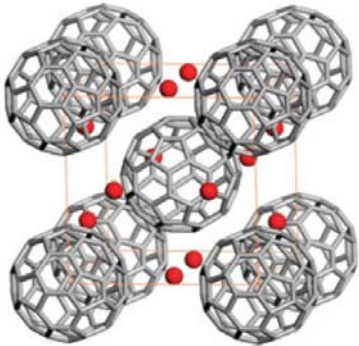


T

Magnetic at ambient pressure
SC and Mott transition

Takabayashi et al, Science 2009

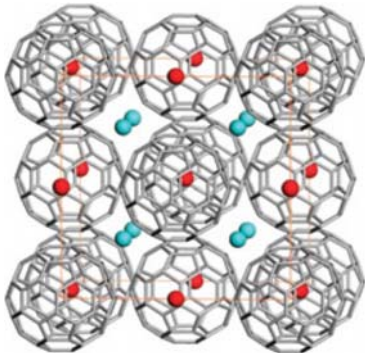
Multiple phases in the samples: ^{133}Cs NMR is very helpful



A15 Cs_3C_{60}

A single Cs site
with non cubic
local symmetry

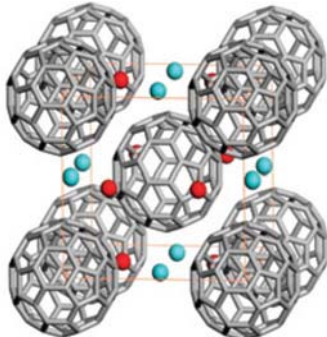
^{133}Cs $I=7/2$



fcc Cs_3C_{60}

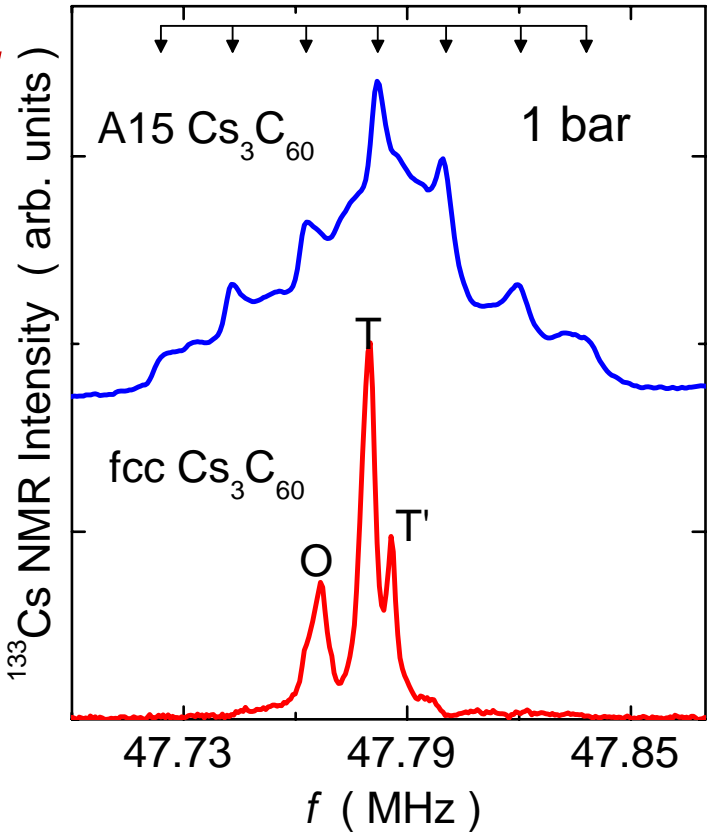
Well known: two Cs sites
O and T = 1 : 2
With cubic local symmetry
T splits at low T

Merohedral disorder of the C_{60}



Cs_4C_{60}

Insulating phase
Eliminated by its very long T_1



A15-rich sample

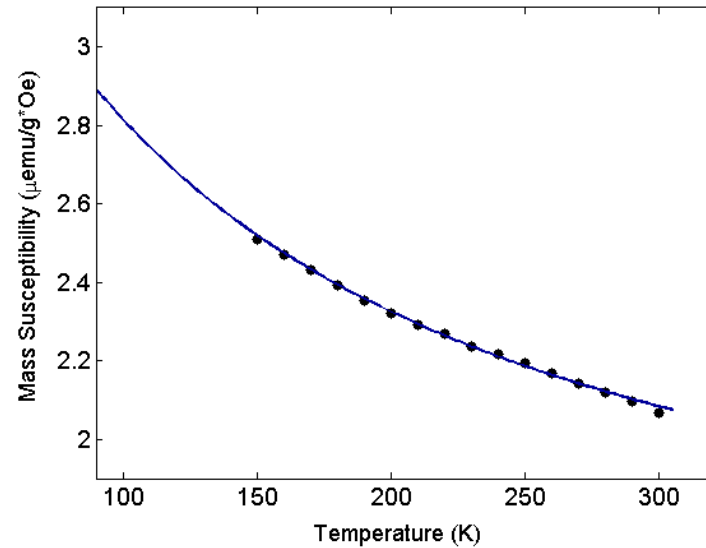
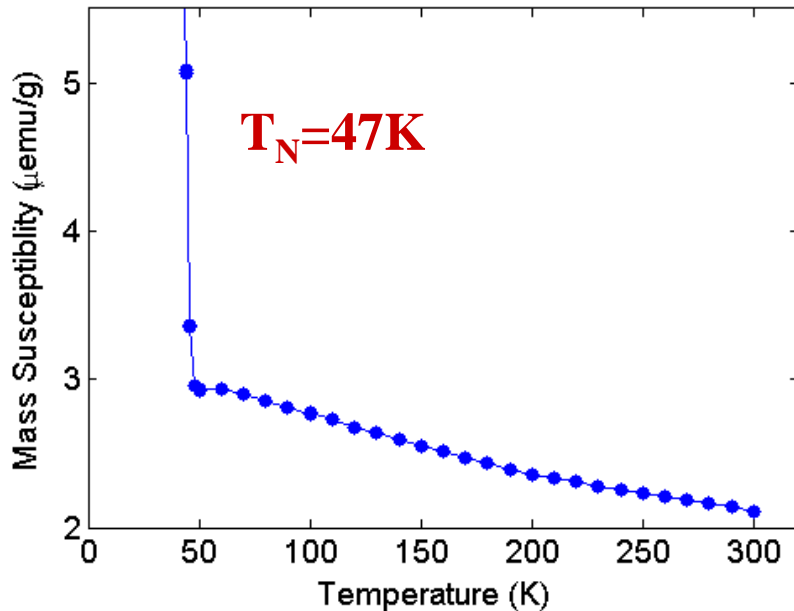
- A15 : **58.4 %**
- FCC : 12 %
- Cs_4C_{60} : 29.5 %

