

Electrostatic Doping in Carbon-based Nanoelectronics Devices

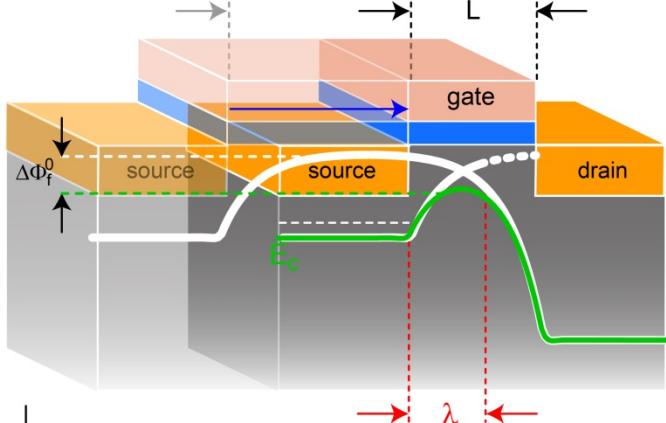
Joachim Knoch

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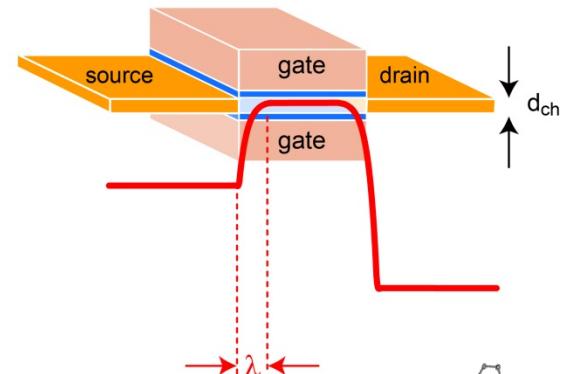


Introduction

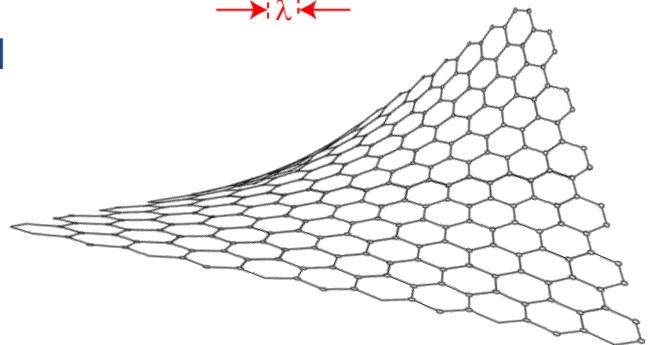
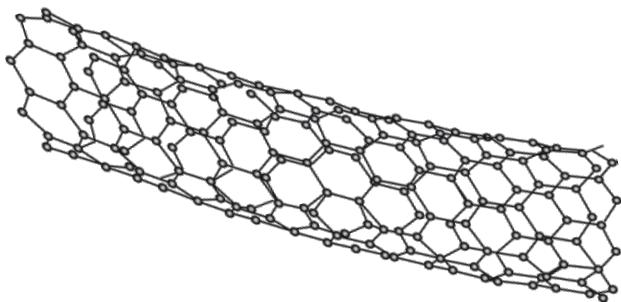
$$\lambda = \sqrt{\frac{\varepsilon_{\text{nw}} d_{\text{nw}}^2}{8\varepsilon_{\text{ox}}} \ln \left(1 + \frac{2d_{\text{ox}}}{d_{\text{nw}}} \right)}$$



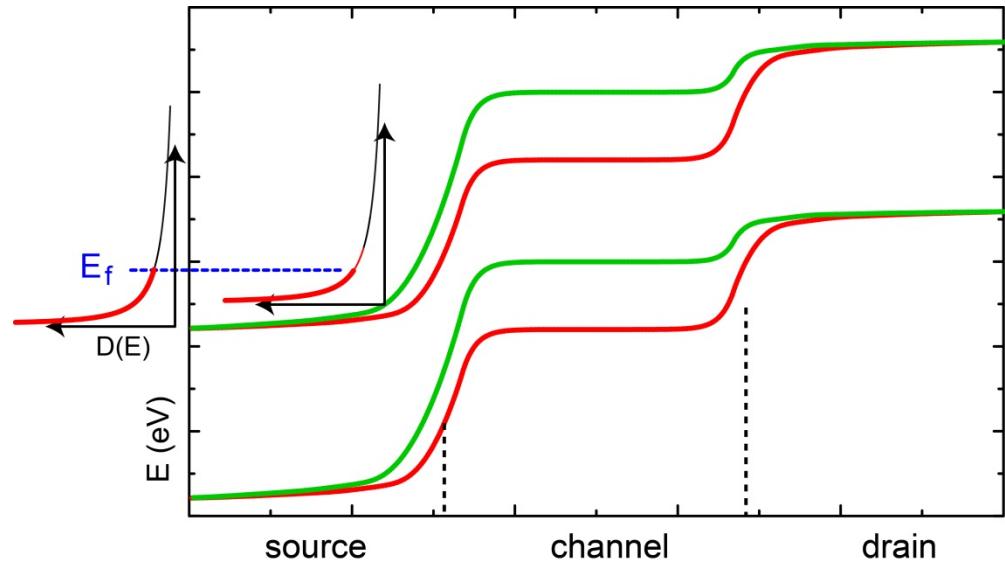
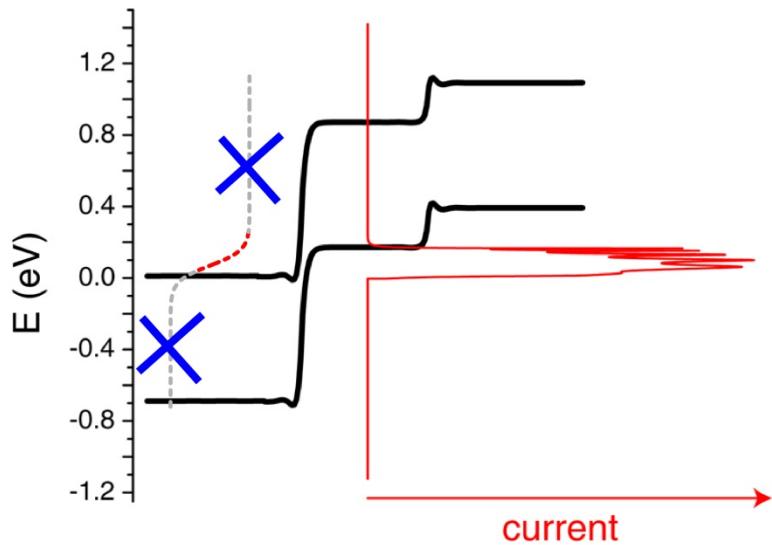
$$\lambda_{\text{ch}} = \sqrt{\frac{\varepsilon_{\text{ch}}}{\varepsilon_{\text{ox}}} d_{\text{ox}} d_{\text{ch}}}$$



carbon-based materials
hold promise to realize
ultimate FETs due to
optimum scalability and
electronic transport
properties



