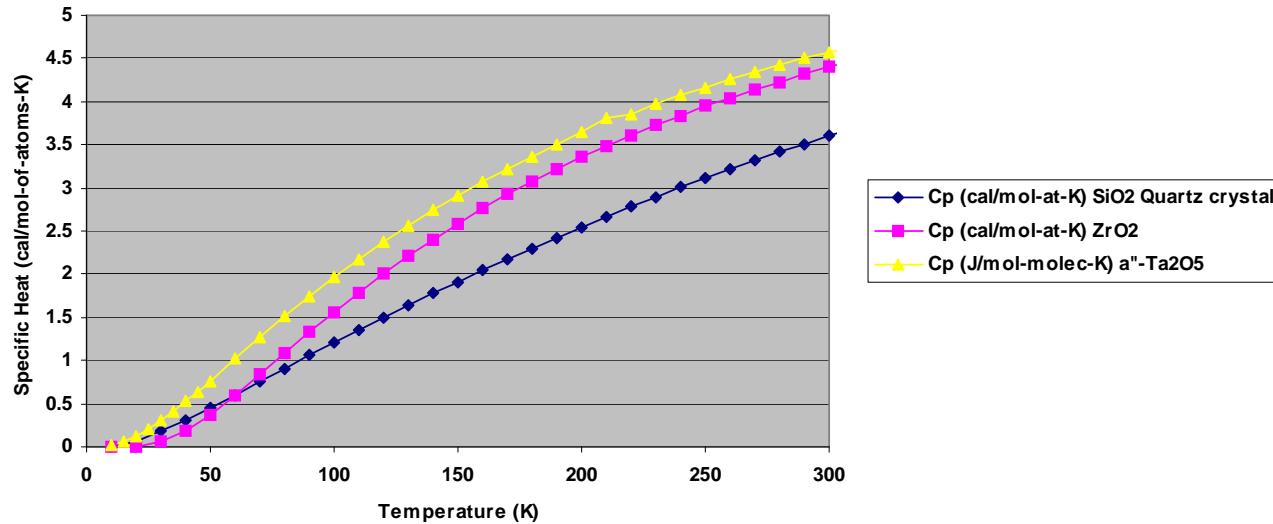
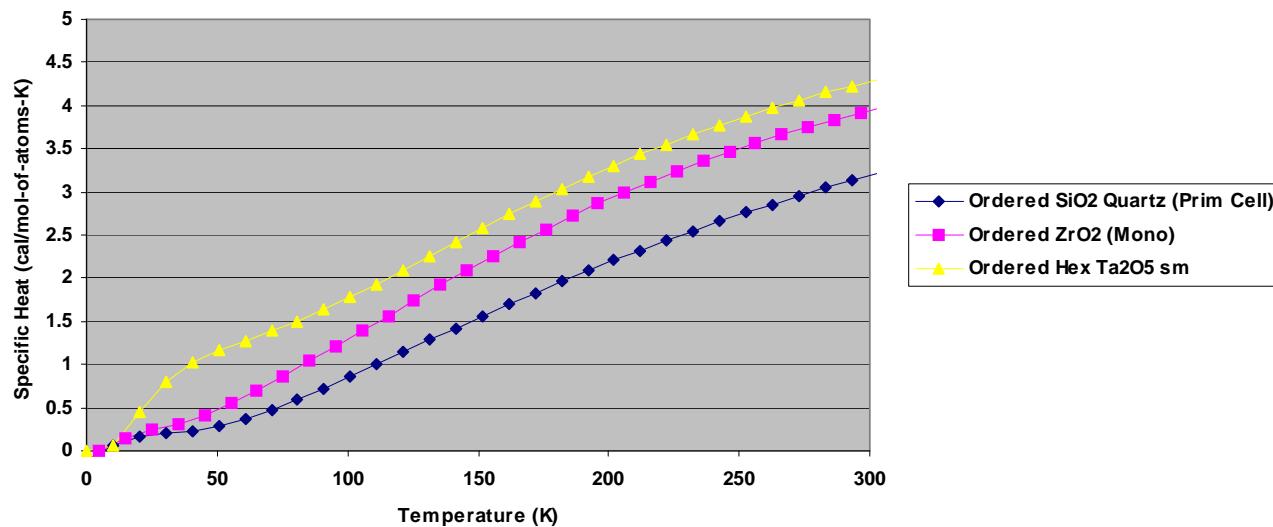