fcc-rich sample

- A15 : 34 %
- FCC : **55 %**
- Cs_4C_{60} : 11 %

Paramagnetic state susceptibility

(30.9% Cs_3C_{60} FCC, 53.6% Cs_3C_{60} A15, 15.5% Cs_4C_{60})



High T Curie-Weiss behavior

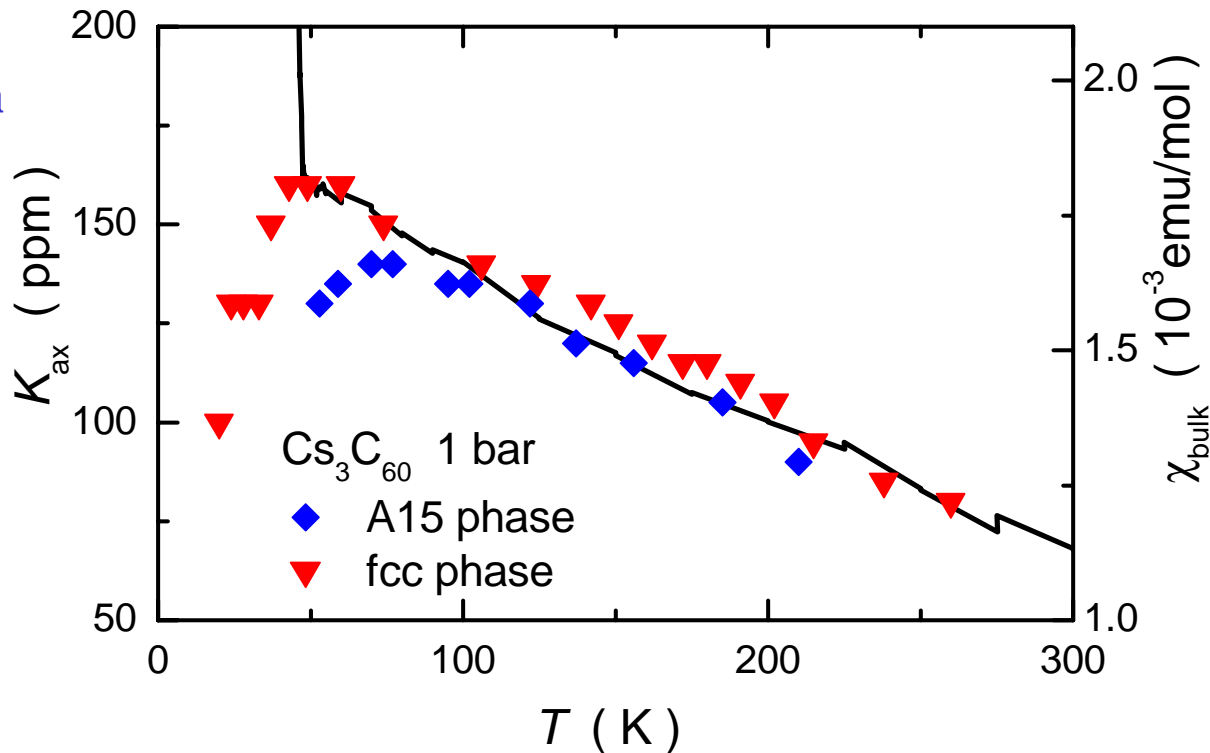
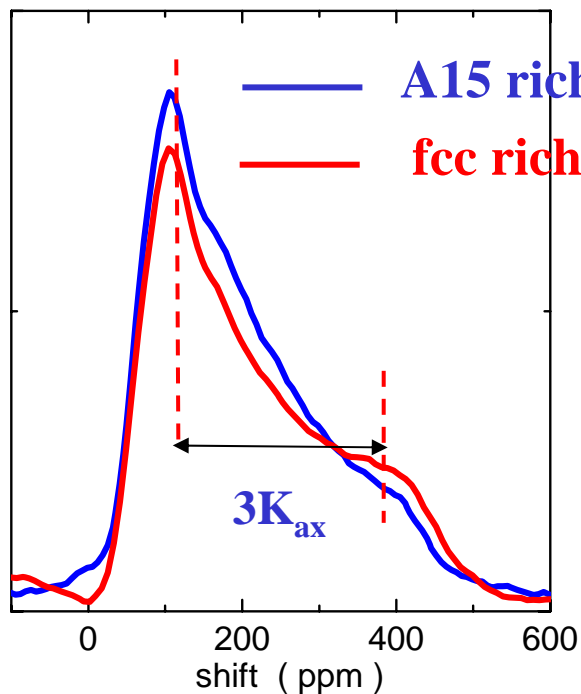
Weiss temperature: $\theta \sim 100\text{ K}$

Effective moment : $p_{\text{eff}} \sim 1.70 \mu_B$

$$\chi^{-1} = p_{\text{eff}}^2 / 3k_B (T + \theta)$$

Local moment $S \sim 1/2$ on the C_{60} balls?