Introduction – Issues with Doping



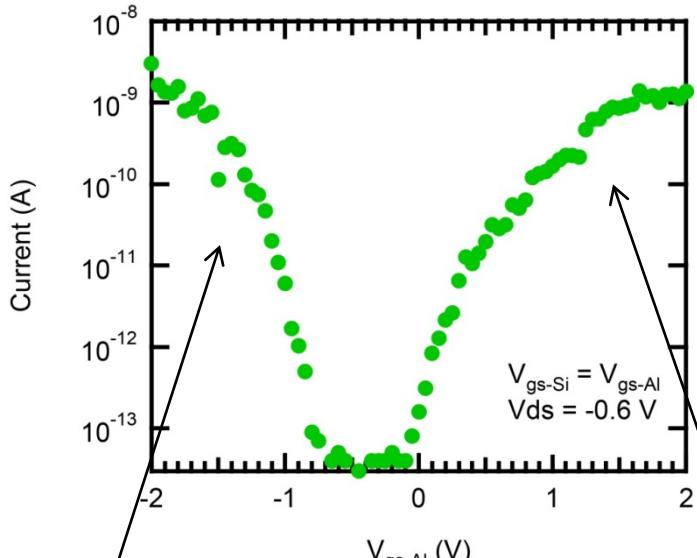
- need excellent gate control of channel and screening of gate impact on source contact
- low density of states such as in carbon nanotubes (1D) or graphene detrimental for screening

→ electrostatic doping

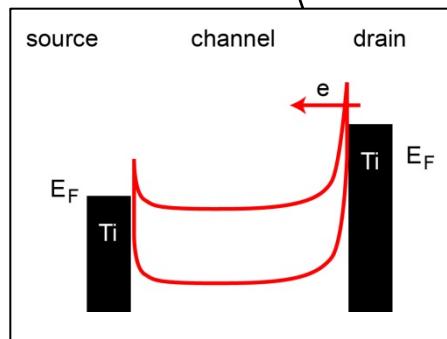
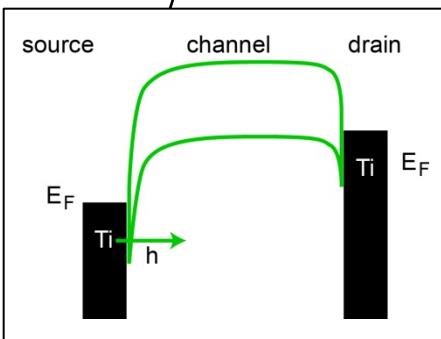
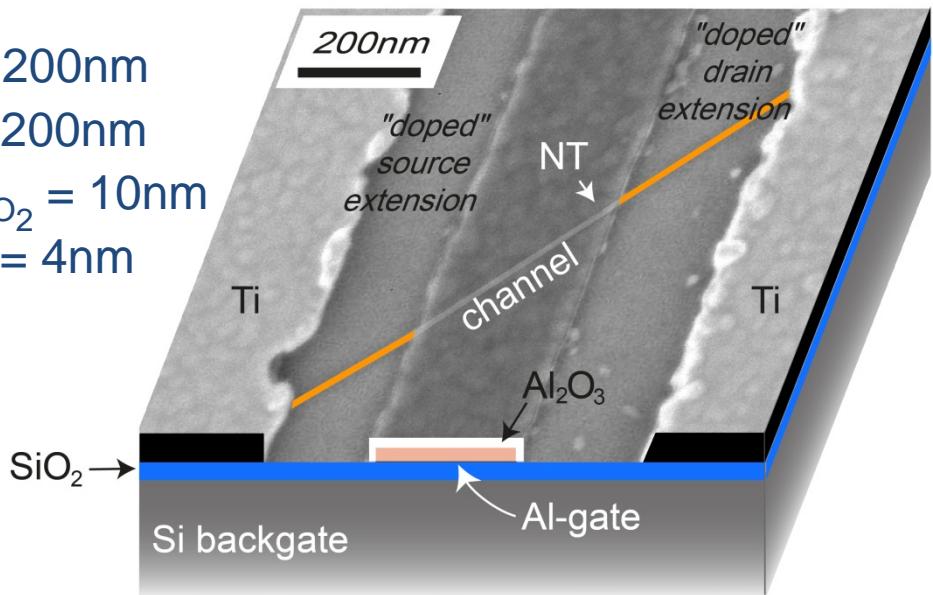
J. Knoch and J. Appenzeller, phys. stat solidi a, **205**, 679 (2008).

J. Knoch, S. Mantl and J. Appenzeller, Solid-State Electron., **51**, 572 (2007).

Tunable Polarity in CNFETs



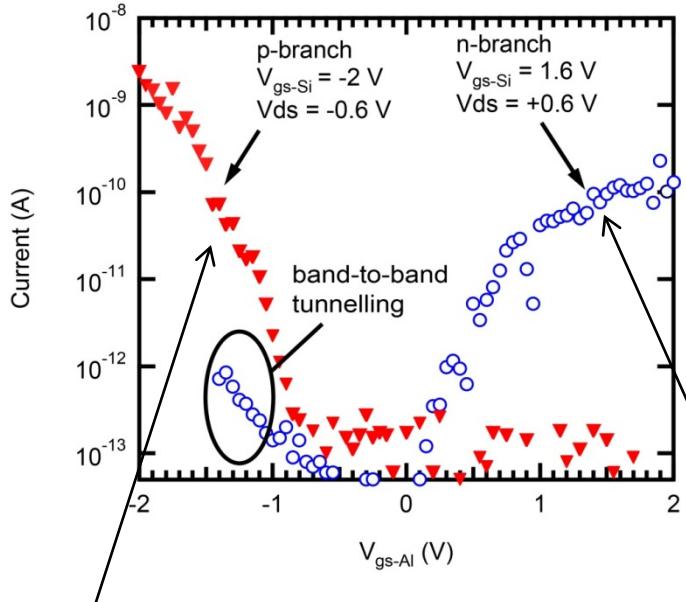
- $L_{bg} \approx 200\text{nm}$
- $L_{ch} \approx 200\text{nm}$
- $d_{ox\text{-SiO}_2} = 10\text{nm}$
- $d_{ox\text{-Al}} = 4\text{nm}$



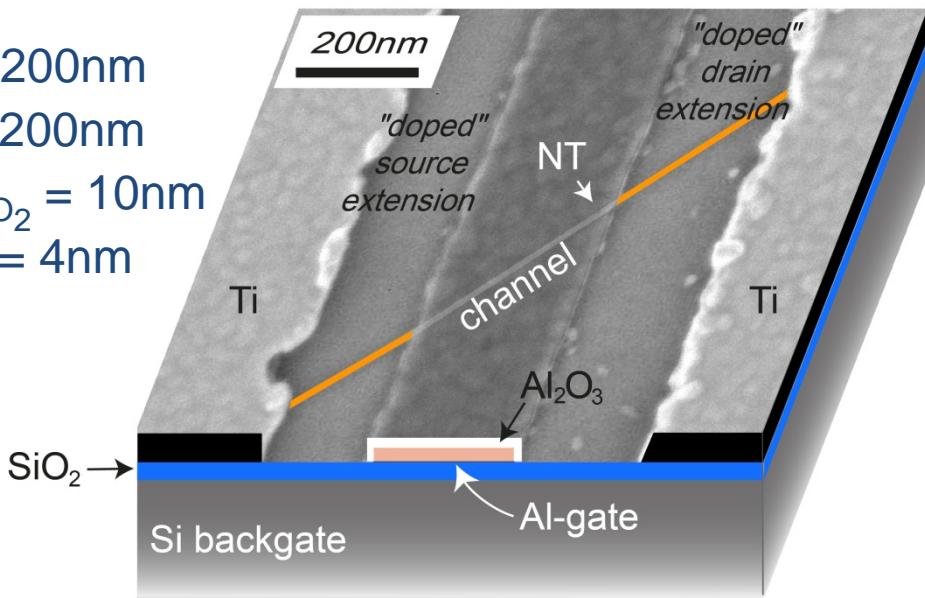
- **ambipolar behavior if both gates biased in the same direction**
- **electron and hole injection through Schottky-barrier at Ti-nanotube interface**

Y.-M. Lin, J. Appenzeller, J. Knoch and Ph. Avouris, IEEE Trans. Nanotechnol., 4, 481 (2005).

