



HELMHOLTZ
ZENTRUM BERLIN
für Materialien und Energie

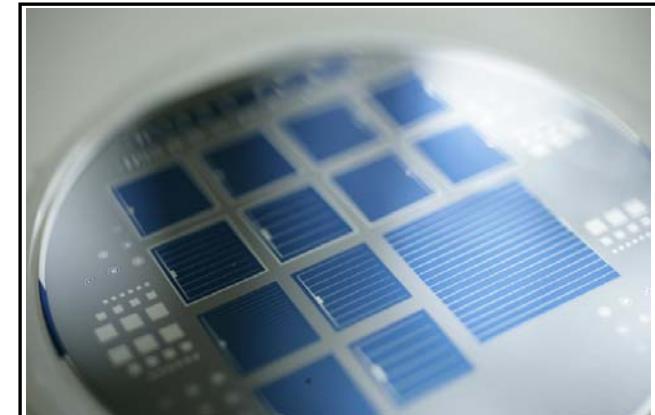
The a-Si:H/c-Si interface - key for high efficiency heterojunction cells

**Lars Korte, Erhard Conrad, Heike Angermann, Rolf Stangl, Tim Schulze,
Manfred Schmidt**

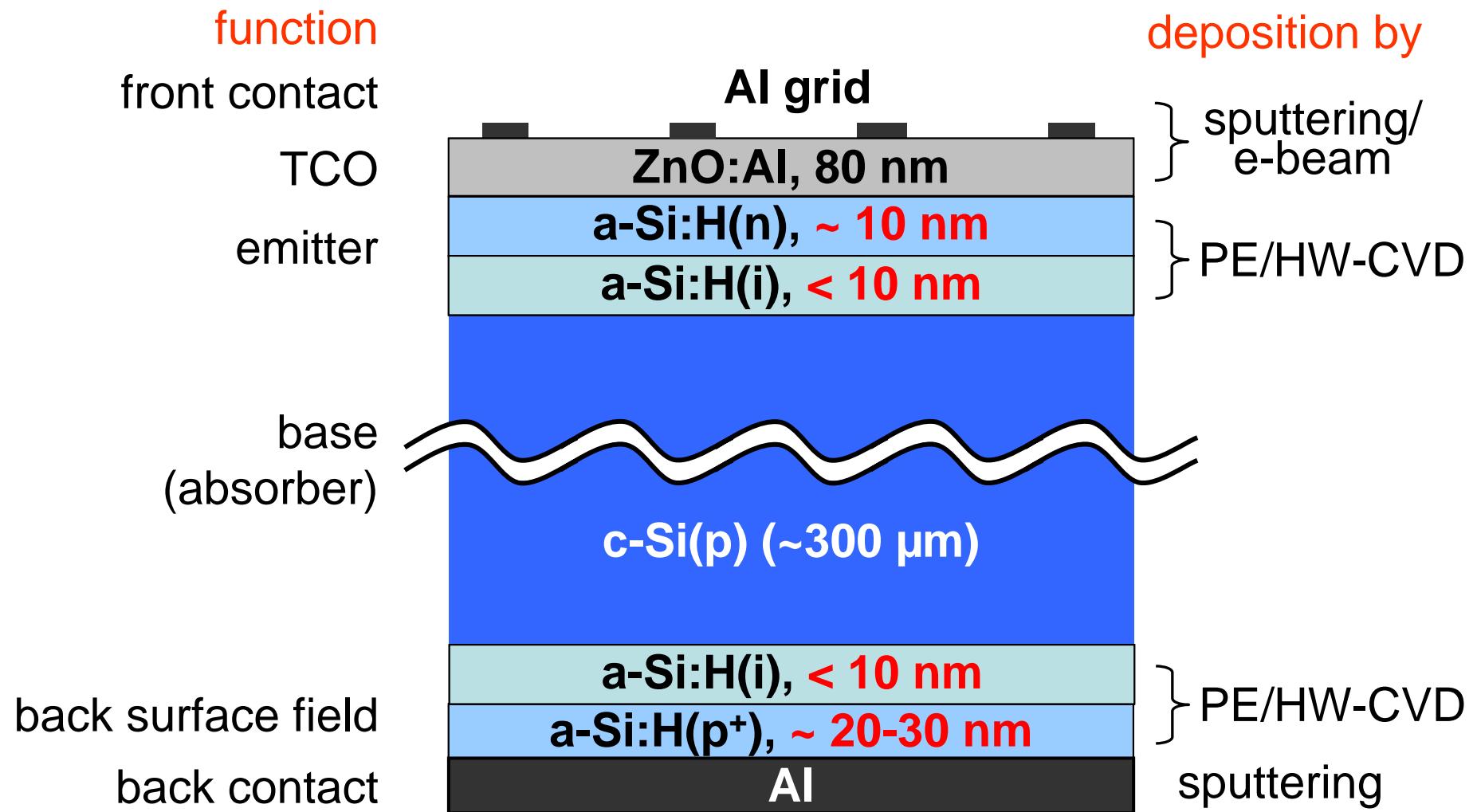
Weierstraß-Institut Berlin, 24. November 2008