Paramagnetic state susceptibility



Anisotropic hyperfine coupling

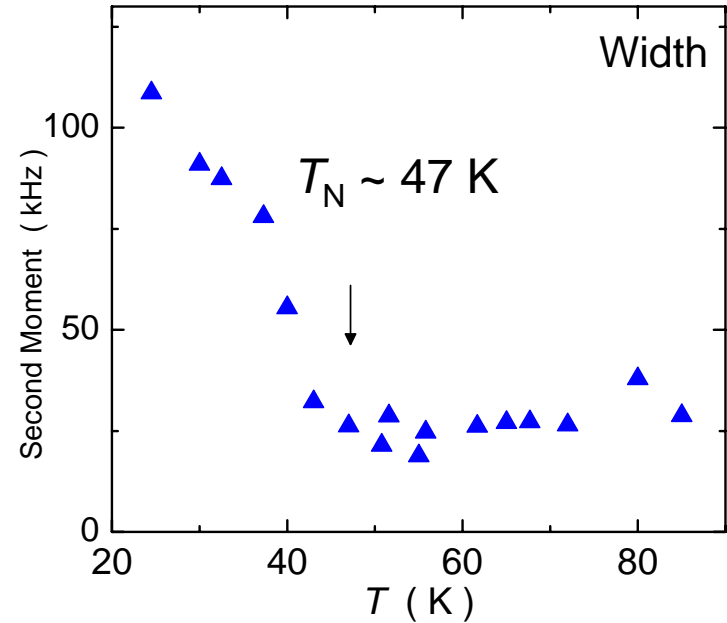
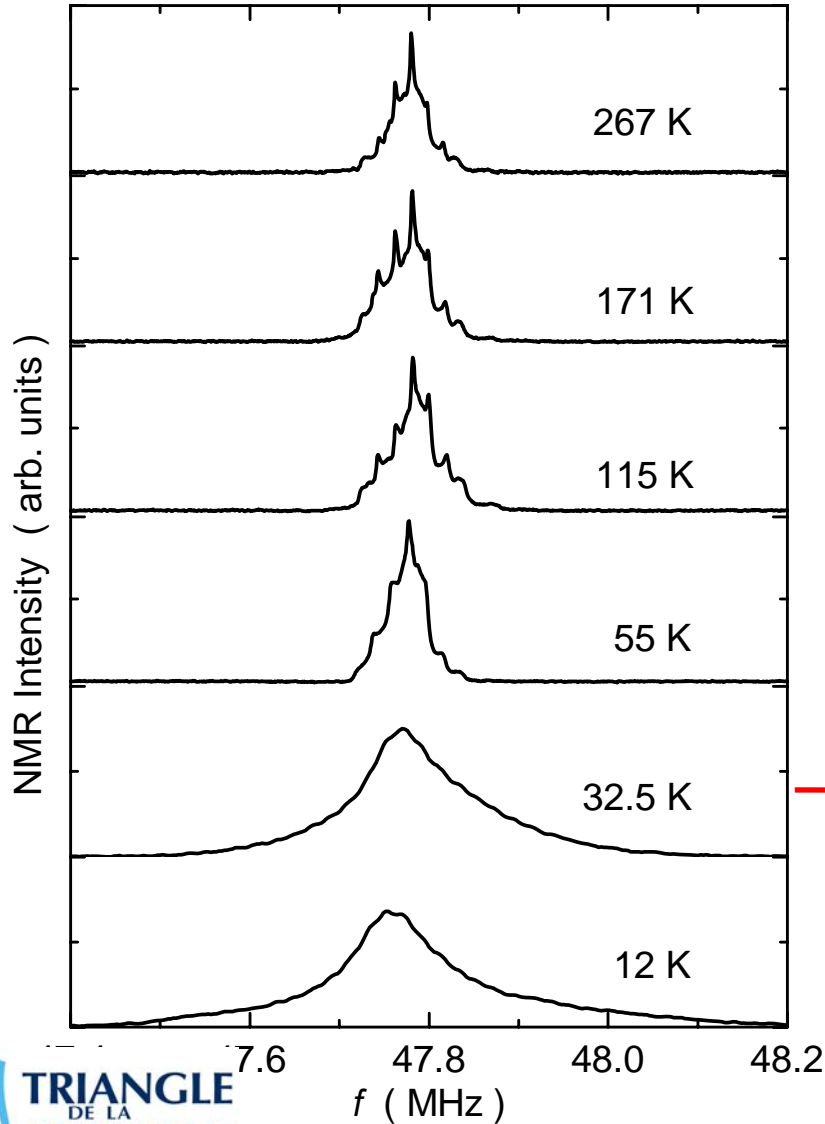
$$K_{ax} = A \chi_s$$

Here A is dipolar $A_{exp} \sim 700 \text{ Oe}/\mu_B$
 Calculated value : $A = 640 \text{ Oe}/\mu_B$

Local moment is indeed on the C₆₀ balls!

