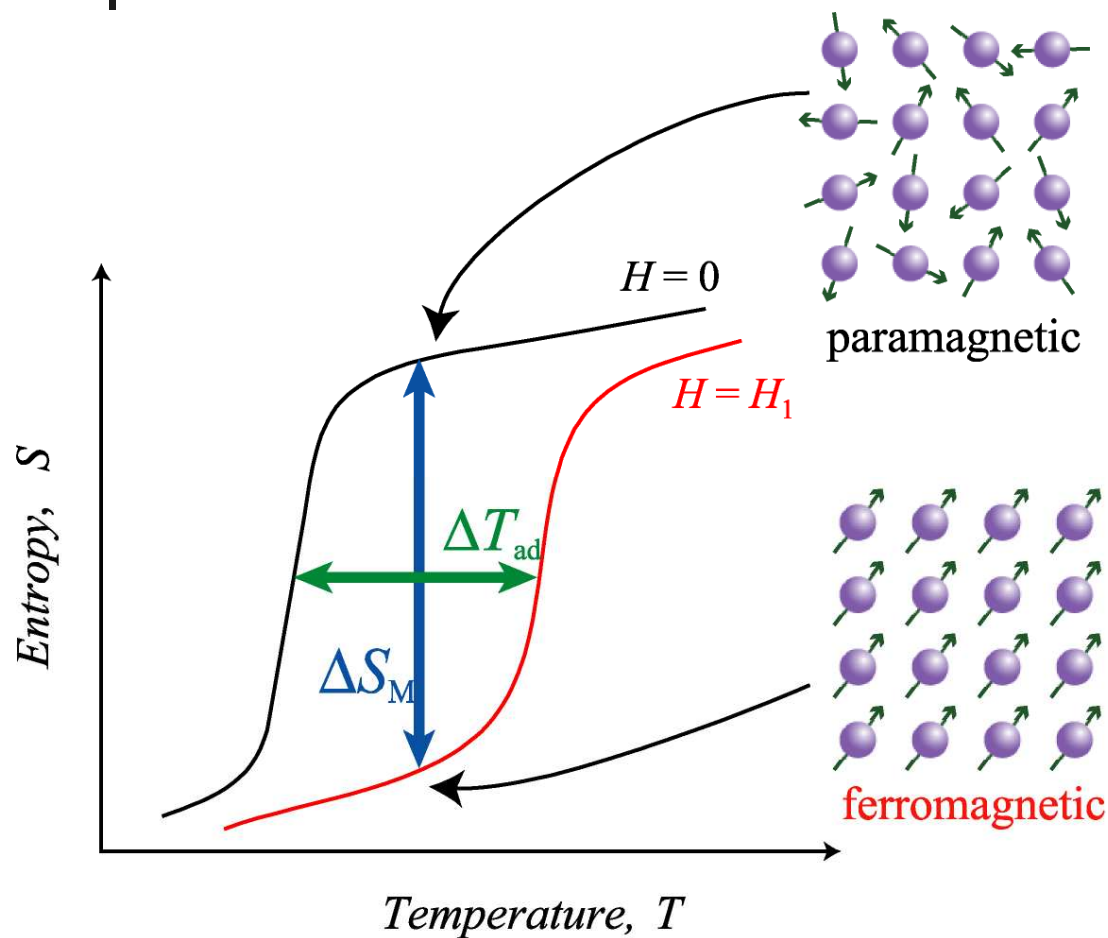


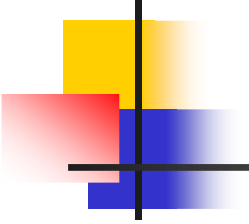
**Pressure dependence of
magnetocaloric effect (MCE) of
 $\text{MnAs}_{1-x}\text{Sb}_x$ and $\text{La}(\text{Fe}_x\text{Si}_{1-x})_{13}$**

H. Wada

*Department of Physics,
Kyushu University, Fukuoka, Japan*

MCE of a first-order magnetic transition (FOMT) system





Advantages and disadvantages of FOMT systems

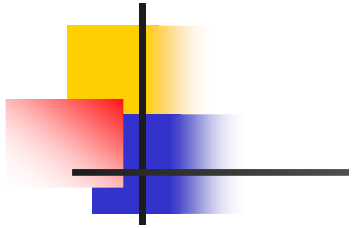
Advantages

- Giant or large MCE in a weak magnetic field.
- Most of materials are transition metal based compounds.

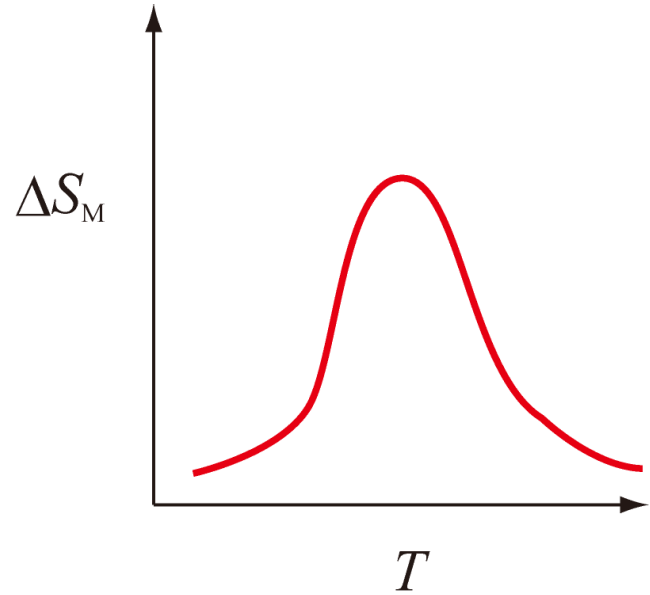
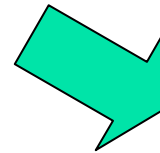
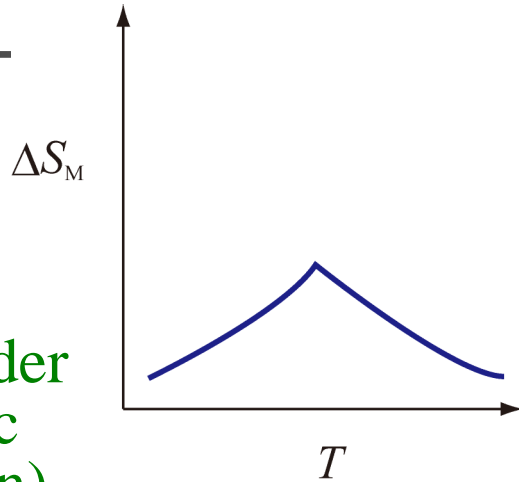
Disadvantages

- MCE is limited in a narrow temperature range.
- Additional properties are not favorable.

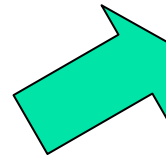
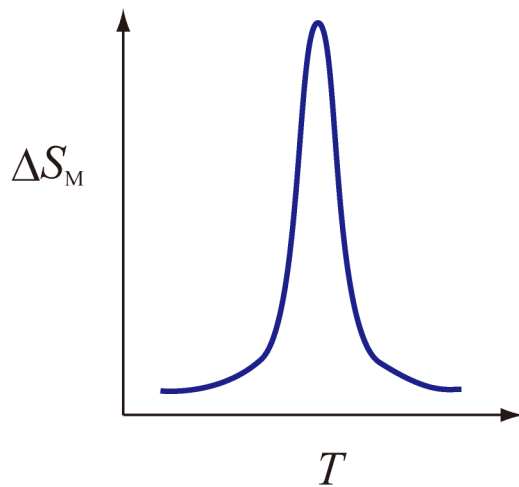
Our new approach



SOMT
(second-order
magnetic
transition)