Outline

- a-Si:H/c-Si heterojunction solar cells
- impact of the heterointerface: passivation and transport
- **experimental results:**
 - “soft deposition” of a-Si:H, initial growth on c-Si
 - optimization of a-Si:H:
 - optimum emitter thickness
 - measuring the density of states in <10 nm a-Si:H
 - optimum doping
 - a-Si:H/c-Si band offsets
- (older) cell results & summary

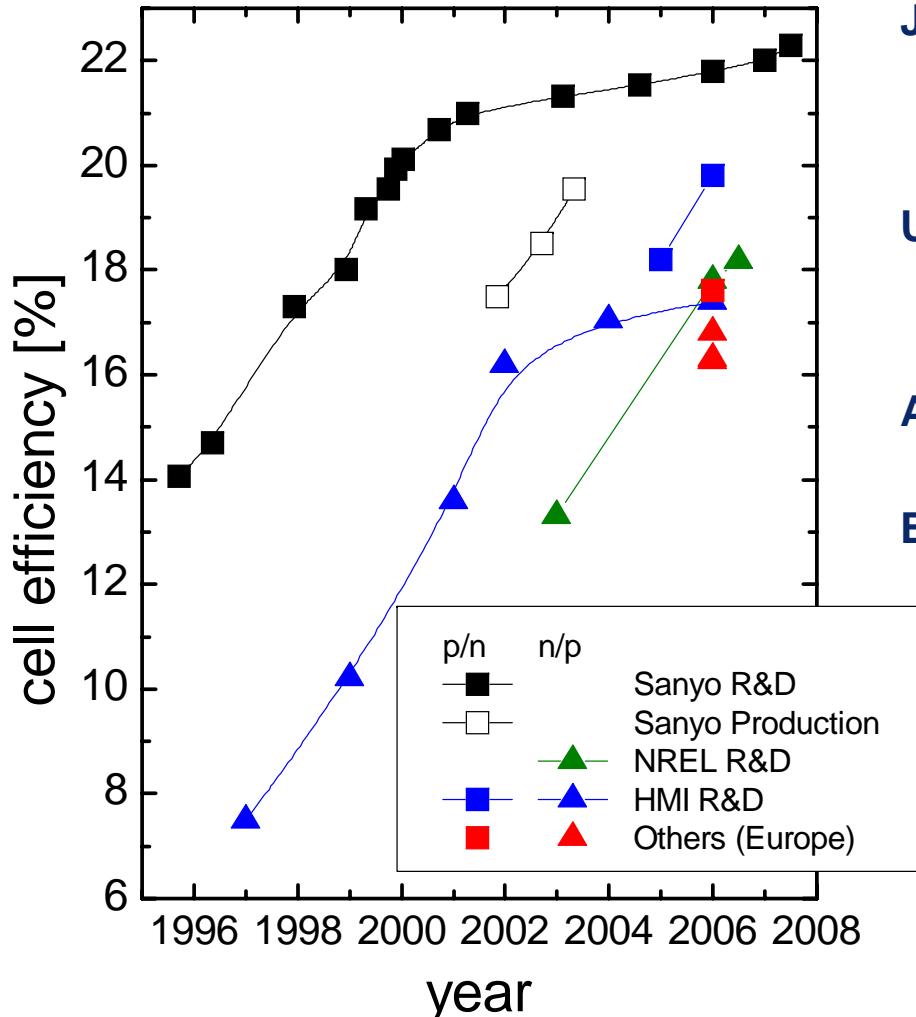


The a-Si:H/c-Si solar cell



Progress in a-Si:H/c-Si cell efficiency

(incomplete) list of R&D on a-Si:H/c-Si cells



Japan

Sanyo:

- (p,i)aSi on (n)cSi (22.3%)
- production: 100cm², 19.5%, 350MWp in '08

US

NREL: p/i on n-type, 18.2%, HW-CVD

UDEL: rear contacted cell, w/ SunPower

Australia

UNSW: p/i on n-type, 17.6% (EPVSEC 22)

Europe

HZB (D):

- n on p-type, 18.4%, and
- p on n-type, 19.8%, **no i-layer**

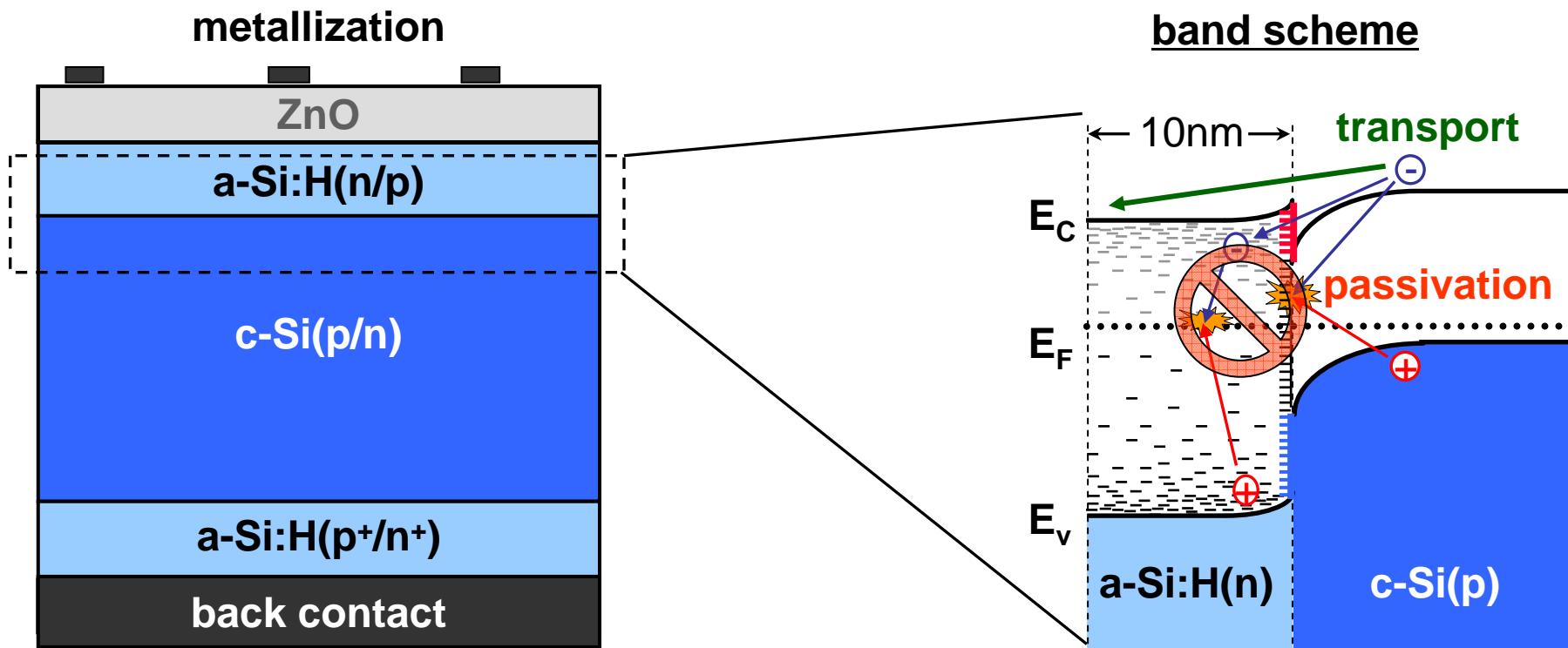
CEA-INES (F): n, (i)pmSi on p, 16.8%, 25cm²

U Neuchâtel (CH): n,i on p, VHF-PECVD

others:

CNRS (F), ENEA (I), IMEC (B),
U Utrecht (NL), IPE (D), FUH (D), ...