Magnetic transition in A15 phase

A15-rich sample



Spectral broadening :
static internal magnetic field.

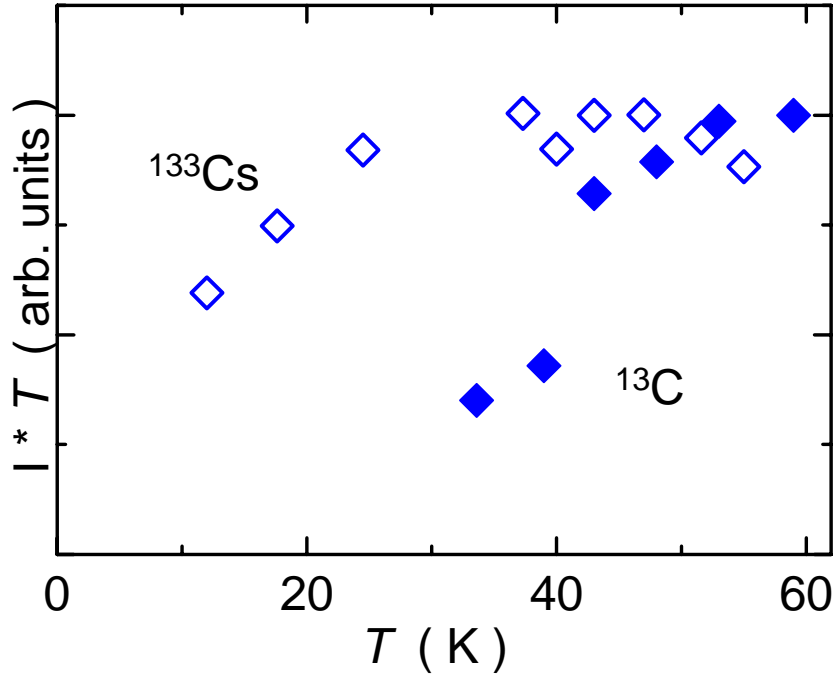
A15 phase shows magnetic order
below $T_N = 47$ K at ambient p .

Magnetism and crystal structure

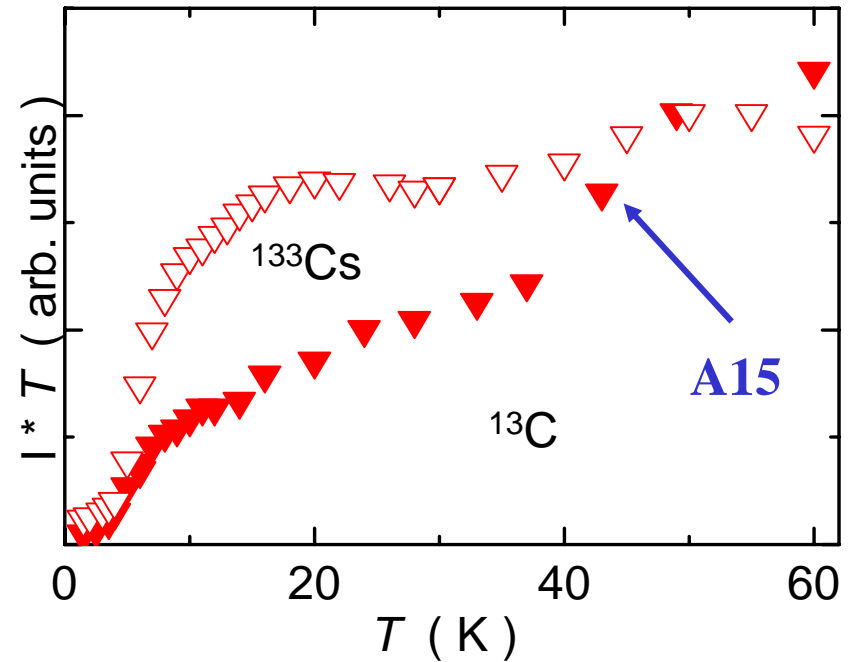
A15 rich

^{13}C and ^{133}Cs NMR intensities

fcc rich



$T_N = 47\text{K}$



Gradual freezing
About 10K ?

Geometrical frustration depresses the magnetic ordering
(fcc not bipartite)