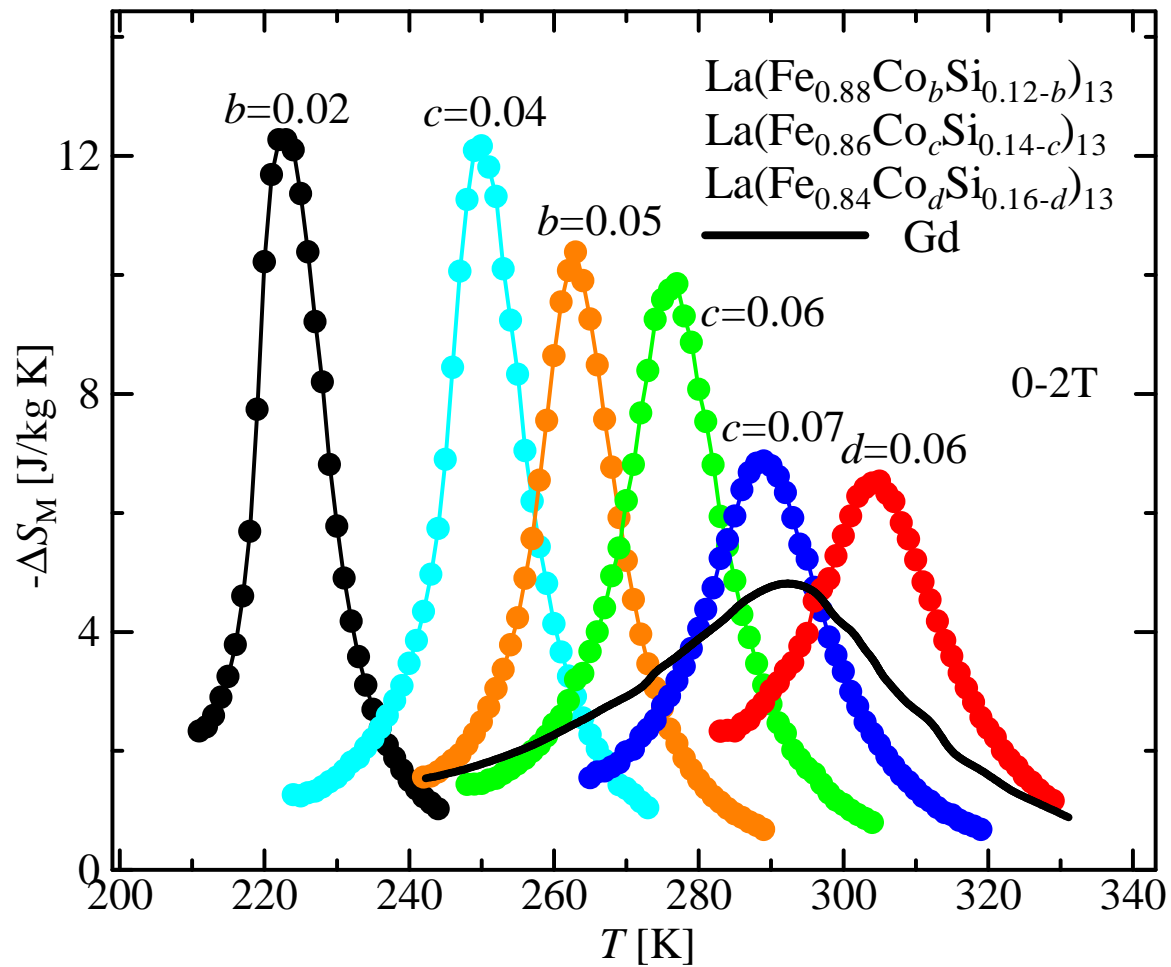


FOMT

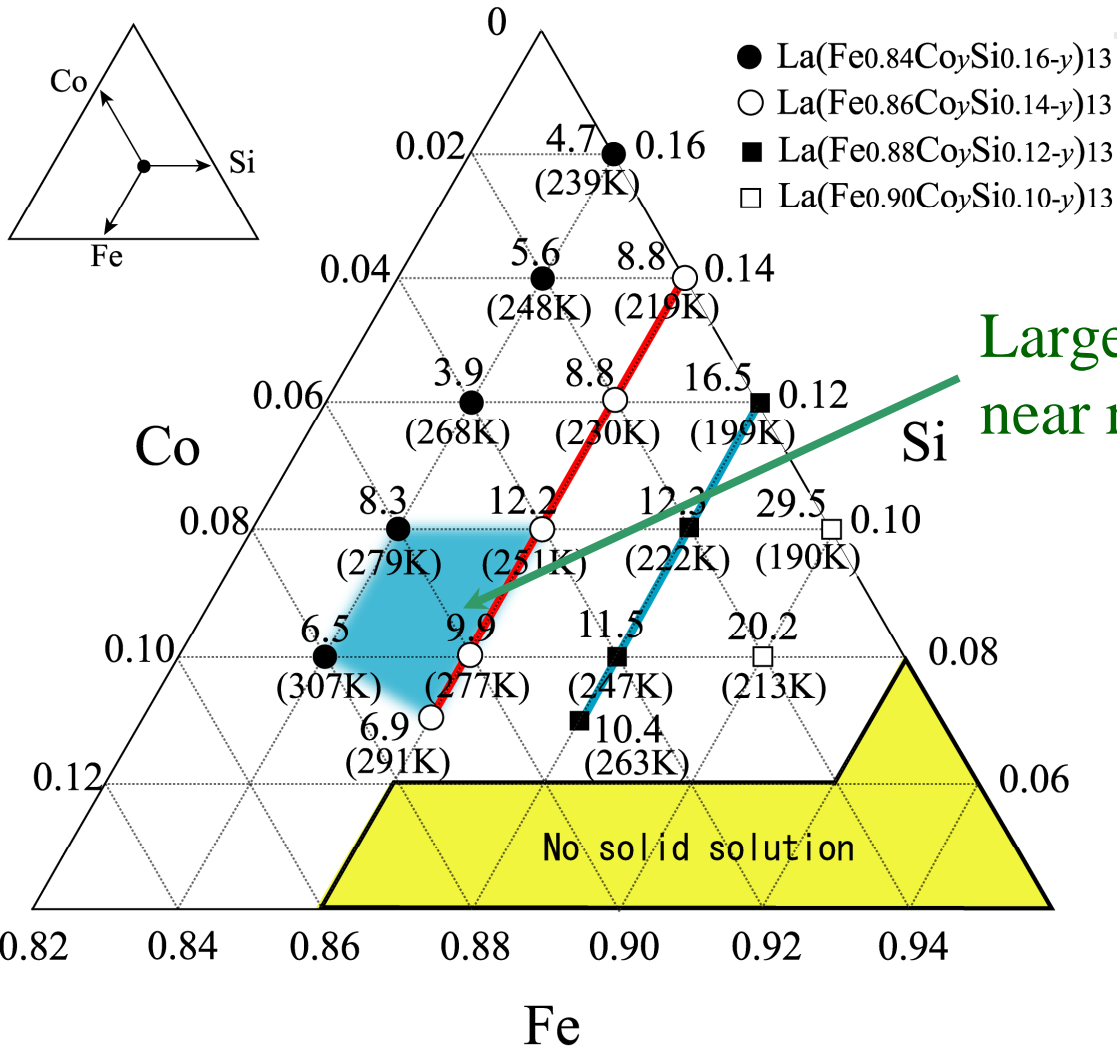
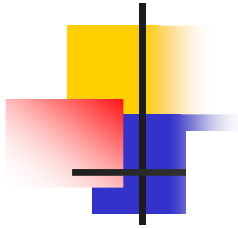


**Broadening
of a FOMT
(continuous)**

Example: $\text{La}(\text{Fe}_x\text{Co}_y\text{Si}_{1-x-y})_{13}$



Concentration dependence of ΔS_{Max} in 0-2T (J/K kg)



Large MCE is obtained near room temperature.

ΔS_{Max} (J/K kg)
(T_C)

Pressure: an effective tool to study the MCE



- Nature of magnetic transition

broadening of a FOMT

- Enhancement of MCE

MnAs?