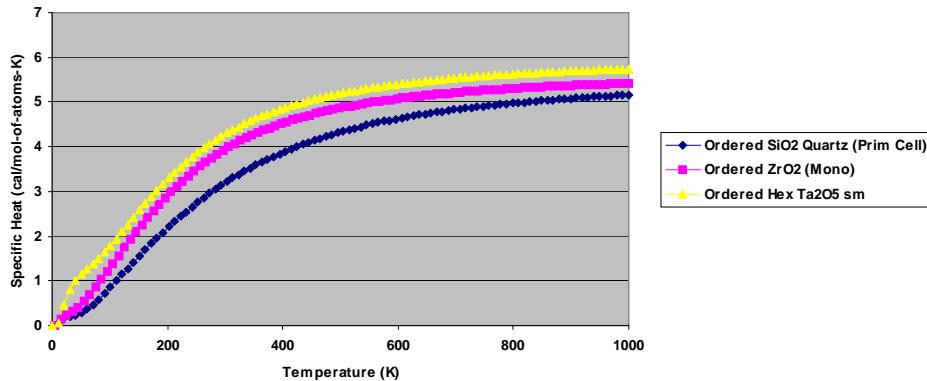
Measured Values from Literature



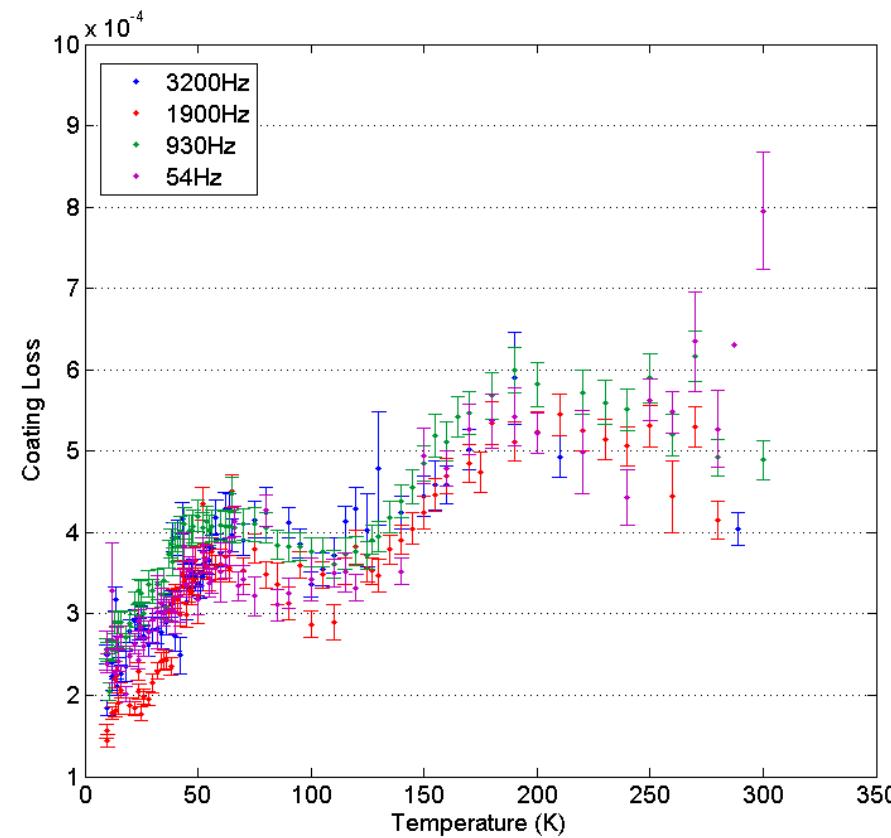
Ab Initio Calculation



Ab Initio Calculation



Loss angle and
per-mole-of-atom specific heat capacity
look a lot alike,
in as far as temperature.



Nawrodt Hafnia Curve

Where I got data.

Coating material	Absorption (ppm)	Mechanical losses
Ta ₂ O ₅	0.5	$\Phi_{Ta_2O_5} = 3.0 \cdot 10^{-4}$
Ta ₂ O ₅ : W	2.45	$\Phi_{Ta_2O_5W} = 7.5 \cdot 10^{-4} *$
Ta ₂ O ₅ : W+Ti	1.65	$\Phi_{Ta_2O_5W+Ti} = 3.2 \cdot 10^{-4}$
Ta ₂ O ₅ : Co	5000	$\Phi_{Ta_2O_5Co} = 1.1 \cdot 10^{-3} *$
ZrO ₂	11	$\Phi_{ZrO_2} = 2.3 \cdot 10^{-4}$
ZrO ₂ : Ti	37	$\Phi_{ZrO_2Ti} = 6.8 \cdot 10^{-4}$
ZrO ₂ : W	10	$\Phi_{ZrO_2W} = 2.8 \cdot 10^{-4}$

**Table 1: optical absorption and mechanical loss of various coating materials
(coating thickness 500nm, * indicates before annealing)**

Found on the internet, but
the data comes from Glasgow.

- HfO₂ loss angle 300K from Nawrodt Hafnia presentation ~ 5E-4.
- Al₂O₃ loss angle 6E-5 from Yamamoto talk, called Glasgow Ta₂O₅ / Al₂O₃. Yamamoto value is 4E-4, off trend !
- sapphire Al₂O₃ Cp 300K from J. Nuc. Sci. Tech 7 [6] 312-316 and also TA Instruments Data Sheet.
- SiO₂ coating loss angle 300K from several places in DCC ~ 5E-5.
- HfO₂ monoclinic 300K Cp from J. Ceram. Soc. 89 12 3751-3578 (2006).
- a"-Ta₂O₅ Cp function from J. Russian Phys. Chem. 61 (2) 361-365 (1987).
- SiO₂, ZrO₂ Cp function from Latin American Applied Research 34 257-265 (2004).
- Ab initio code for phonon calculations provided by Accelrys, \$150K worth of code on a temp license. CASTEP, all GGA-PW91 functional.
- Accelrys also pointed out that Cv output is per cell and needs to be divided by atoms per cell – trend then obvious.