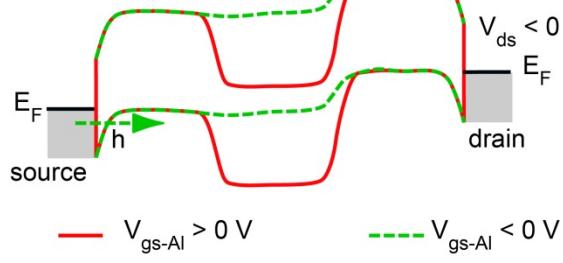
Tunable Polarity in CNFETs



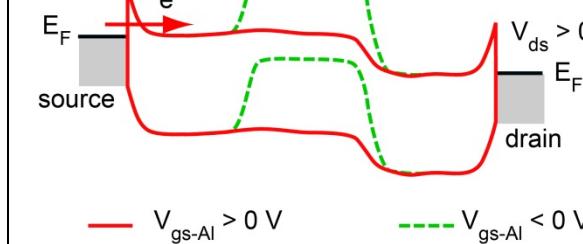
- $L_{bg} \approx 200$ nm
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$V_{gs\text{-Si}} < 0$ (p/i/p) p-FET



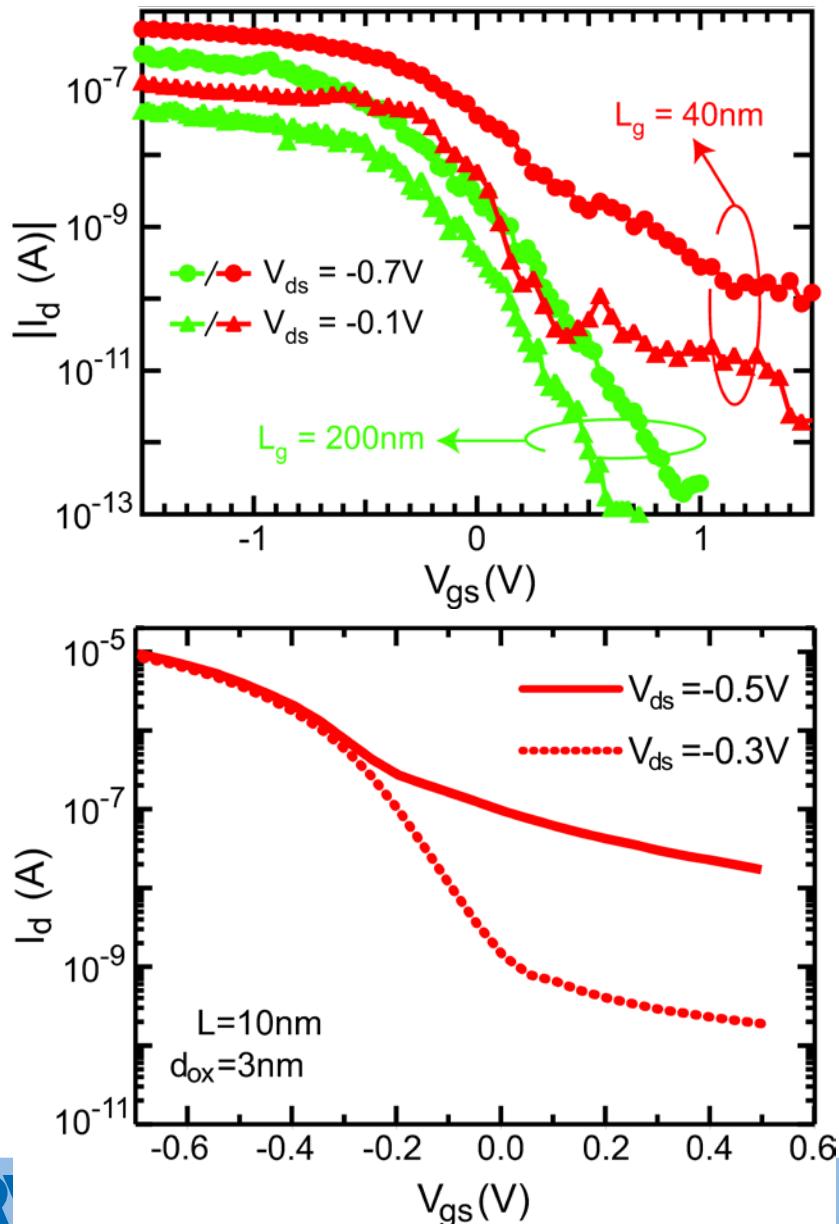
$V_{gs\text{-Si}} > 0$ (n/i/n) n-FET



- **unipolar device behavior if both gates are biased with opposite polarity**
- **n- and p-type CNFETs possible due to excellent gate control**

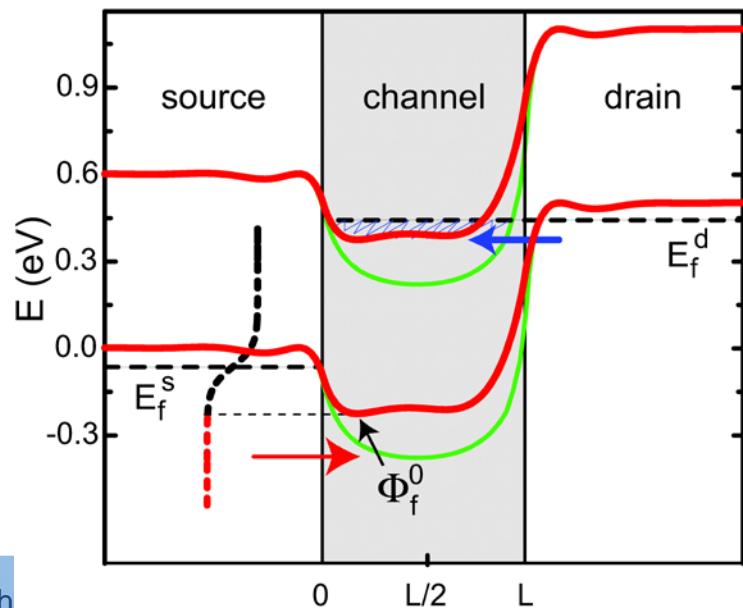
Y.-M. Lin, J. Appenzeller, J. Knoch and Ph. Avouris, IEEE Trans. Nanotechnol., 4, 481 (2005).