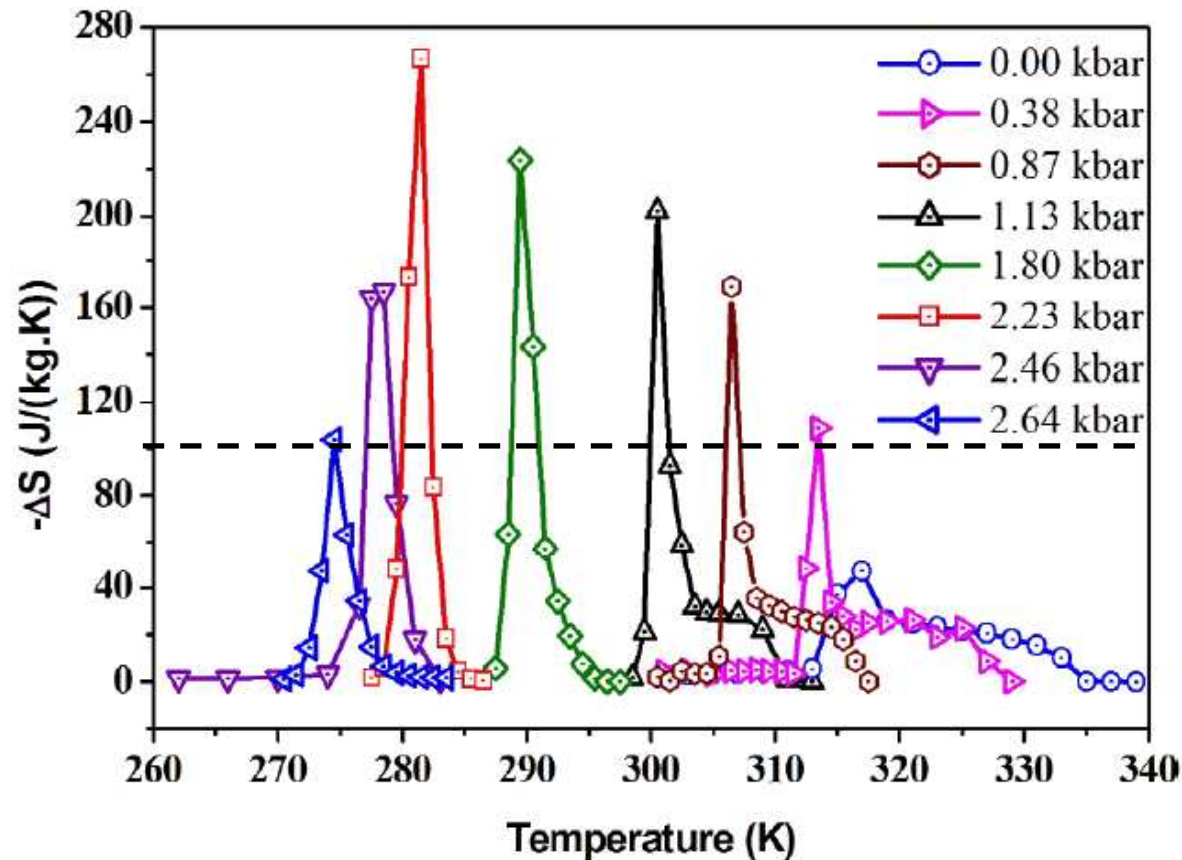
- Origin of MCE

magnetic entropy change

lattice entropy change

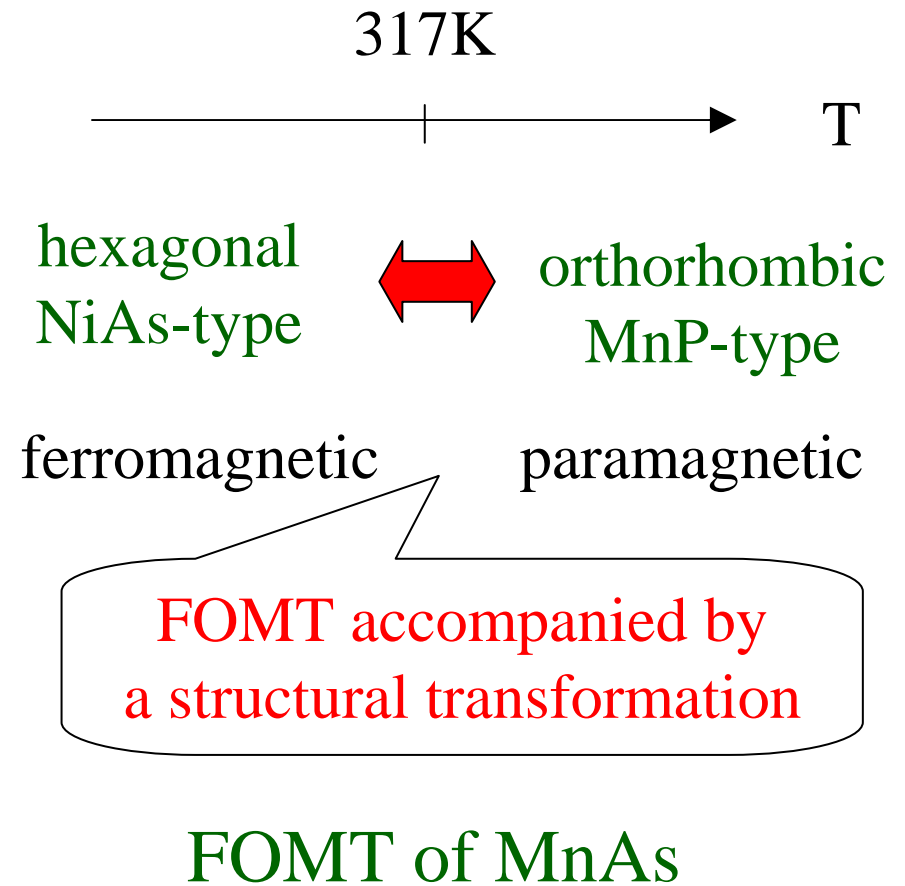
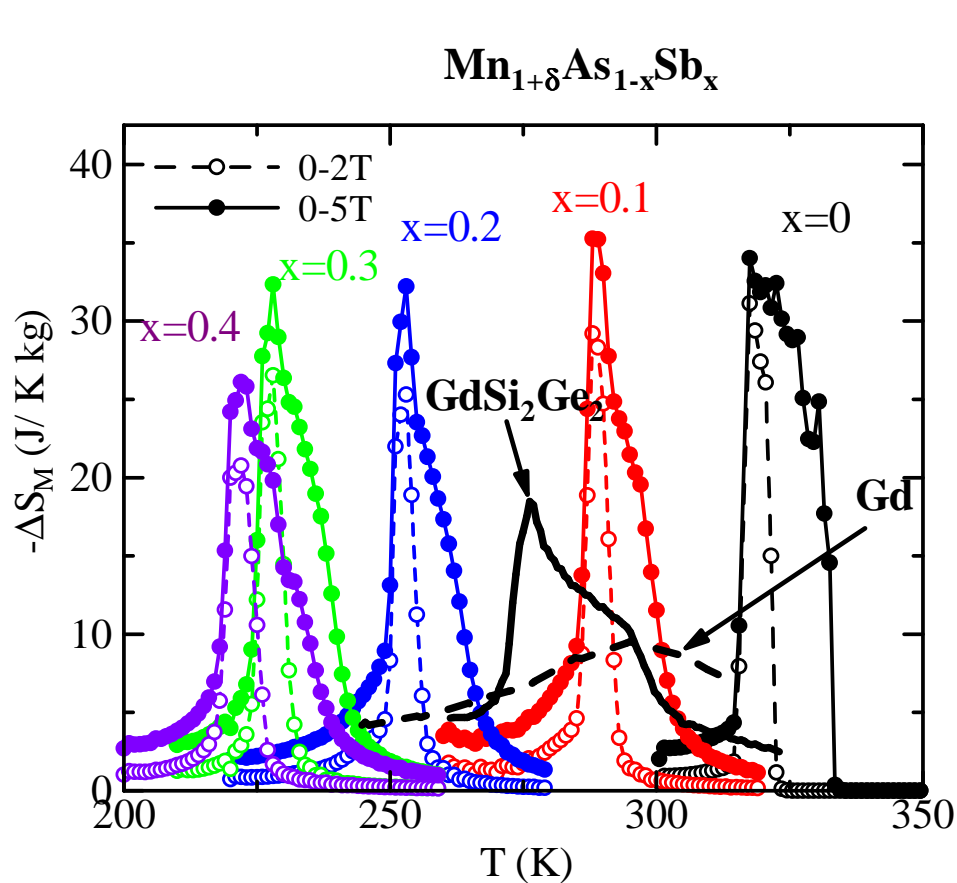
Colossal MCE of MnAs