Spin dynamics and crystal structure

^{133}Cs NMR T_1

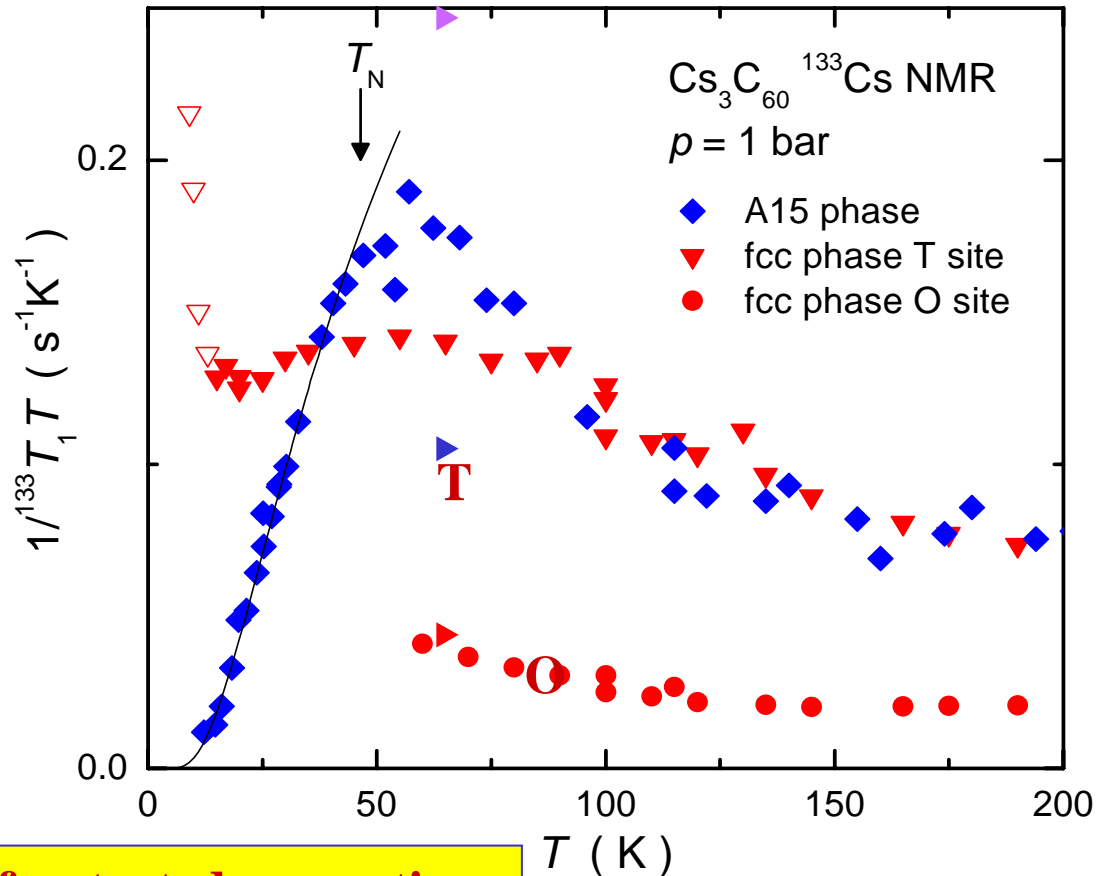
A15 phase
magnetic gap
 $\Delta \sim 50$ K

fcc phase
No magnetic gap.
Still some slow dynamics
below 10 K.

O site has twice smaller coupling
constant than that for T site.

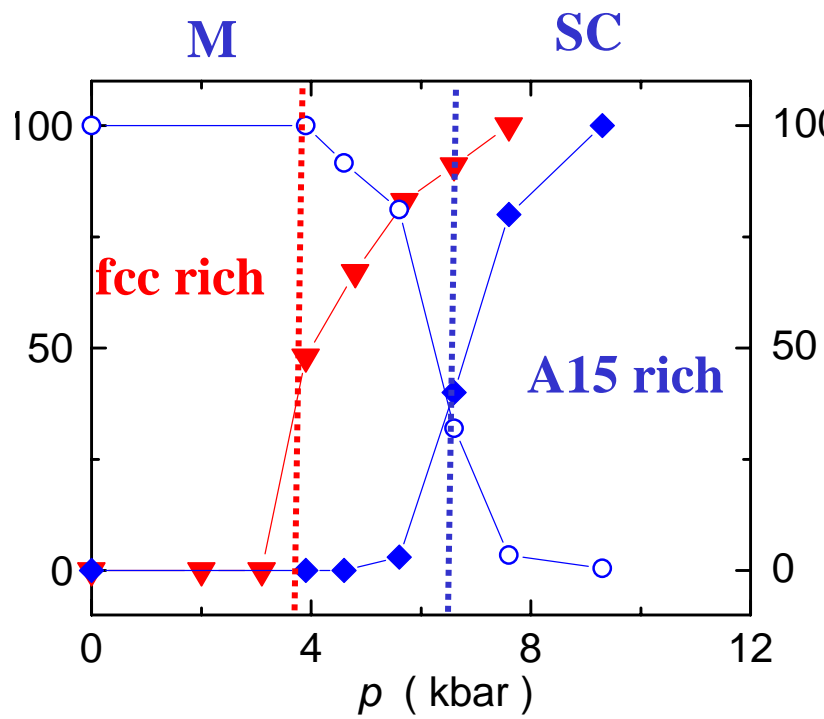
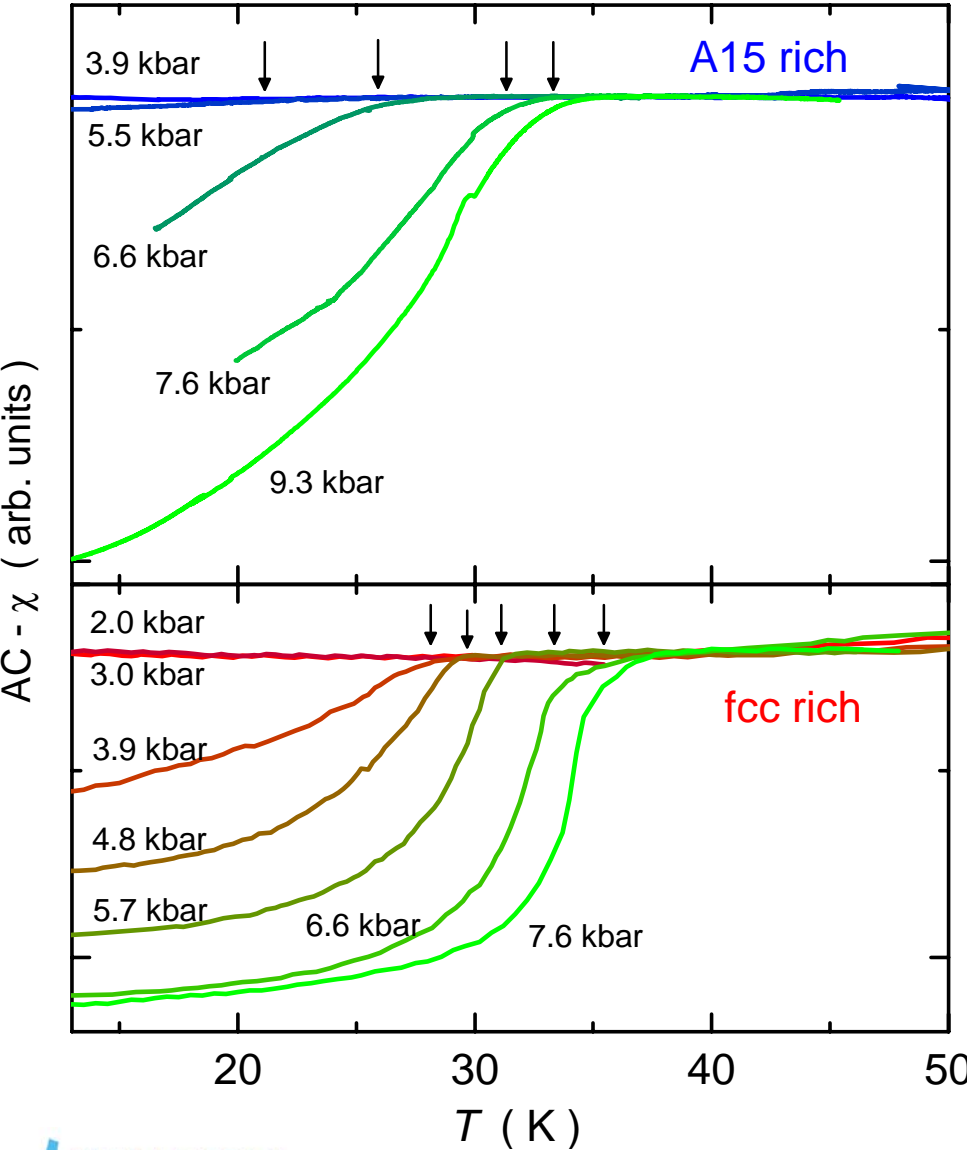
Both phases