Conventional CNFETs

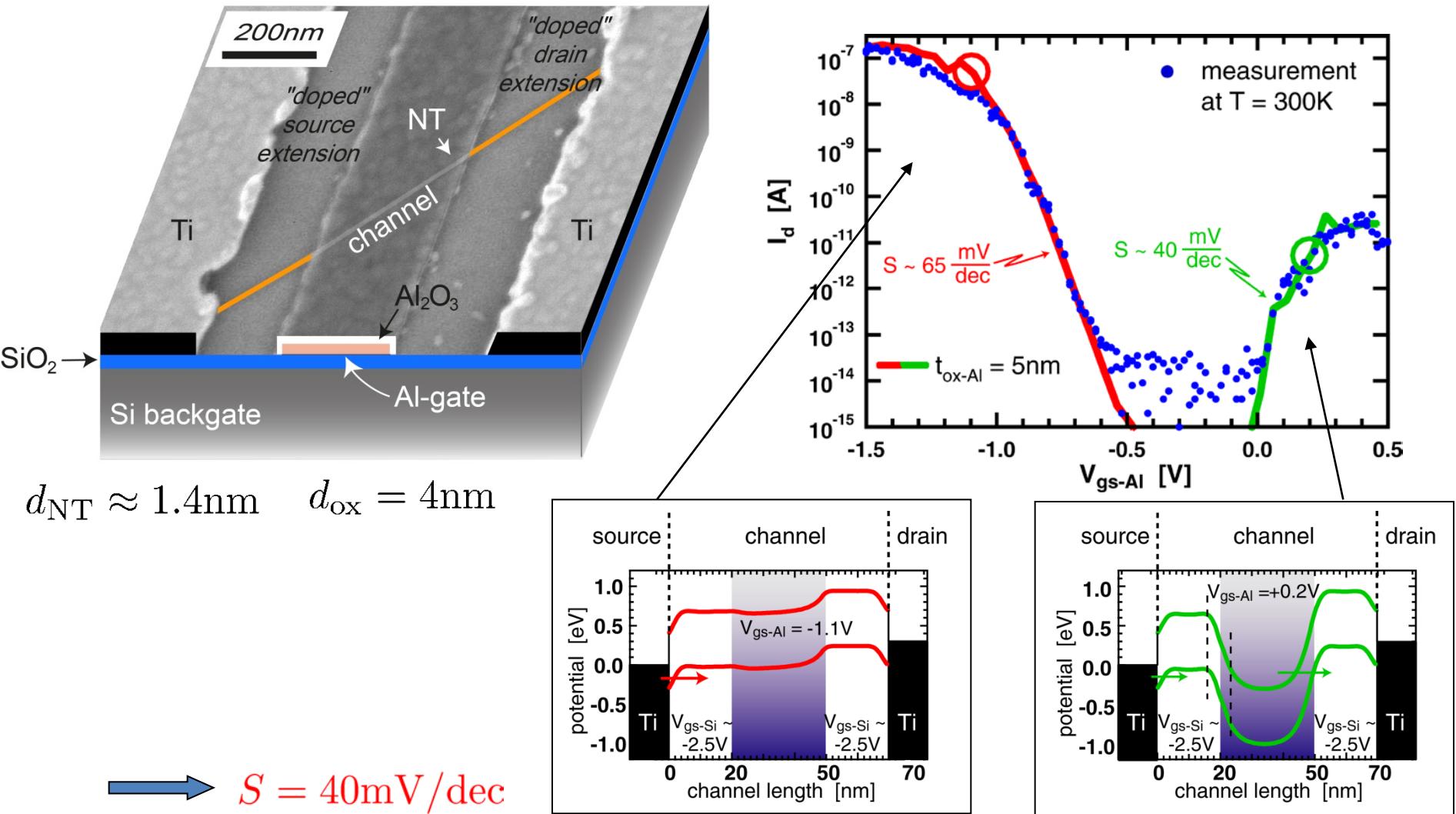


- device should be scalable to $L < 20\text{nm}$ without SCE
- increasing off-state leakage for short conventional CNFETs

J. Knoch, S. Mantl and J. Appenzeller,
Solid-State Electron., **49**, 73 (2005)



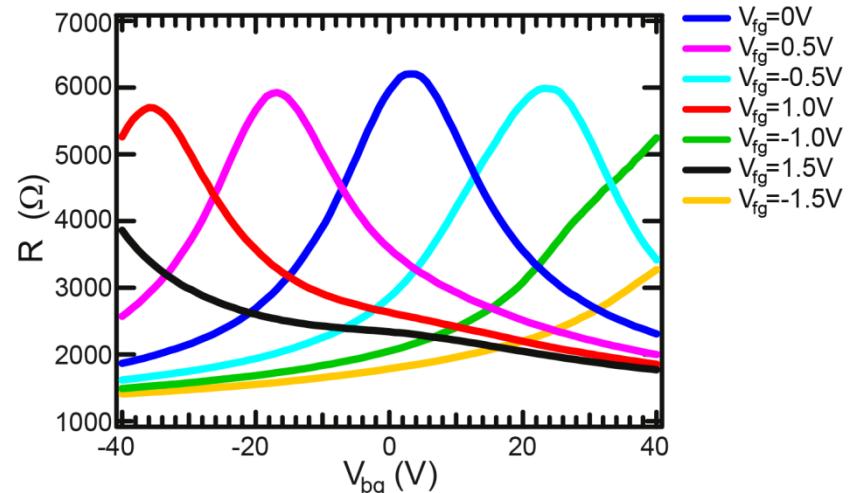
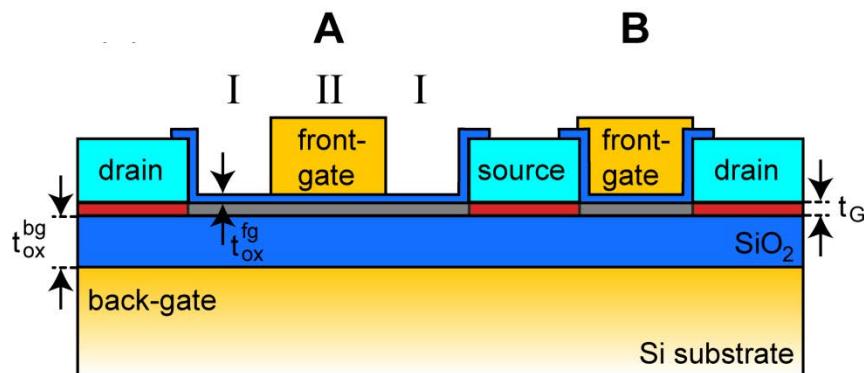
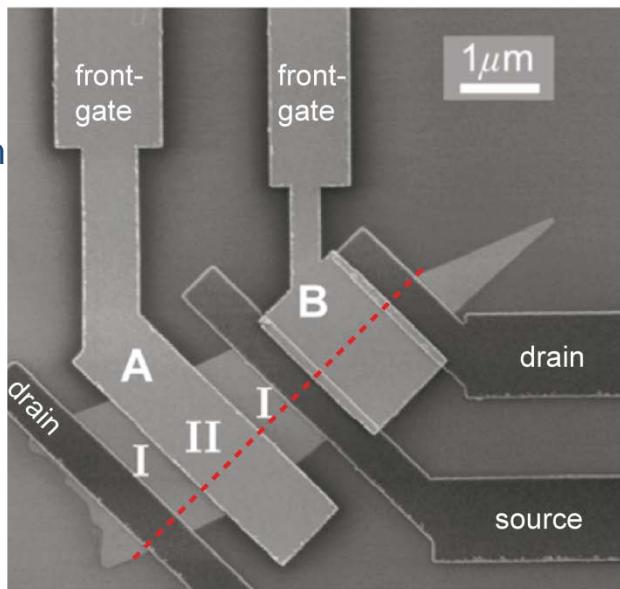
Tunnel-FET based on CNTs