Gama *et al.* 2004

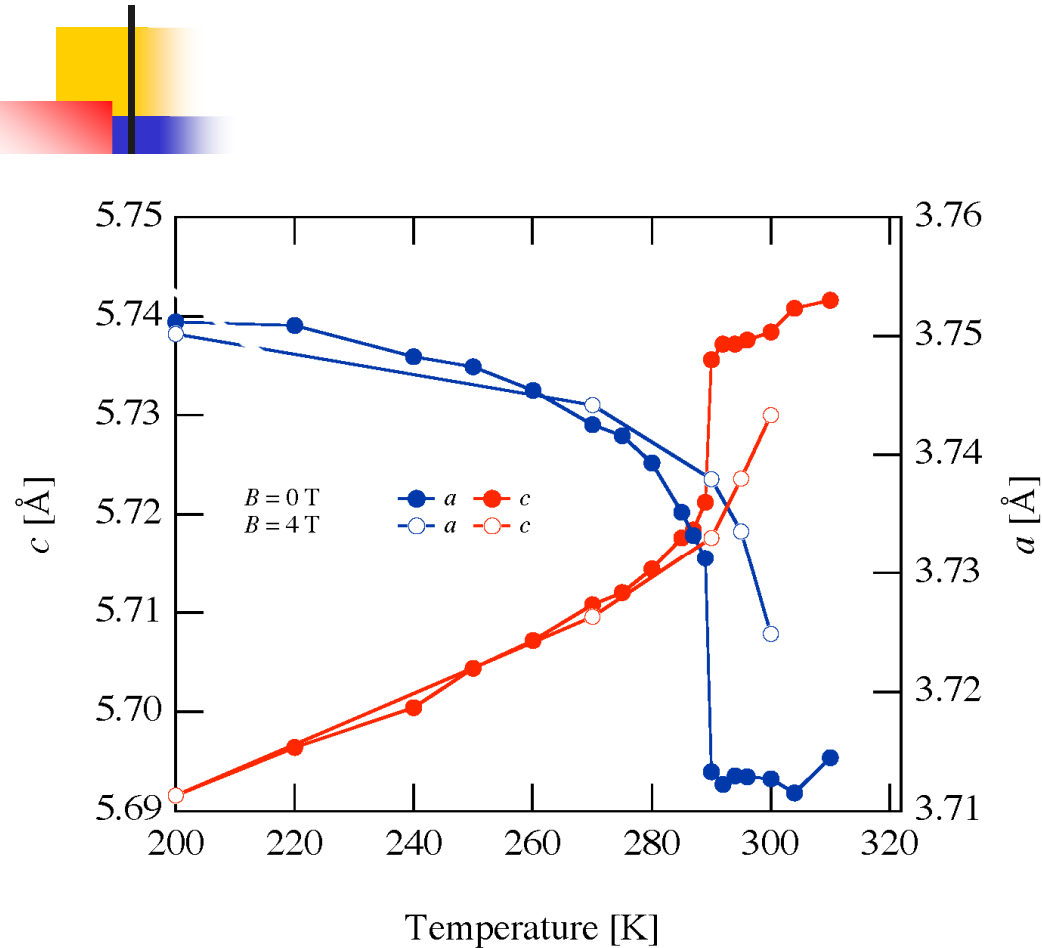


$$R \log(2S+1) = 103 \text{ J/kgK with } S=2.$$

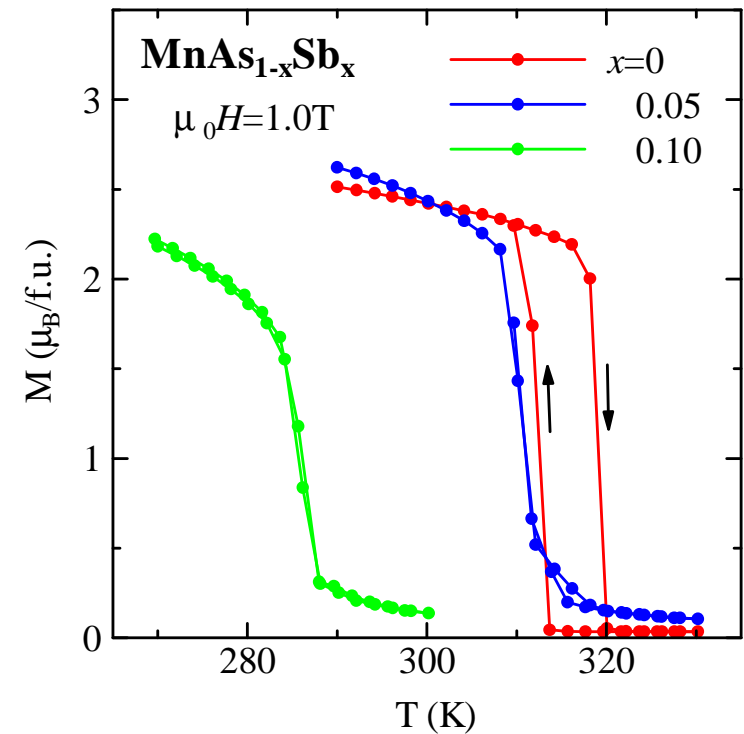
MCE of $\text{MnAs}_{1-x}\text{Sb}_x$ at ambient pressure



No structural transformation in $\text{MnAs}_{1-x}\text{Sb}_x$



No structural transformation
Sb stabilizes the NiAs-type structure.



Thermal hysteresis
is reduced

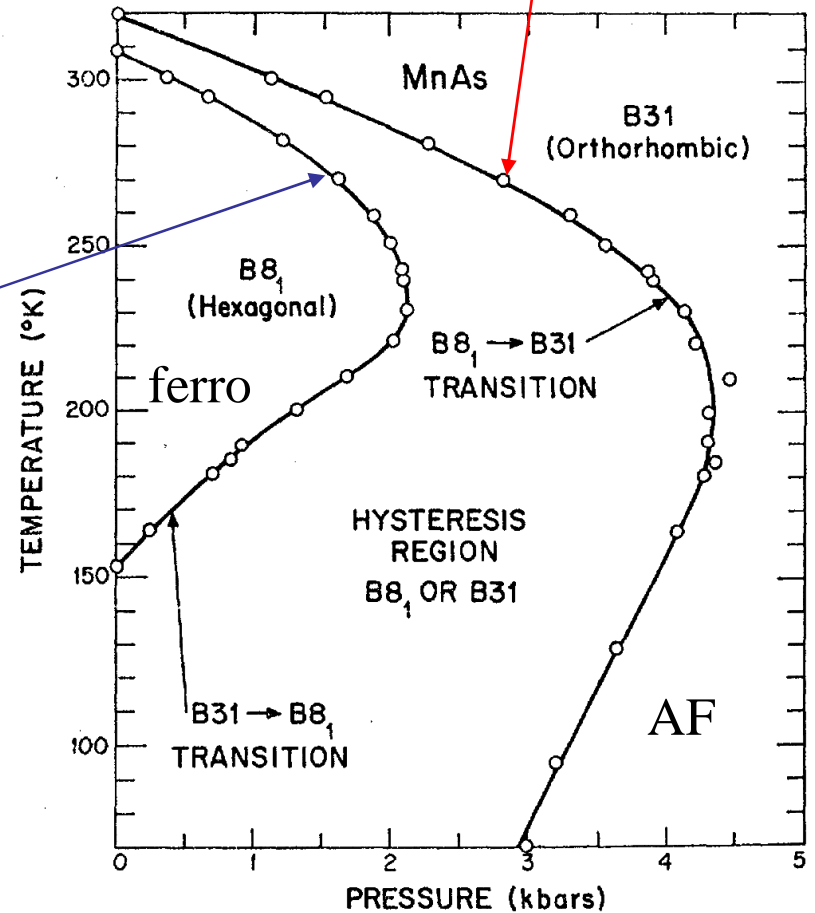
$T - p$ phase diagram of MnAs

Pressure stabilizes the MnP-type structure.

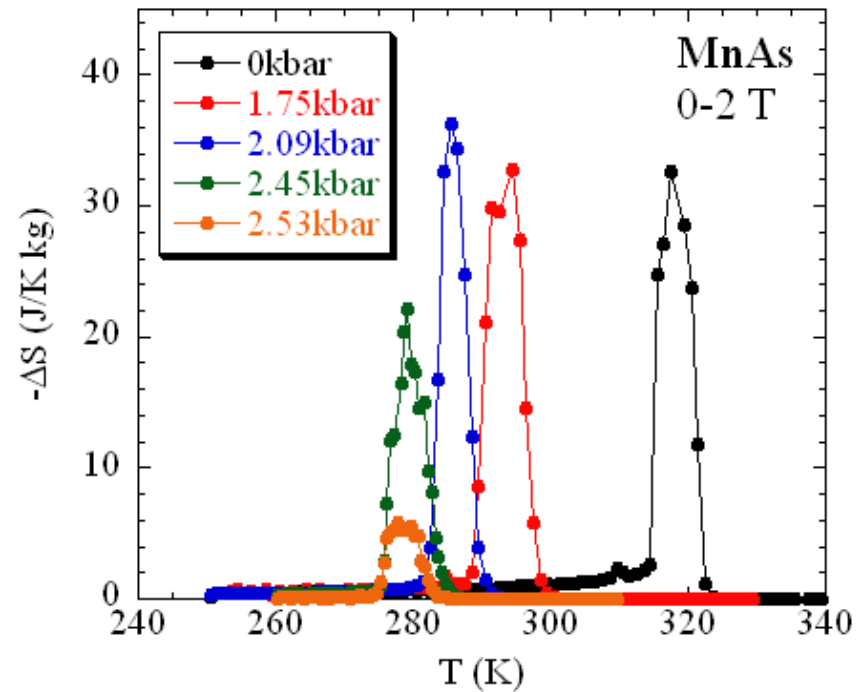
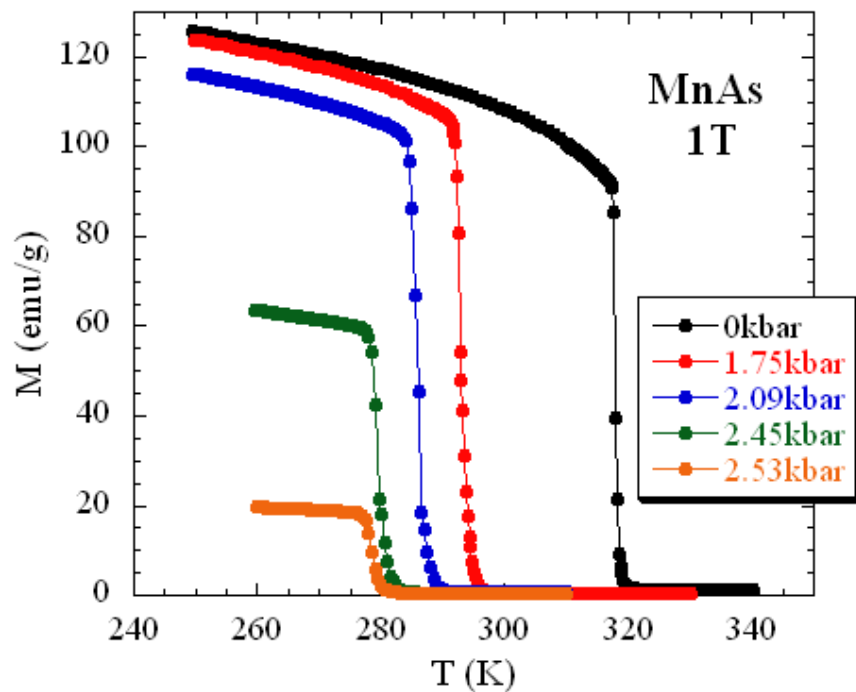
decreasing p