Enhanced magnetic fluctuations
in the paramagnetic state.



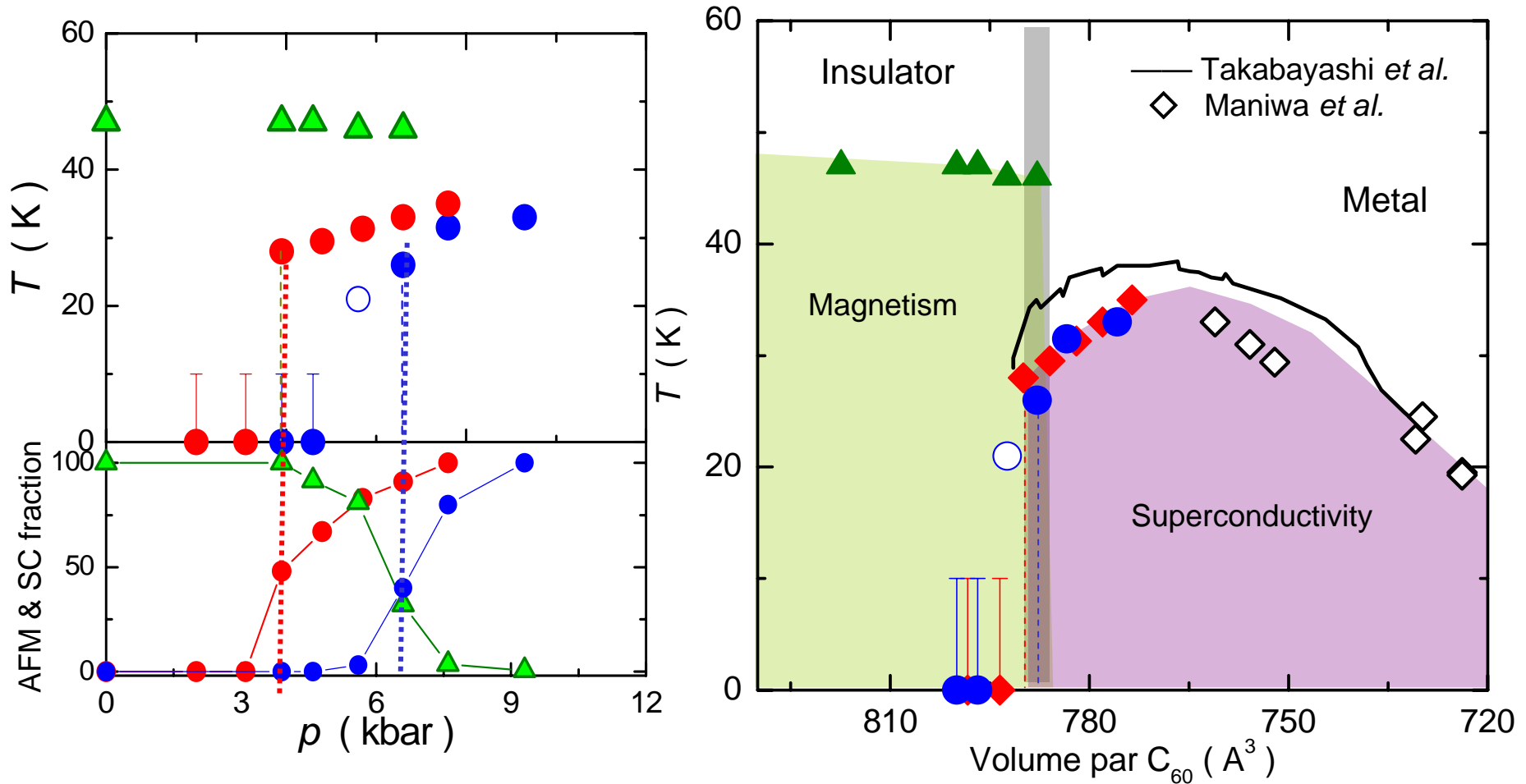
Ordered versus frustrated magnetism
(fcc not bipartite)?