J. Appenzeller, Y.-M. Lin, J. Knoch, Z. Chen and Ph. Avouris, IEEE Trans. Electron Dev., **52**, 2568 (2005).

Electrostatic Doping in Graphene

$L=1\text{mm}$
 $W=1\text{mm}$
 $t_{\text{ox},\text{bg}}=300\text{nm}$
 $t_{\text{ox},\text{fg}}=10\text{nm}$
Pd contacts



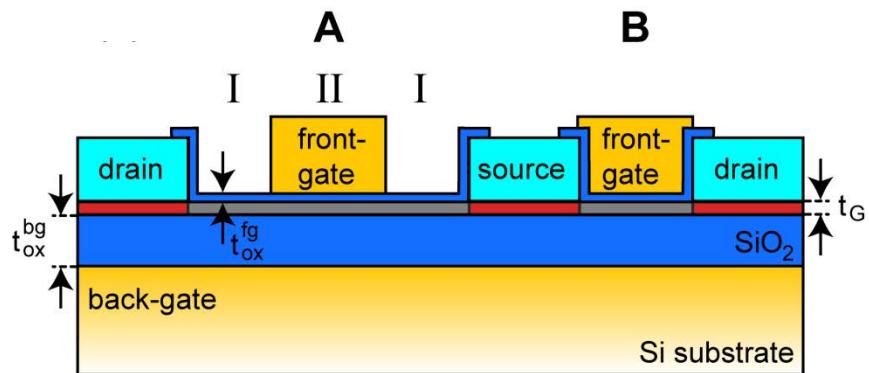
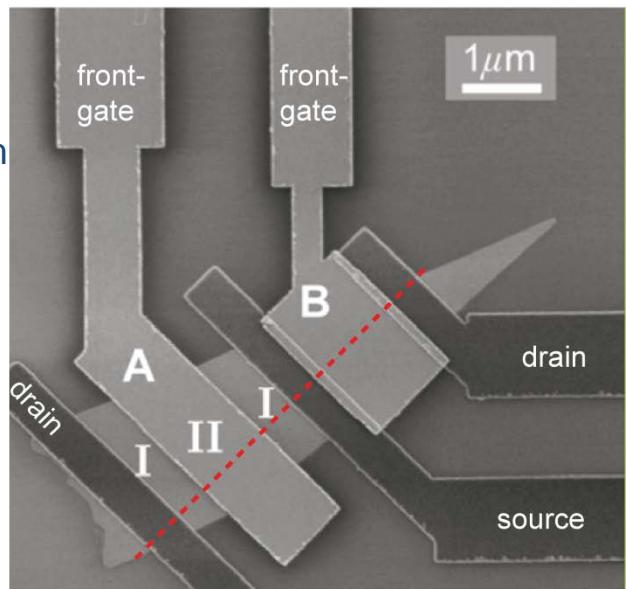
- **back-gate characteristics of device „B“ show double peak-structure**
- **second peak much less pronounced than main peak**