B8₁: NiAs-type
B31: MnP-type

increasing p

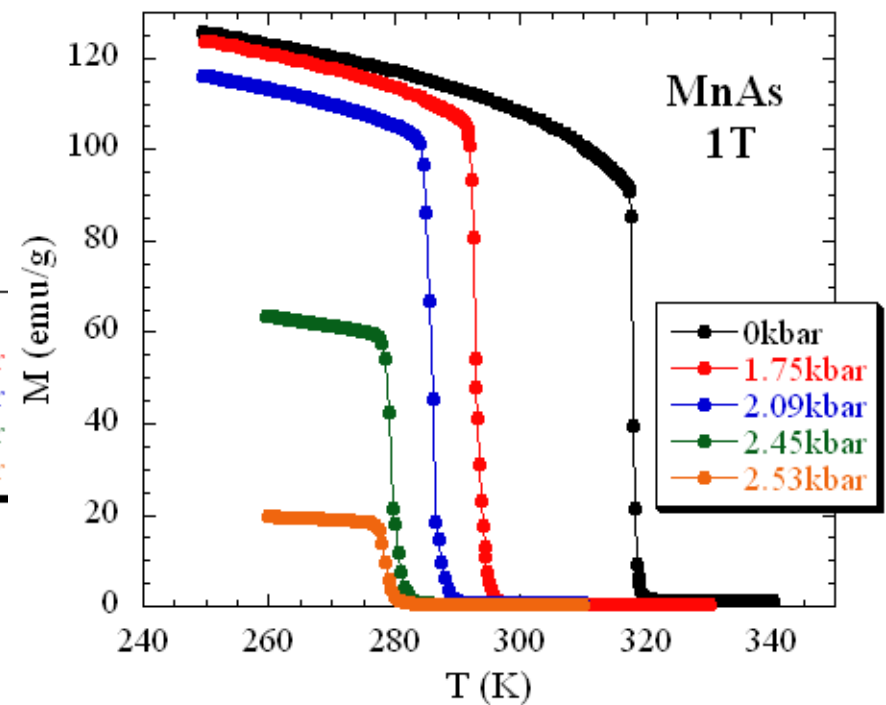
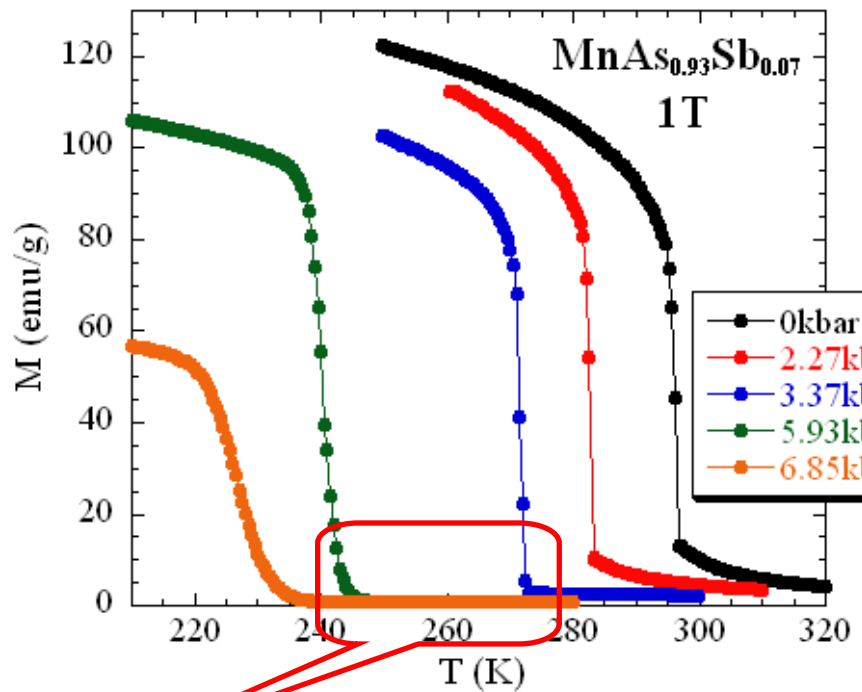


MCE of MnAs under pressure



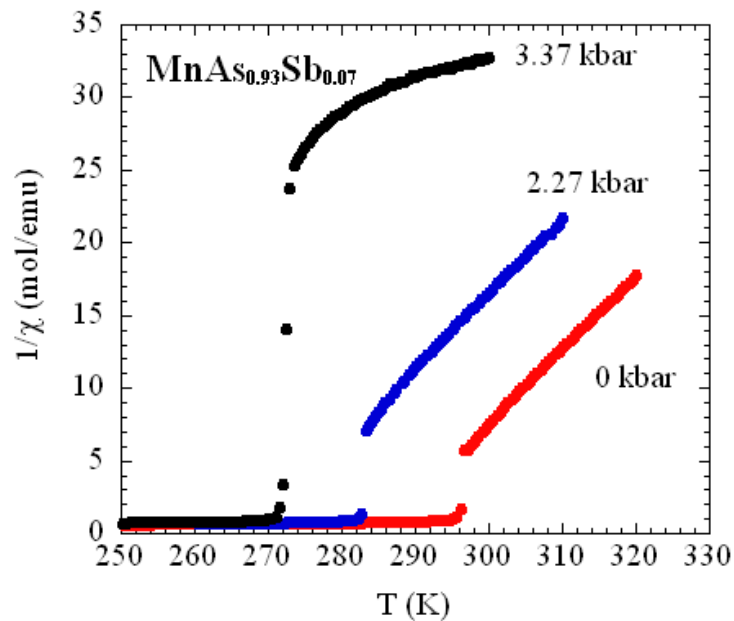
No strong enhancement of MCE by pressure was observed.