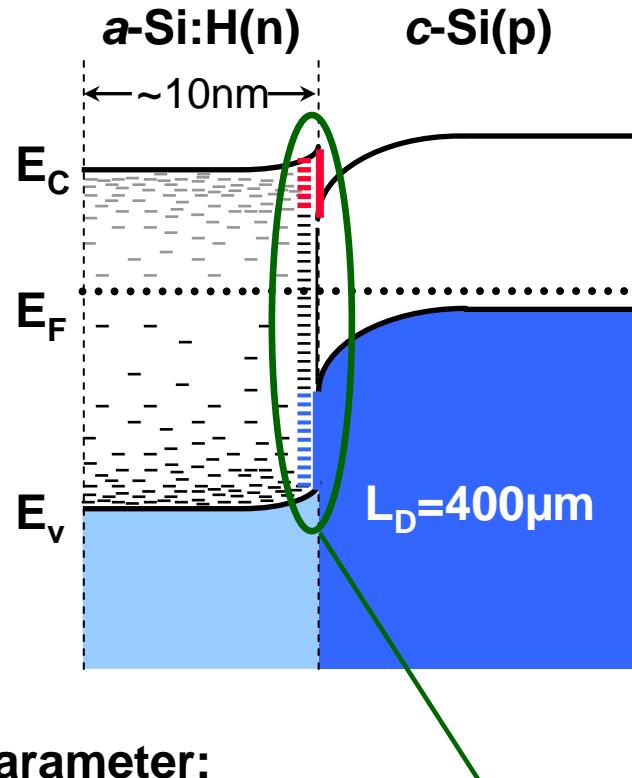
Impact of the heterointerface: Recombination and transport



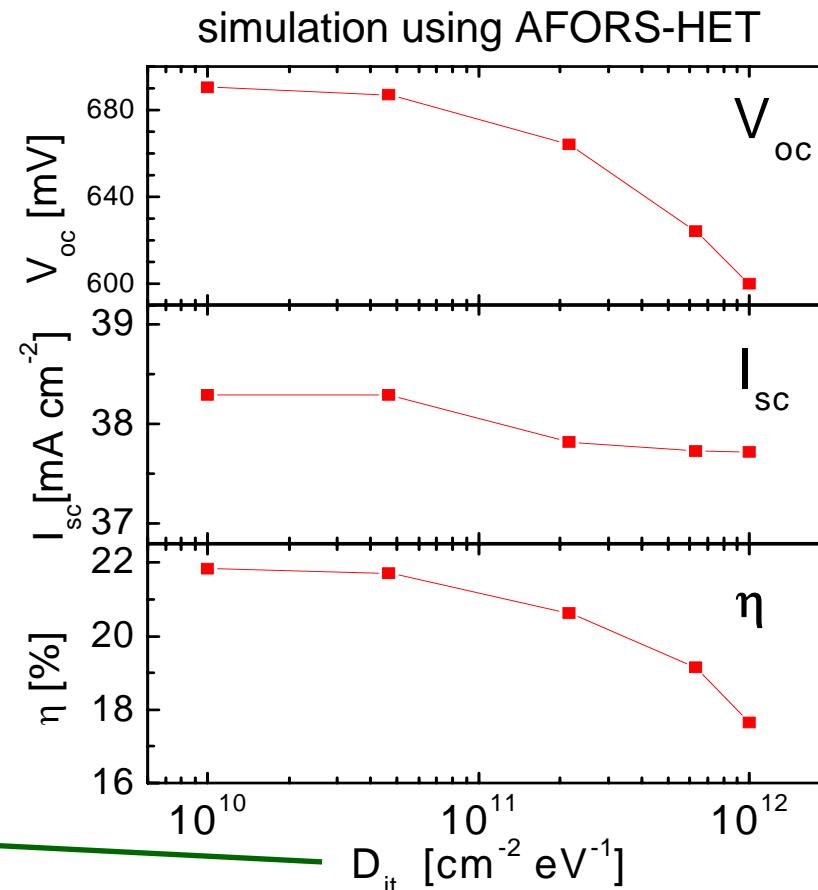
Tasks:

- minimize recombination losses at/near a-Si:H/c-Si interface
- maximize efficiency of charge carrier transport over heterointerface

influence of interface defects on cell parameters - simulation study



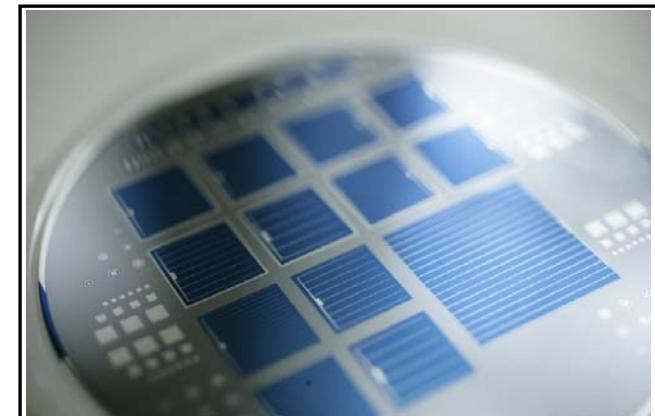
parameter:
interface state density $D_{it}(E)$