J. Knoch, Z. Chen and J. Appenzeller, IEEE Trans. Nanotechnol., 11, 513 (2012).

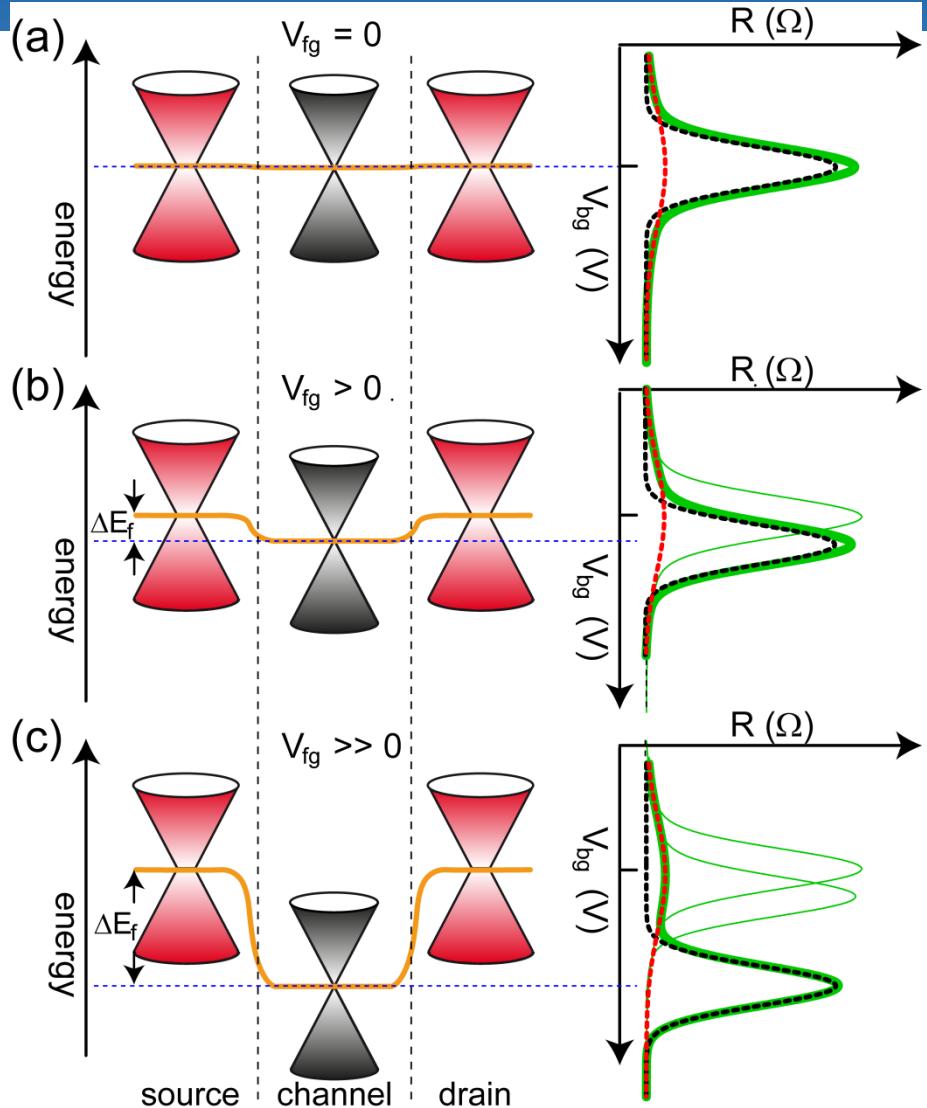
Electrostatic Doping in Graphene

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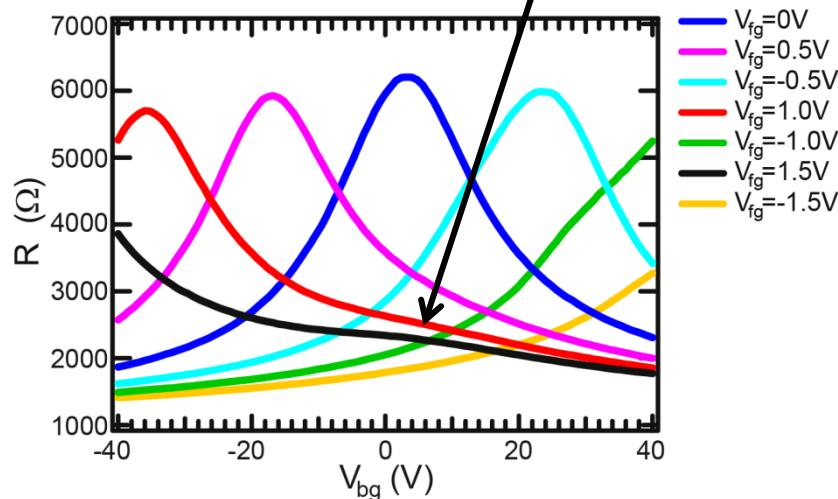
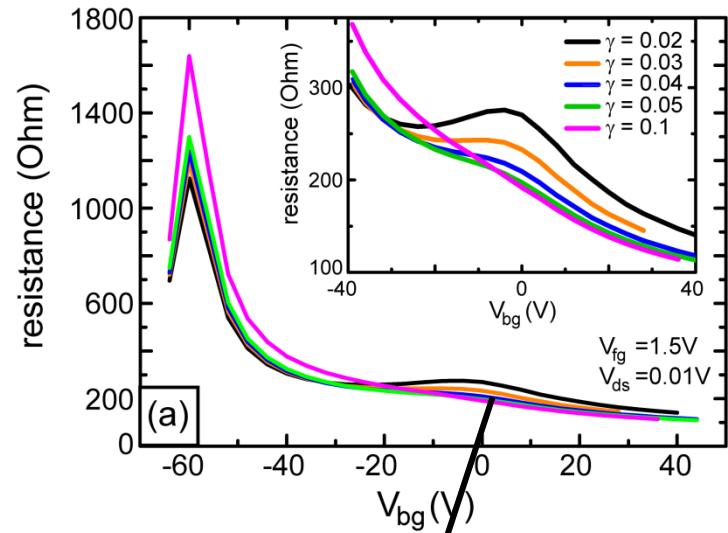
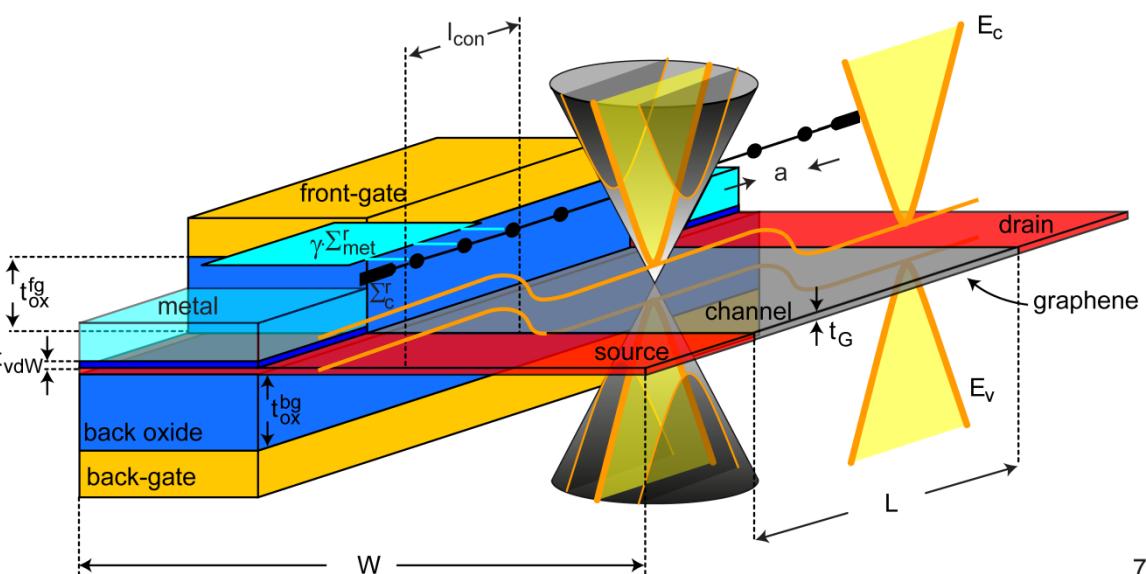


J. Knoch, Z. Chen and J. Appenzeller, IEEE Trans. Nanotechnol., 11, 513 (2012).



- second peak observable

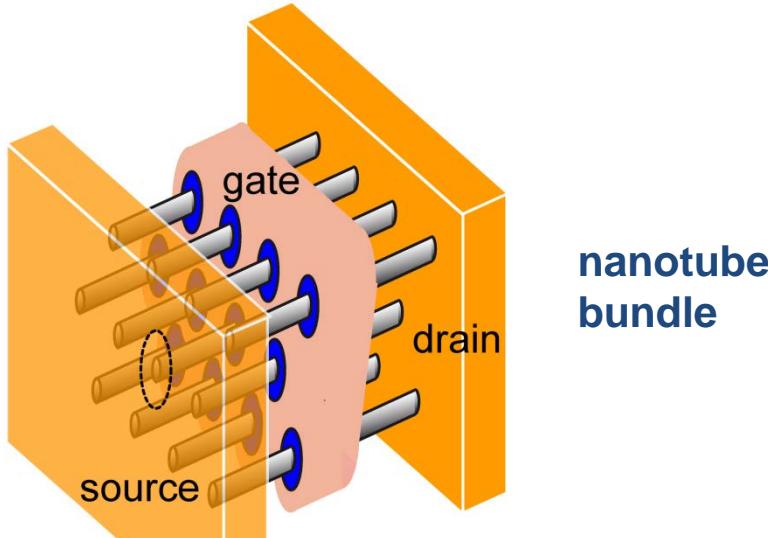
Electrostatic Doping in Graphene



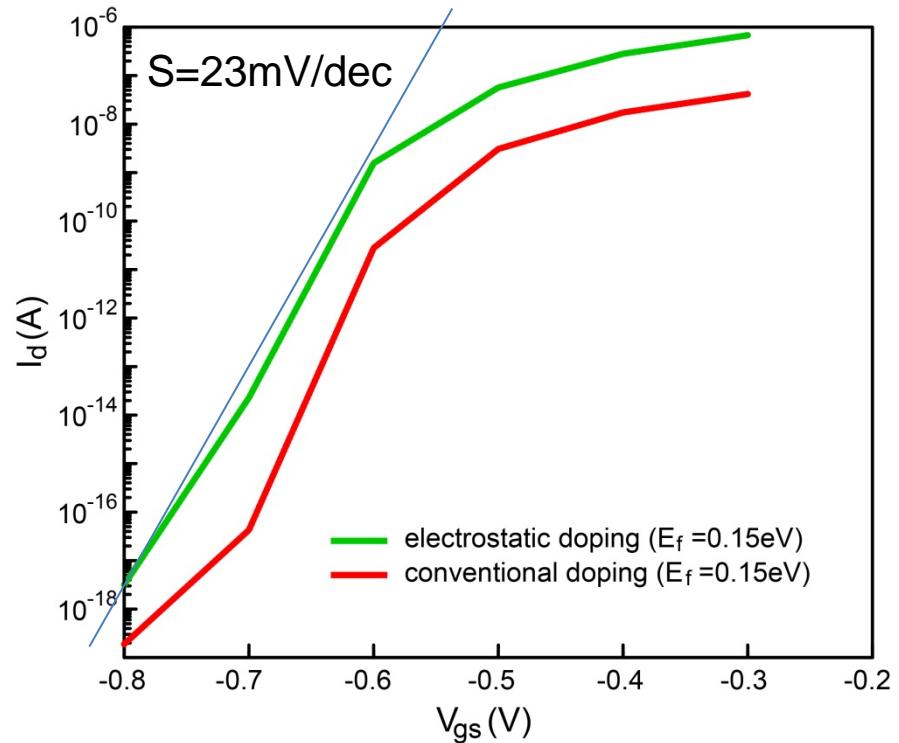
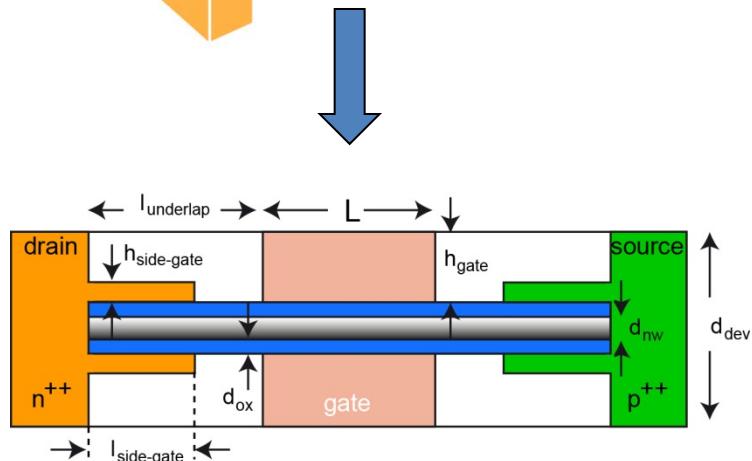
- back-gate voltage characteristics for different coupling strengths
- second peak unobservable if coupling too large