Magnetic transition of $\text{MnAs}_{0.93}\text{Sb}_{0.07}$ under pressure

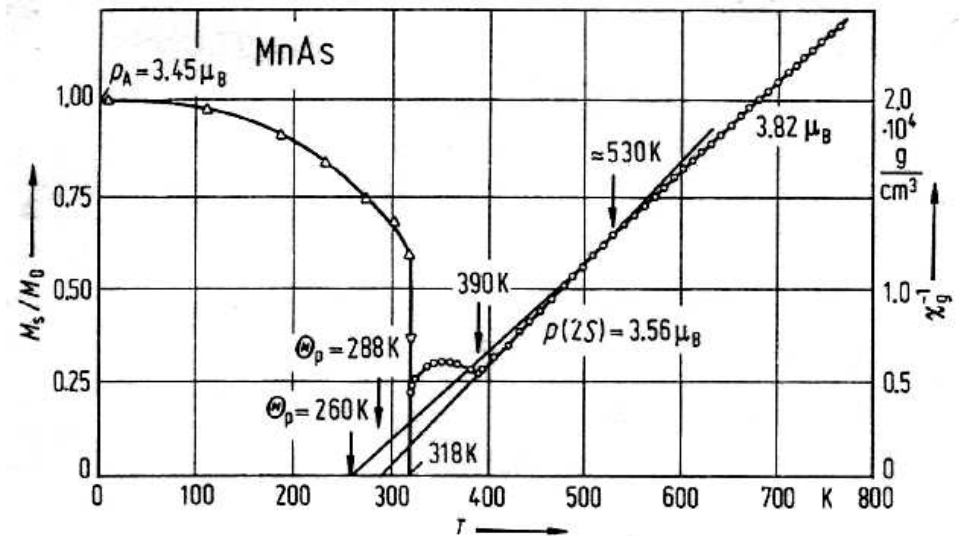


suggesting a structural transformation

χ^{-1} - T curve of $\text{MnAs}_{0.93}\text{Sb}_{0.07}$



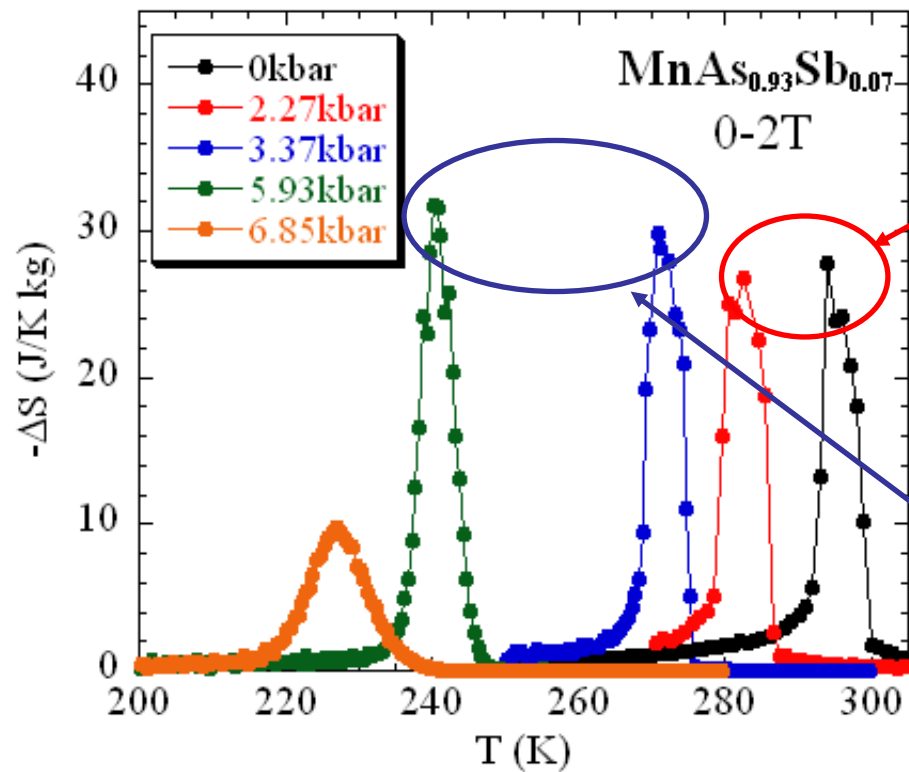
$\mu_{\text{eff}} = 4.0\mu_{\text{B}}/\text{Mn}$ for 0 kbar



$\mu_{\text{eff}} = 4.45\mu_{\text{B}}/\text{Mn}$ for MnAs

$\text{MnAs}_{0.93}\text{Sb}_{0.07}$ undergoes the first-order NiAs-type \rightarrow MnP-type transition above 3.37 kbar.

MCE of $\text{MnAs}_{0.93}\text{Sb}_{0.07}$ under pressure



NiAs → NiAs

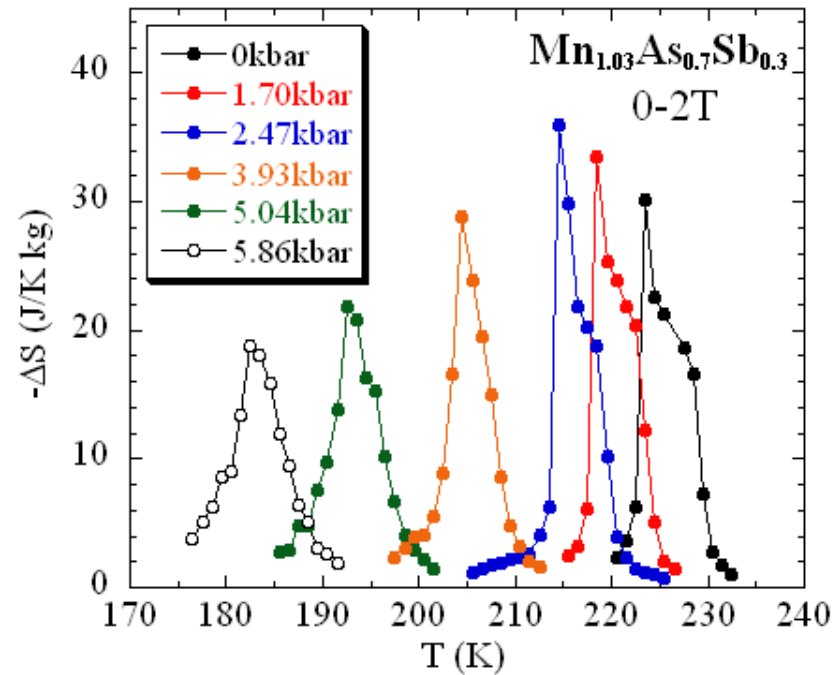
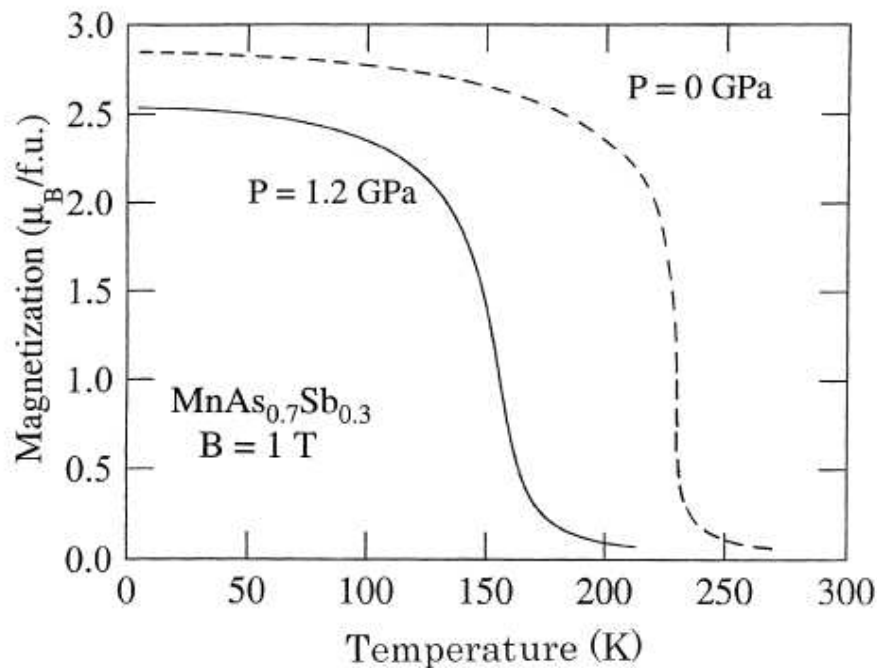
$$\Delta S = \Delta S_M + \Delta S_{st}$$

NiAs → MnP

- Lattice entropy change due to a structural transformation, ΔS_{st} is small in the $\text{MnAs}_{1-x}\text{Sb}_x$ system.

MCE of $\text{MnAs}_{0.7}\text{Sb}_{0.3}$ under pressure

Goto *et al.*, 2001



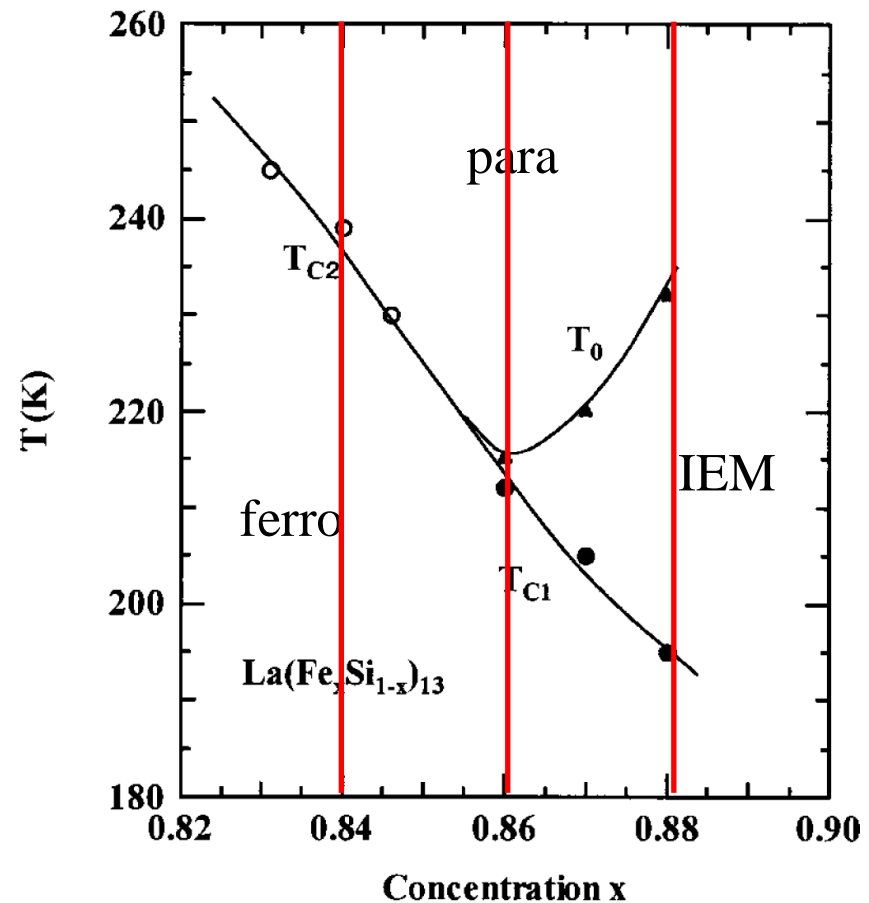
The NiAs-type structure is stable up to 12 kbar.