Pressure induced superconductivity



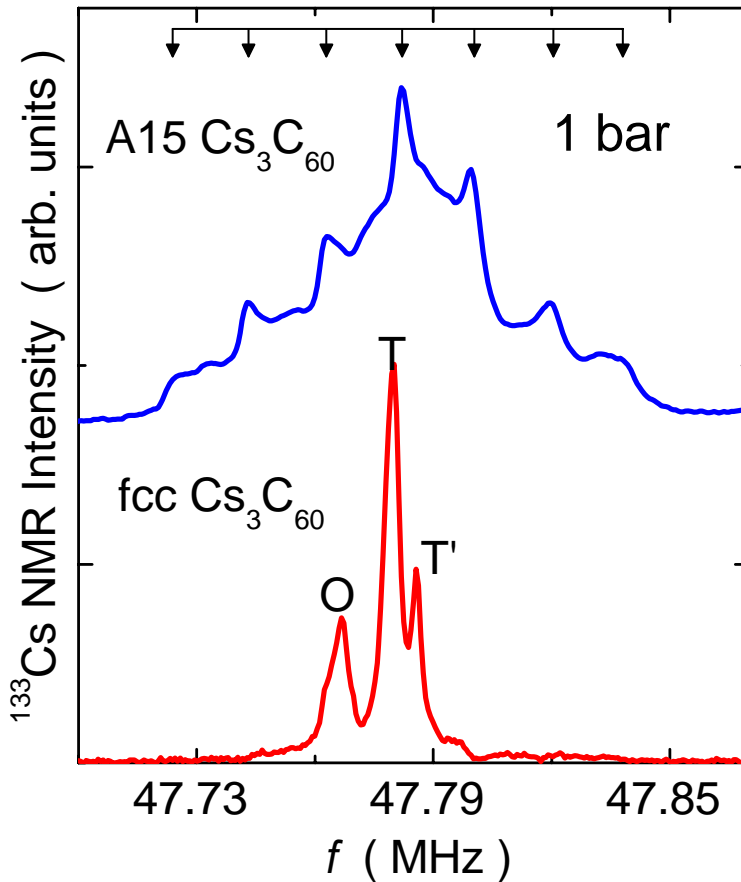
Distinct (p, T) phase diagrams for the two phases