J. Knoch, Z. Chen and J. Appenzeller, IEEE Trans. Nanotechnol., 11, 513 (2012).

Electrostatic Doping in CNT Tunnel-FETs



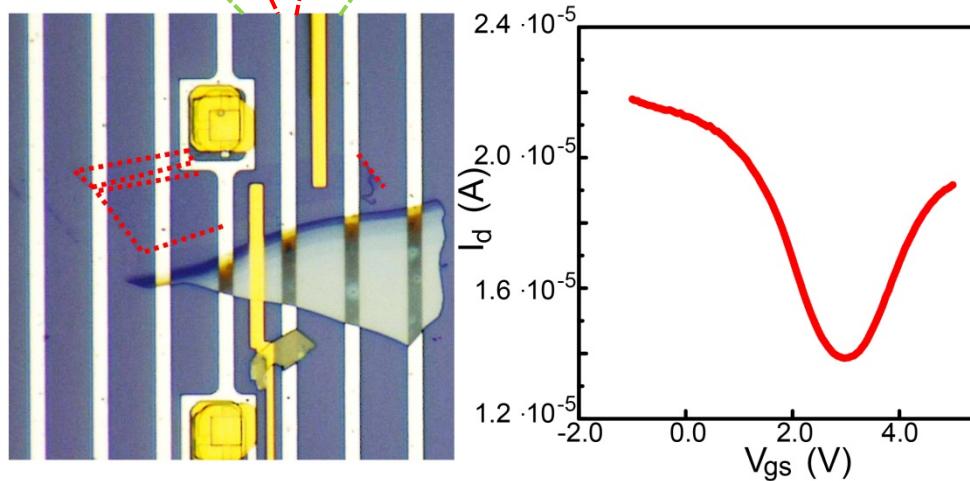
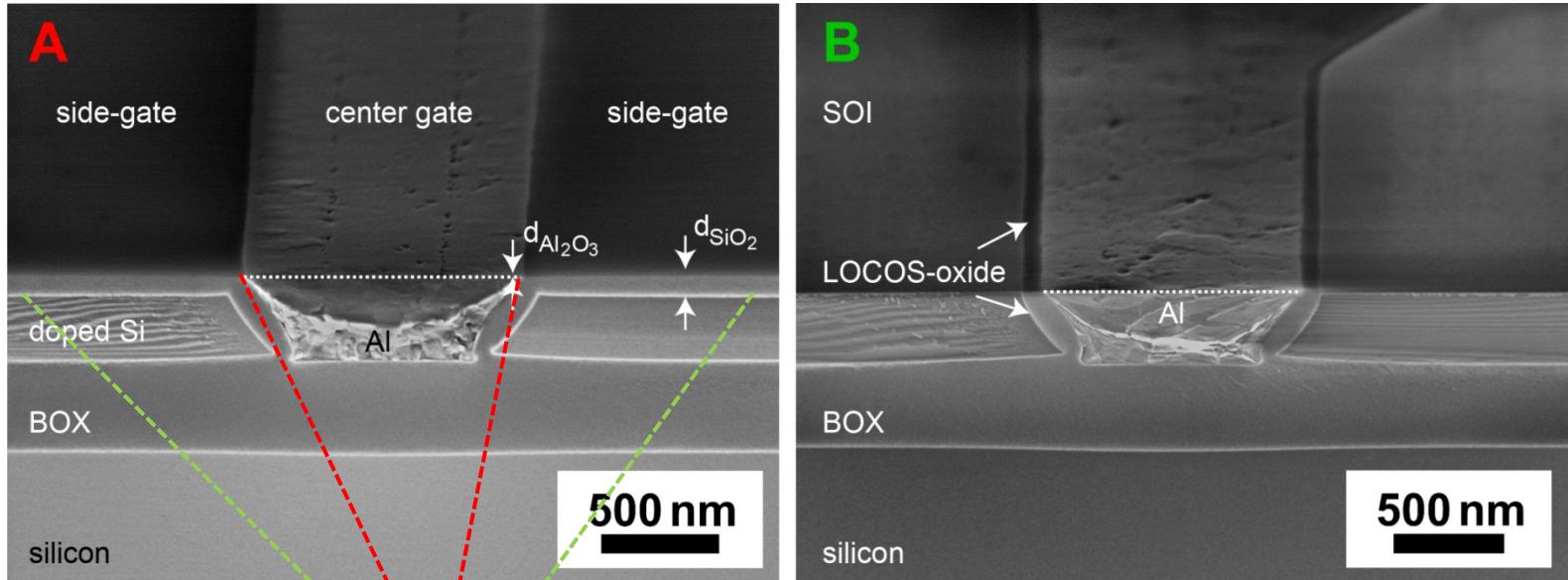
nanotube
bundle



- **significantly improved on-state due to improved screening**
- **steep inverse subthreshold slope of 23mV/dec over several orders of magnitude**