Magnetic phase diagram of $\text{La}(\text{Fe}_x\text{Si}_{1-x})_{13}$

Compounds are formed
in $0.81 \leq x \leq 0.90$.

The FOMT appears in
 $x > 0.86$.

We studied MCE
under pressure for
 $x=0.84$ (SOMT)
 0.86 (critical)
 0.88 (FOMT)



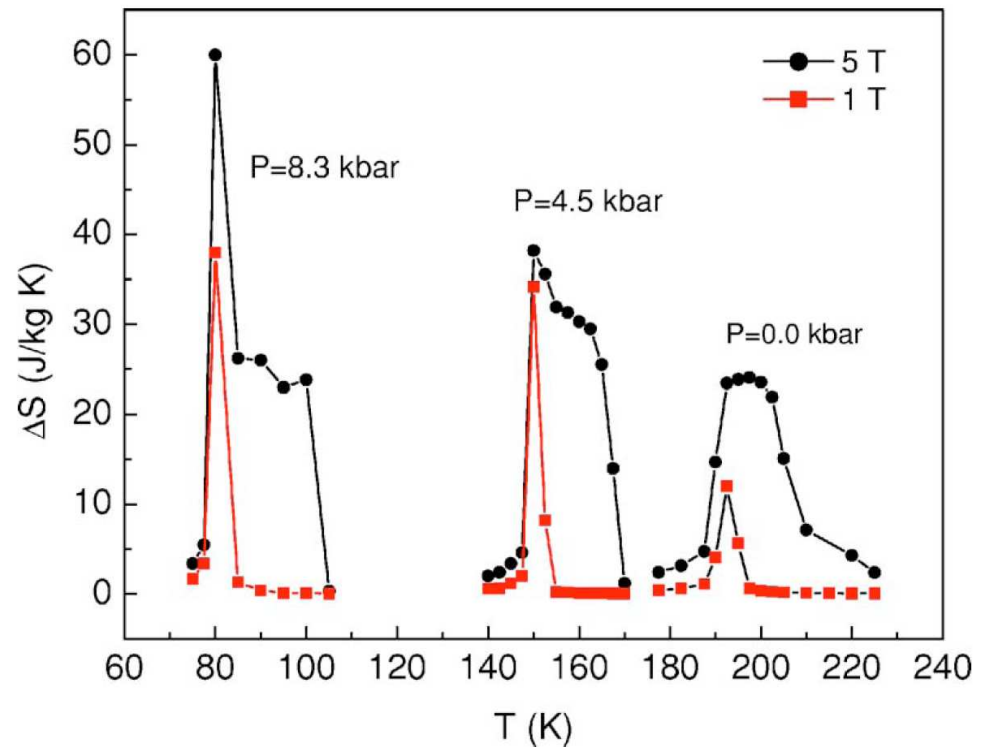
Fujita *et al.*, 2001

Previous reports

Previous measurements
are concentrated on the
FOMT.

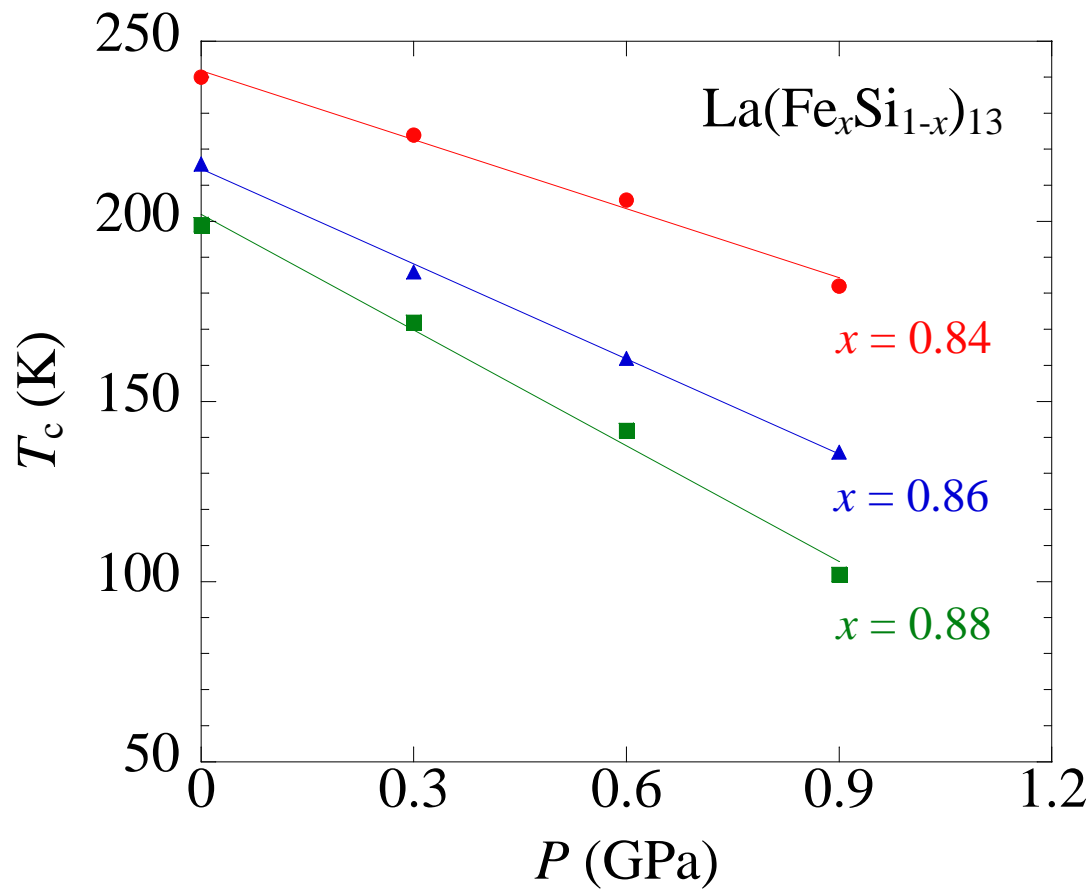
$\text{LaFe}_{11.6}\text{Si}_{1.4}$ ($x=0.892$)
(Sun *et al.*, 2006)

$\text{LaFe}_{11.5}\text{Si}_{1.5}$ ($x=0.885$)
(Jia *et al.*, 2007).



