

Addendum

Addendum to “Novel features in solid phase transition behavior of TII [E.A. Secco, *Thermochim. Acta* 342 (1999) 161]”

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A report by Samara et al. on the transformation of yellow  $\beta$ -TII, orthorhombic structure (*Cmcm* space group), to red  $\alpha$ -TII, cubic CsCl-type structure (*Pm3m* space group) was missed in our literature search [1]. This report contains the heat of transition  $\Delta H_t$ , and the transition temperature of TII,  $T_t$ , on a powdered sample obtained from a single crystal of TII.

The transition temperature was determined at 1 bar to be 170°C. The quantitative value for the heat of transition at atmospheric pressure was obtained by comparing the peak areas of the DTA signals of the sample with the literature values for the two standard transitions in  $\text{AgNO}_3$ , viz., the solid–solid transition and the solid–liquid (fusion) transition. From these two standard values the  $\Delta H_t$  of TII was calculated to be 1120 and 1226 J/mol, respectively, giving an average value  $\Delta H_t=1173$  J/mol. This  $\Delta H_t$  value is in excellent agreement with our  $\Delta H_t=1180\pm 50$  J/mol cited in the above reference.

The existing literature data on the solid phase transition of TII from different sources are collected in Table 1. The data suggest an involvement of two factors each independently affecting the transition temperature  $T$  and the transition enthalpy  $\Delta H_t$ . One

Table 1  
Solid phase transition data of TII

Transition	Reference
Temperature, $T$ (°C)	Enthalpy, $\Delta H_t$ (vol./mol)
170	1173±50 [1]
172±2	875±40 [2,3]
181±2	1180±50 [4]
175	– [5]

factor relates to the sample, viz. particle size or structure (metastable) and the other implicates the calorimetric quantitative analysis method.

References

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