# Thermoelectricity at nanoscale: theoretical models



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### Joint work with Oleg Zhirov (BINP) Europhys. Lett. 103, 68008 (2013)



Time evolution of figure of merit *ZT* (center) (from A.Majumdar Science **303**, 777 (2004))

## **Early works**



from Y.Imry (Weizmann Inst) talk at Inst. H. Poincaré, Paris (2012)

# **Early works**





Oil burning lamp powering a radio using the first commercial thermoelectric generator containing ZnSb built in USSR, circa 1948

from Y.Imry (Weizmann Inst) talk at Inst. H. Poincaré, Paris (2012)

# **Main characteristics**

### Seebeck coefficient:

 $S = \Delta V / \Delta T = \pi^2 k_B^2 T[(d \ln \sigma / dE)]|_{E_F} / e \text{ (Mott relation (1958))}$ 

For 2DEG with Wiedemann-Franz law:  $S = 2\pi k_B^2 Tm/(3eh^2 n_e)$ ; typical value  $S \approx 10 \mu V/K$  at T = 0.3K,  $n_e = 4 \cdot 10^{10} cm^{-2}$ 



Thermoelectric figure of merit  $ZT = \sigma S^2 T/\kappa$ , thermal conductivity  $\kappa = \kappa_{el} + \kappa_{phonon}$  (heat flux  $Q = -\kappa \bigtriangledown T$ )

# **Experiments on Seebeck coefficient for 2DEG**



R.Fletcher, V.M.Pudalov et al. Semicond. Sci. Technol. 15, 386 (2001)

# Experiments on Seebeck coefficient for 2DEG



FIG. 5. (Color online) (a) S vs  $n_t$  for 0.28 K < T < 0.7 K. The broken green line shows  $S_d$  [Eq. (2)] at 0.28 K. Inset:  $\rho_{2DES}$  vs  $n_s$  at the same T values; there is little T dependence in this range. (b) Low-T linear variation of S. Inset: Descriptions based on variable-ranged hopping, where S is expected to decay to zero as  $T^{1/3}$ , do not adequately describe the observed data.

V.Narayan, S.Goswami, M.Pepper et al. PRB **85**, 125406 (2012) In dimensionless units  $S = 10 mV/K \approx 100 \gg 1$ 

# **ZT** in various materials



from http://www.thermoelectrics.caltech.edu/thermoelectrics/

# ZT in p-type $Bi_2Te_3/Sb_2Te_3$ superlattices

#### Phonon-blocking/electron-transmitting structures

The results obtained with the  $10\text{\AA}/50\text{\AA}$  Bi<sub>2</sub>Te<sub>3</sub>/Sb<sub>2</sub>Te<sub>3</sub> superlattices indicate that we can fine-tune the phonon and hole (charge carriers)



Figure 3 Temperature dependence of ZT of 10Å/50Å p-type  $Bi_2Te_3/Sb_2Te_3$  superlattice compared to those of several recently reported materials.

## R.Venkatasubramanian et al. (N.Carolina) Nature 413, 597 (2001)

# ZT in SnSe



Figure 1 | SnSe crystal structure *Pnma* and *ZT* values. a, Crystal structure along the *a* axis: grey, Sn atoms; red, Se atoms. b, Highly distorted SnSe<sub>7</sub> coordination polyhedron with three short and four long Sn–Se bonds. c, Structure along the *b* axis. d, Structure along the *c* axis. e, Main panel, *ZT* values along different axial directions; the *ZT* measurement uncertainty is about 15% (error bars). Inset images: left, a typical crystal; right, a crystal cleaved along the (000) plane, and specimens cut along the three axes and corresponding measurement directions. Inset diagram, how crystals were cut for directional measurements; *ZT* values are shown on the blue, red and grey arrows; colours represent specimens oriented in different directions.

#### Li-Dong Zhao et al. (Illinois) Nature 508, 373 (2014)

# **Research at CNRS Grenoble (O.Bourgeois Refs.3,4**)



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# Applications

#### TE Applications are mostly 'Niche' Applications



# Thermoelectricity of Wigner crystal in a periodic potential

Hamiltonian  $H = \sum_{i} \left( \frac{p_{i}^{2}}{2} + K \cos x_{i} + \frac{1}{2} \sum_{j \neq i} \frac{1}{|x_{i} - x_{j}|} \right)$ Dynamic equations  $\dot{p}_{i} = -\partial H / \partial x_{i} + E_{dc} - \eta p_{i} + g\xi_{i}(t)$ ,  $\dot{x}_{i} = p_{i}$ Here the Langevin force is given by  $g = \sqrt{2\eta T}$ ,  $\langle \xi_{i}(t)\xi_{j}(t') \rangle = \delta_{ij}\delta(t - t')$ ;  $n_{e} = \nu/2\pi$ ,  $\nu = \nu_{g} = 1.618...$  Fibonacci rational approximates. Aubry transition at  $K = K_{c} = 0.0462$ . I.Garcia-Mata, O.Zhirov, DLS EPJD **41**, 325 (2007)

# Time evolution of Wigner crystal



Electron density variation in space and time from one Langevin trajectory at  $K/K_c = 2.6$ ,  $T/K_c = 0.11$ ,  $\eta = 0.02$ , N = 34,  $M = L/2\pi = 21$ ; density changes from zero (dark blue) to maximal density (dark red); only a fragment of *x* space is shown.