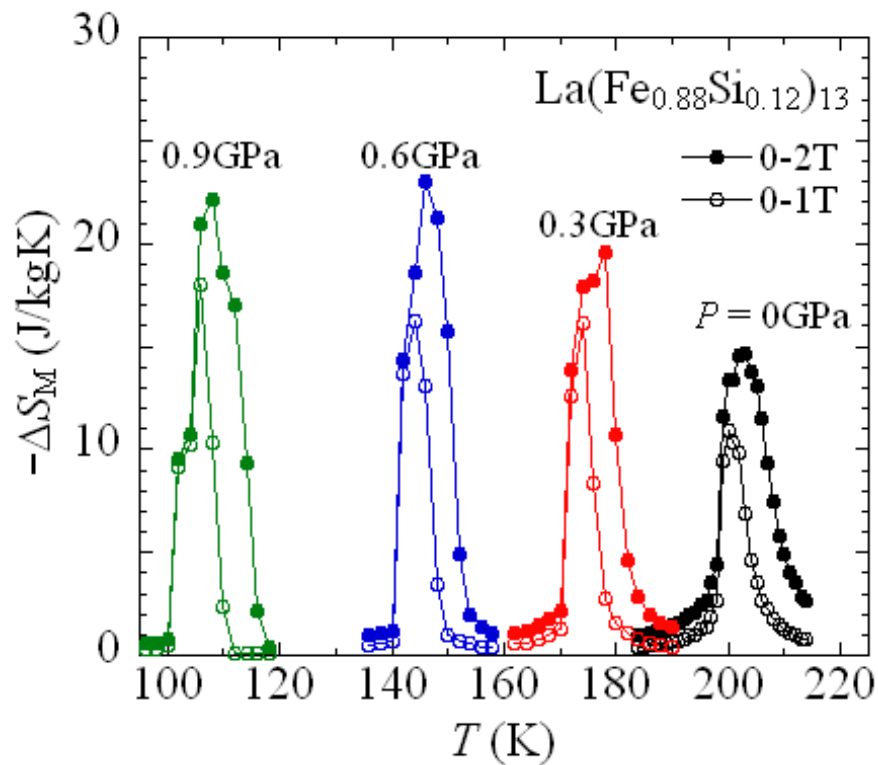
Sun *et al.*, 2006

Pressure dependence of T_C

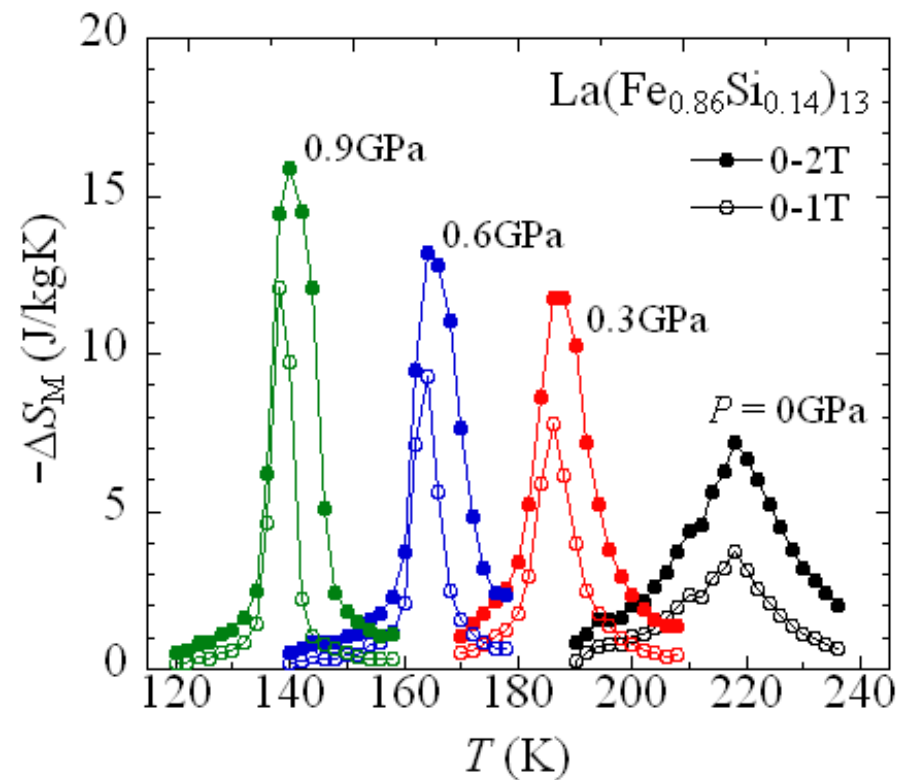


The results are in agreement with the previous reports by Fuita *et al.* (2001).

Pressure dependence of MCE for $x=0.88$ and 0.86

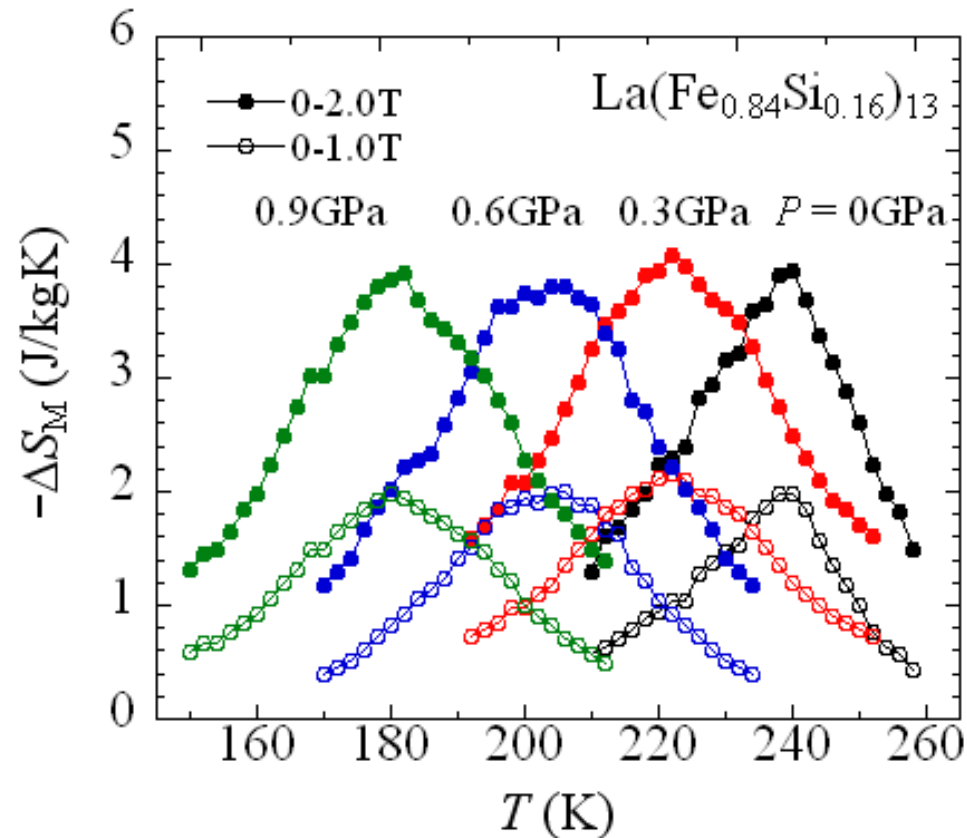


FOMT



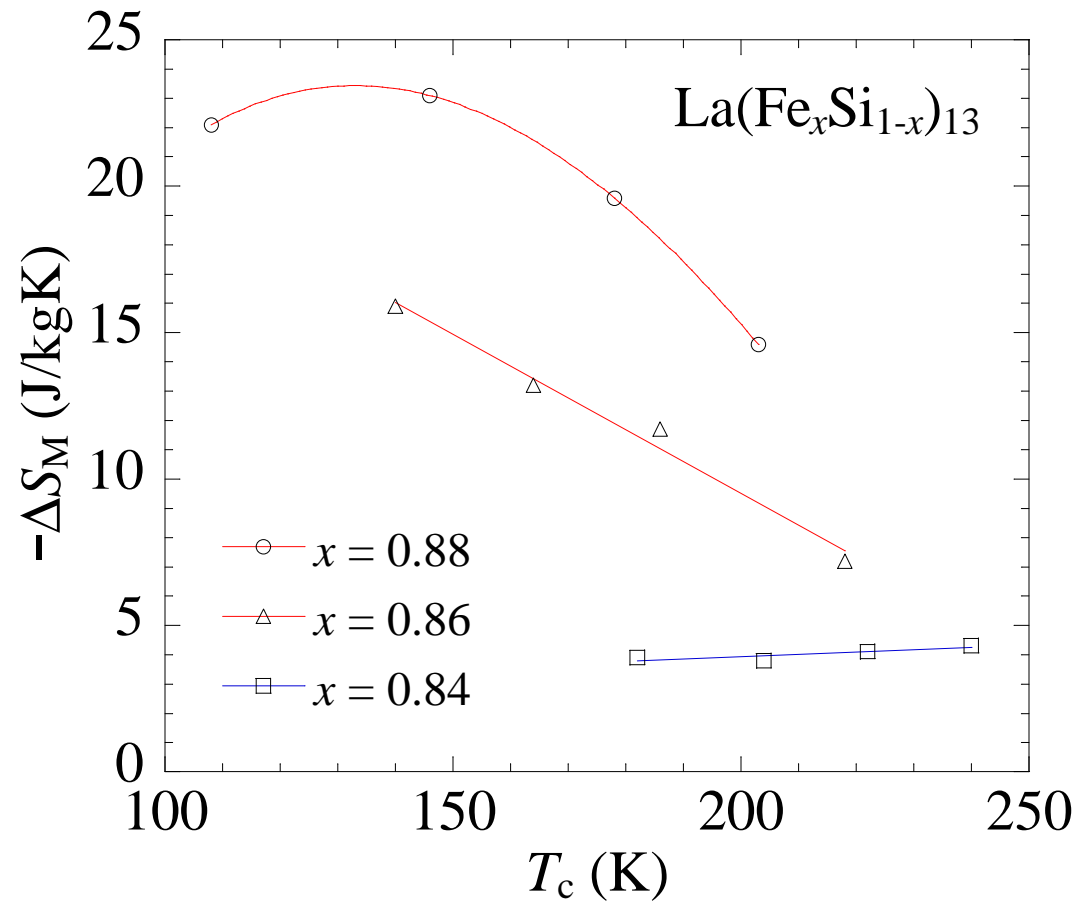
SOMT \rightarrow FOMT

Pressure dependence of MCE for $x=0.84$



The SOMT persists up to 0.9 GPa.

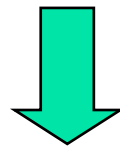
Summary of ΔS of $\text{La}(\text{Fe}_x\text{Si}_{1-x})_{13}$ under pressure





Conclusions : $\text{MnAs}_{1-x}\text{Sb}_x$

- No strong enhancement of the MCE was observed.
- Pressure stabilizes the MnP-type structure.
- $\text{MnAs}_{0.93}\text{Sb}_{0.07}$ undergoes the first-order NiAs-type \rightarrow MnP-type transition above 3.37 kbar.
- The lattice entropy change due to a structural transformation is small.



Broadening of the FOMT seems difficult.



Conclusions : $\text{La}(\text{Fe}_x\text{Si}_{1-x})_{13}$

- The pressure effect on the MCE strongly depends on x .
- A FOMT is induced by pressure for $x = 0.86$, which enhances the peak value of ΔS .



Broadening of the FOMT seems possible.