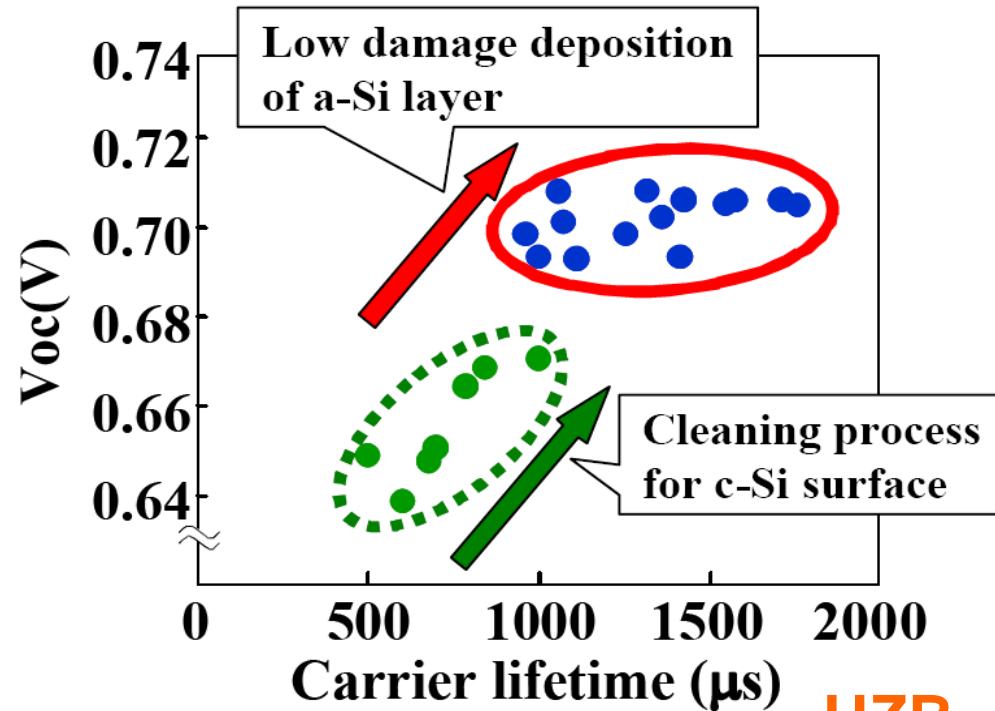
- strong influence of interface recombination at front (and rear) side
- $D_{it} < 10^{10} \text{cm}^{-2} \text{eV}^{-1} \rightarrow$ IF recombination plays no role in cell

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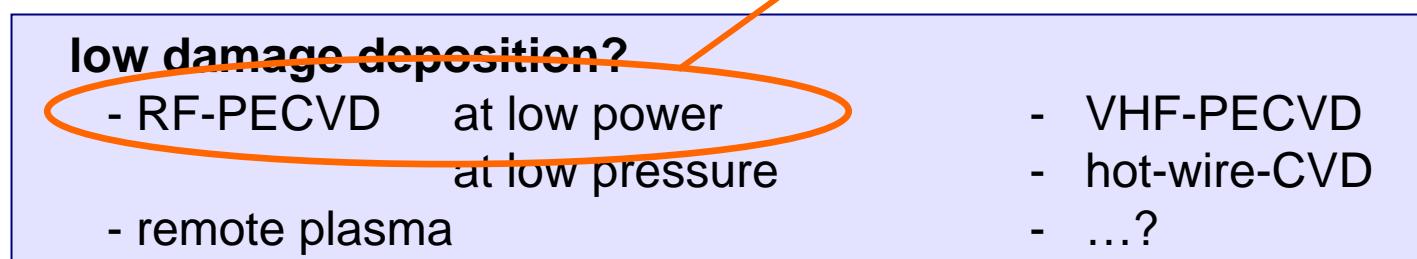


soft deposition of a-Si:H - the Sanyo „mantra“



Maruyama *et al.*, 4th IEEE WCPEC (2006)
and also

- Taira *et al.*, 22nd EPVSEC (2007),
- H. Kanno *et al.*, 23rd EPVSEC (2008),
- ...