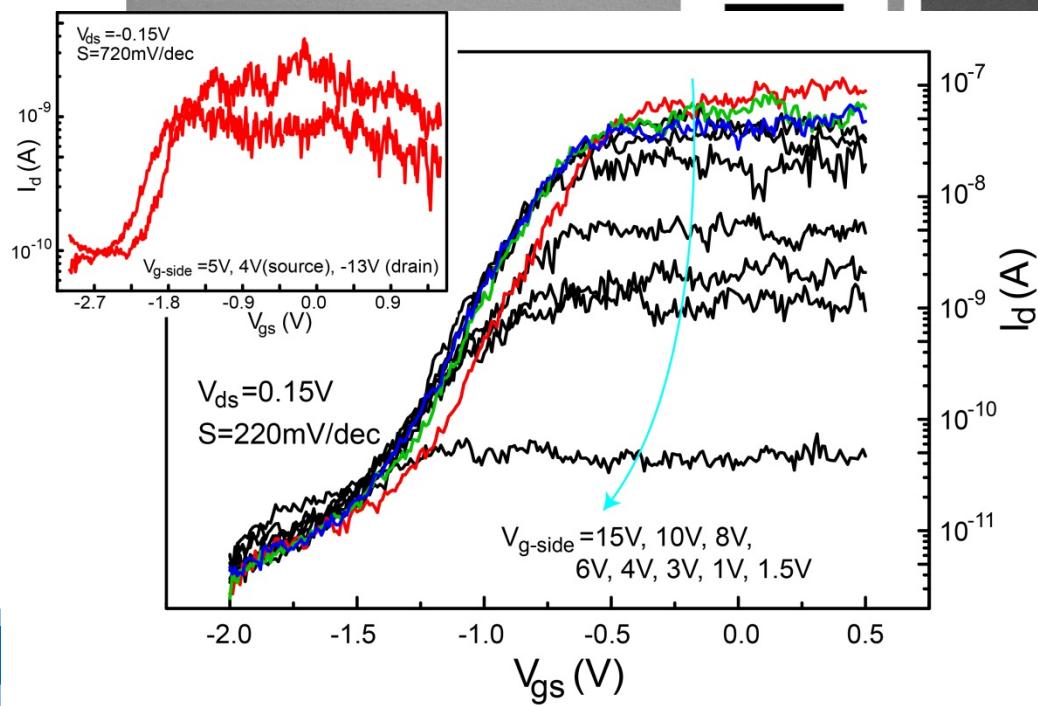
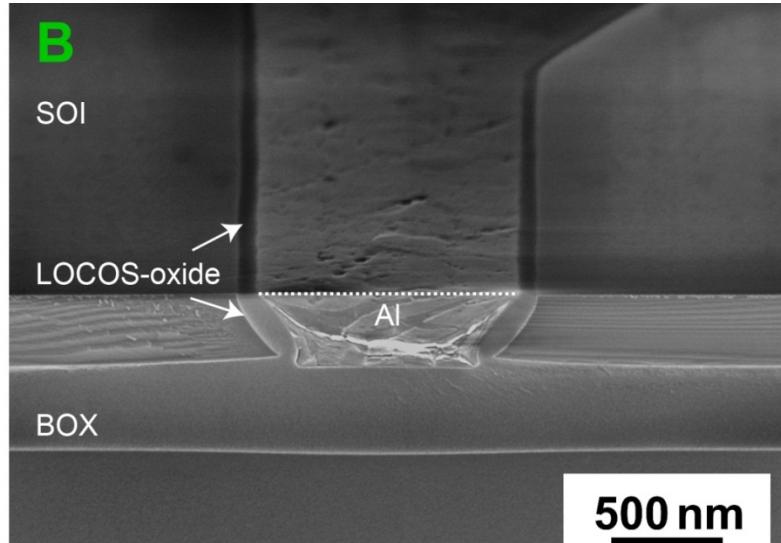
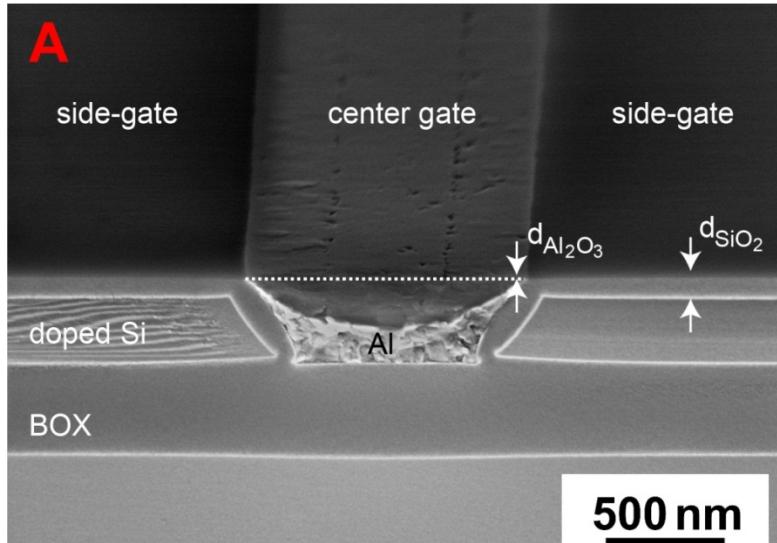
T. Grap and J. Knoch, submitted for publication

Electrostatic Doping

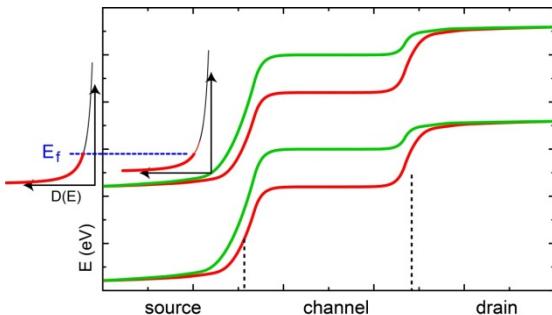


C. Kontis, M. Müller, K. Kallis and
J. Knoch, Appl. Optics, **51**, 385
(2012).

Electrostatic Doping

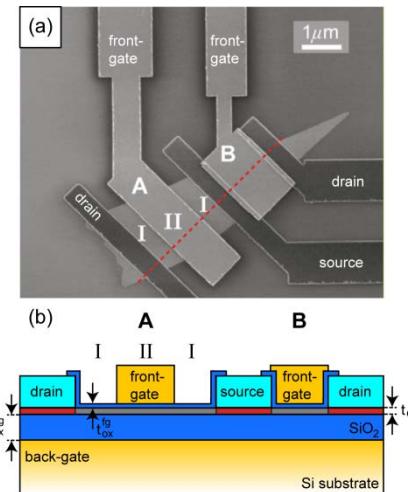
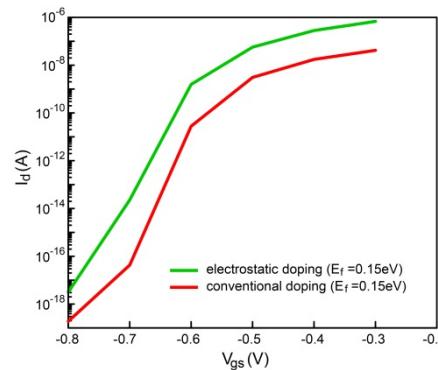
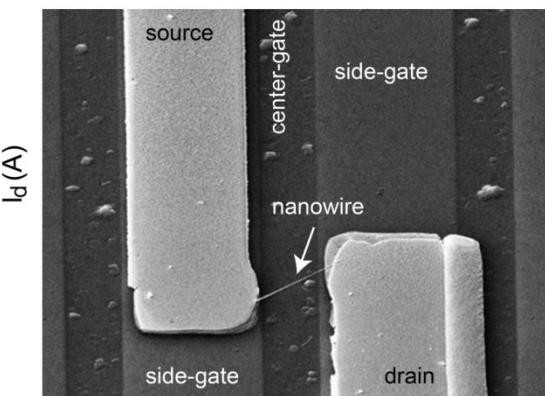


Conclusion



- low density of states detrimental for novel transistor concepts such as tunnel FETs due to insufficient screening

- electrostatic doping avoids issues related to doping *and* adds flexibility for general investigations of carbon based nanostructures



Acknowledgements

- M. Müller – RWTH Aachen and TU Dortmund
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