## **Numerical fits**



Left panels: dependence of electron temperature  $T_e(x)$  (*top, blue points*) and rescaled density  $\nu(x)$  (*bottom, black points*) on distance x along the chain placed on the Langevin substrate with a constant temperature gradient (it is shown by the blue line) at average temperature  $\overline{T} = 0.01$  and temperature difference  $\Delta T = 0.2\overline{T}$ ; black line shows the fit of density variation in the bulk part of the sample. *Right panel*: density variation produced by a static electric field  $E_{dC} = 4 \times 10^{-4}$  at a constant substrate temperature T = 0.01; black line shows the fit of gradient in the bulk part of the sample. Here N = 34, M = 21,  $K = 1.52K_C$ ,  $\eta = 0.02$ , averaging is done over time interval  $t = 10^{-7}$ ; S = 3.3 at  $T = 0.01 \approx 0.22K_C$ .

## Seebeck coefficient



Left panel: Dependence of the Seebeck coefficient *S* on rescaled potential amplitude  $K/K_c$  at temperatures  $T/K_c = 0.065, 0.11, 0.22$  and 0.65 shown by black, blue, green and red colors, respectively from top to bottom. The full and open symbols correspond respectively to chains with N = 34, M = 21and N = 55, M = 34. *Right panel*: Dependence of *S* on  $T/K_c$  at different  $K/K_c = 0, 0.75, 1.5, 2.2, 3$  shown respectively by black, violet, blue, green and red points; N = 34, M = 21; the dashed gray line shows the case K = 0 for noninteracting particles. The stars show corresponding results from left plane at same *N*, *M*. Dotted curves are drown to adapt an eye. Here and in other Figs. the statistical error bars are shown when they are larger than the symbol size. Here  $\eta = 0.02$ .

# Conductivity and thermal conductivity



*Left panel*: Rescaled electron conductivity  $\sigma/\sigma_0$  as a function of  $K/K_c$  shown at rescaled temperatures  $T/K_c = 0.065$ , 0.22, 0.65 by black, green and red points respectively. *Right panel*: Rescaled thermal conductivity  $\kappa/\kappa_0$  as a function of  $K/K_c$  shown at same temperatures and colors as in left panel. Here we have N = 34, M = 21,  $\eta = 0.02$ ,  $\sigma_0 = \nu_a/(2\pi\eta)$ ,  $\kappa_0 = \sigma_0 K_c$ .

# **ZT** dependence on parameters



*Left panels*: Dependence of *ZT* on  $K/K_c$  at temperatures  $T/K_c = 0.11$  (top panel) and  $T/K_c = 0.65$  (bottom panel); the black points and open triangles correspond respectively to  $\eta = 0.02$  and  $\eta = 0.05$  at N = 34, M = 21. *Right panels*: Dependence of *ZT* on  $T/K_c$  for  $K/K_c = 0.75$  at  $\eta = 0.02$ , N = 34, M = 21. *Bottom right panel*: Same as in top right panel at  $K/K_c = 2.6$  and N = 34, M = 21 (black points); N = 89, M = 55 (green circles); N = 144, M = 89 (red stars).

# **ZT** dependence on parameters



Dependence of ZT on  $K/K_c$  and  $T/K_c$  shown by color changing from ZT = 0 (black) to maximal ZT = 4.5 (light rose); contour curves show values ZT = 1, 2, 3, 4. Here  $\eta = 0.02$ , N = 34, M = 21.

# Other quantities dependences



*Left panel*: Rescaled thermal conductivity  $\kappa/\kappa_0$  as a function of rescaler temperature  $T/K_c$ , to adapt an eye the straight dashed line shows the dependence  $\kappa/\kappa_0 = 0.6T/K_c$ ; *right panel*: same as in left panel for  $S^2\sigma/\sigma_0$ . Data are obtained at  $K/K_c = 2.6$ ,  $\eta = 0.02$ , N = 34, M = 21,  $\sigma_0 = \nu_g/(2\pi\eta)$ ,  $\kappa_0 = \sigma_0 K_c$ .

In physical units we can estimate the critical potential amplitude as  $U_c = K_c e^2/(\epsilon d)$ , where  $\epsilon$  is a dielectric constant,  $\Delta x$  is a lattice period and  $d = \nu \Delta x/2\pi$  is a rescaled lattice constant Ref.5. For values typical for a charge density wave regime Ref.6 we have  $\epsilon \sim 10$ ,  $\nu \sim 1$ ,  $\Delta x \sim 1 nm$  and  $U_c \sim 40 mV \sim 500K$  so that the Aubry pinned phase should be visible at room temperature. The obtained  $U_c$  value is rather high that justifies the fact that we investigated thermoelectricity in the frame of classical mechanics of interacting electrons.

# Message from Novosibirsk GES



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R2. H.J. Goldsmid, *Introduction to thermoelectricity*, Springer, Berlin (2009).
R3. Nanoelectronics : Concepts, Theory and Modeling. Network meeting and workshop on thermoelectric transport 21-27 October 2012, Cargese, Corsica http://iramis.cea.fr/meetings/nanoctm/program.php
R4. Slides at R3 by O.Bourgeois, L.W.Molenkamp, S.Voltz
R5. I.Garcia-Mata, O.V.Zhirov, D.L.Shepelyansky, EPJD 41, 325 (2007)
R6. S.Brazovskii *et al.*, Phys. Rev. Lett. 108, 096801 (2012)