Phase diagrams

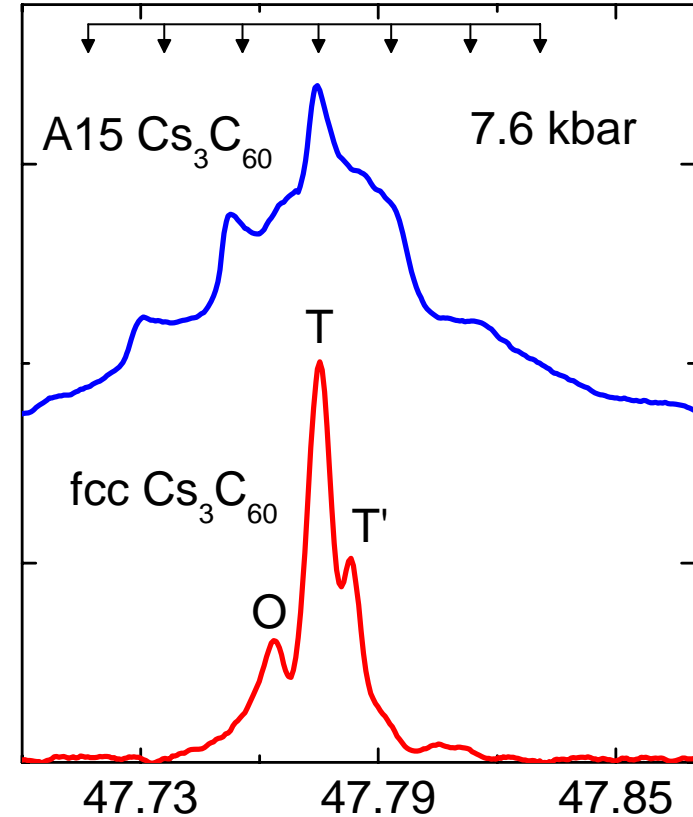


No difference with crystal structure on the SC side

Mott transitions to the metallic state in the A15 phase



$p=1\text{bar}$

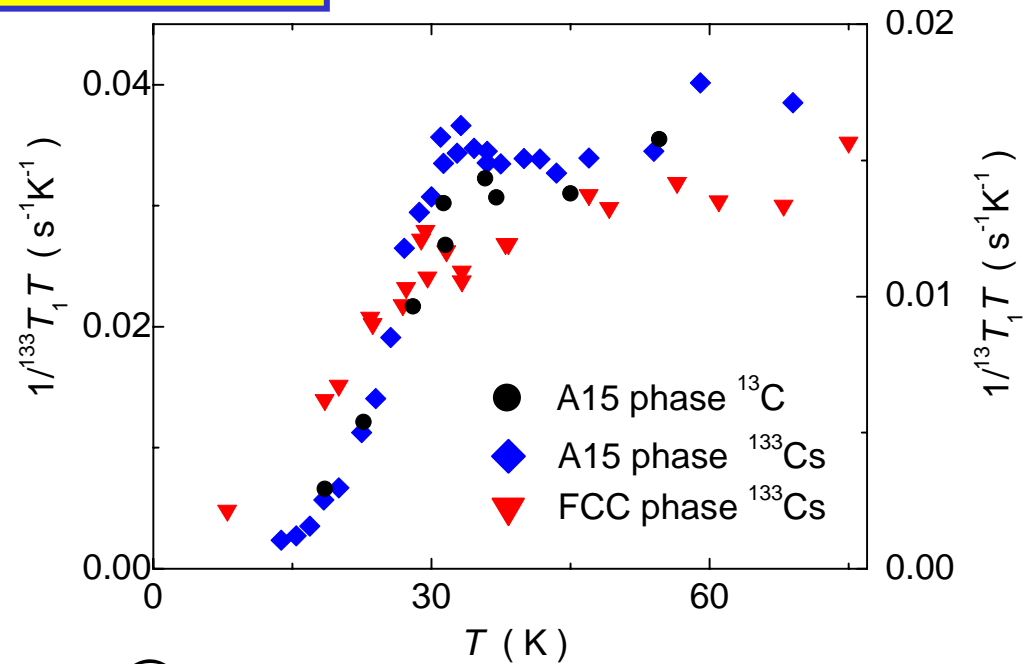
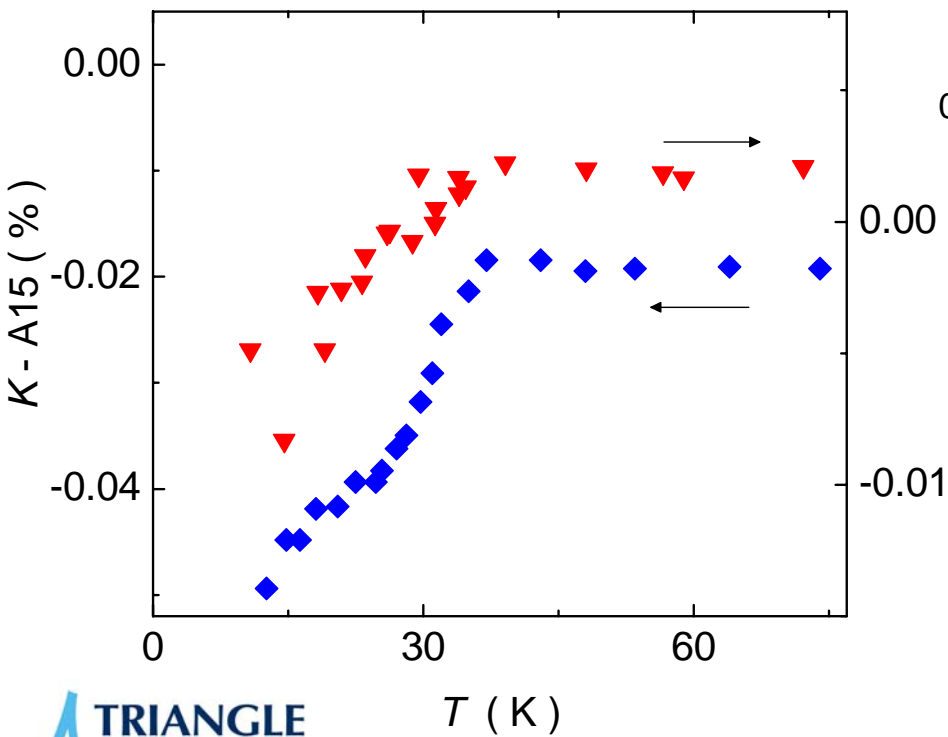


$p=7.6\text{ kbar}$

No change of crystal structures

Superconductivity

$p \gg p_c$



T_1 data

Knight shift
(singlet pairing)

Summary

- **Fulleride compounds are peculiar correlated electron systems**
- **Originalities** associated with their nanostructure
 - Icosaedric symmetry of the soccer ball.
 - Orientalional disorder
 - Internal degrees of freedom of the molecule:
 - Phonons, Molecular Jahn-Teller distortions**
- **They display many effects driven by correlations**
 - Static charge segregation in Cs_1C_{60}
 - High T_c superconductivity near a MIT (Cs_3C_{60})**
 - Molecular excitations survive in the solid.
- **Those are strongly influenced by molecular Jahn Teller effects**
 - Favor singlet formation in Cs_1C_{60}
 - Jahn Teller Mott insulating states in K_4C_{60} or Na_2C_{60}
 - Jahn-Teller effects explain why electronic correlations appear smaller in A_3C_{60} but remain sizable as Cs_3C_{60} is magnetic
- **Excellent possibility to study multiorbital Mott transitions**