Influence of epitaxy at the a-Si:H/c-Si interface

detrimental effect of epi-growth at the a-Si:H/c-Si interface

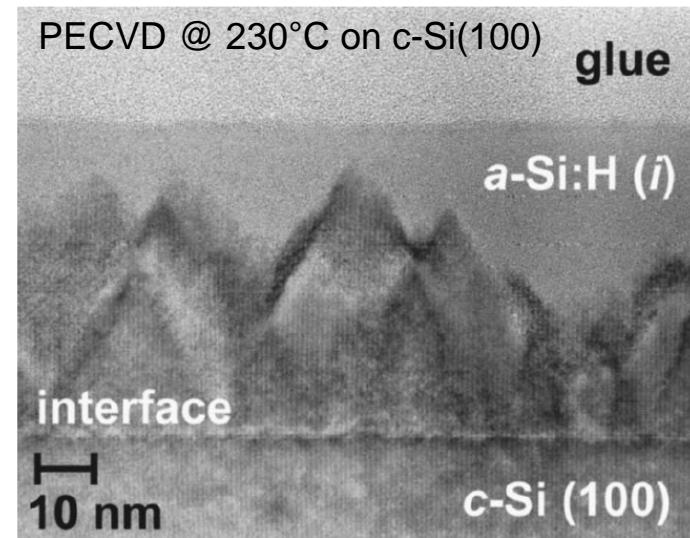
reported many times:

- E. Centurioni *et al.*, IEEE Trans. El. Devices & 19th EPVSEC, Paris (2004) 1285
- T.H. Wang *et al.*, ibd., p. 1269
- de Wolf & Kondo, Appl. Phys. Lett. 90 (2007) 042111
- ...

usual result: epitaxy up to doped layer $\rightarrow V_{oc} < 600\text{mV}$

possible explanations:

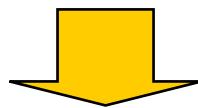
- epitaxy extends through i-layer
 \rightarrow surface defect passivation by *doped* a-Si:H:
much less effective than (i)a-Si:H
- partial epitaxy/mixed phase (i)a-Si:H
 \rightarrow increased interface area betw. a-Si & c-Si
- poor conditions for epitaxy
 \rightarrow epi highly defective



de Wolf & Kondo, Appl. Phys. Lett. 90 (2007) 042111

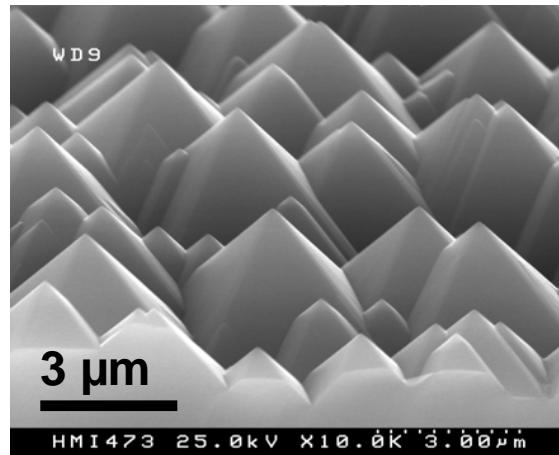
Smoothing and passivation of pyramids

additional
complication:
random pyramid
surface texture
(light trapping)



need optimized
chemical pre-
treatment of c-Si
wafer

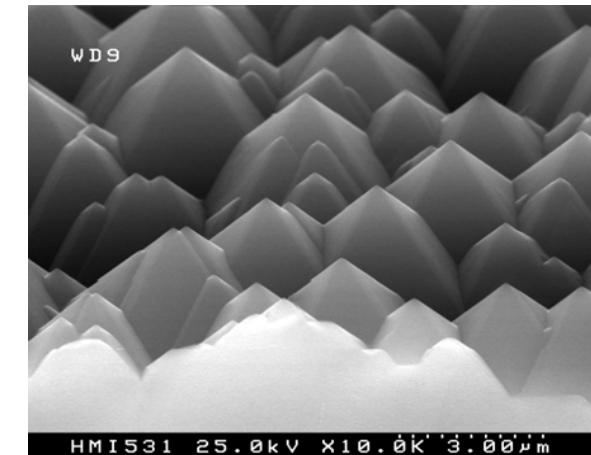
anisotropic etching



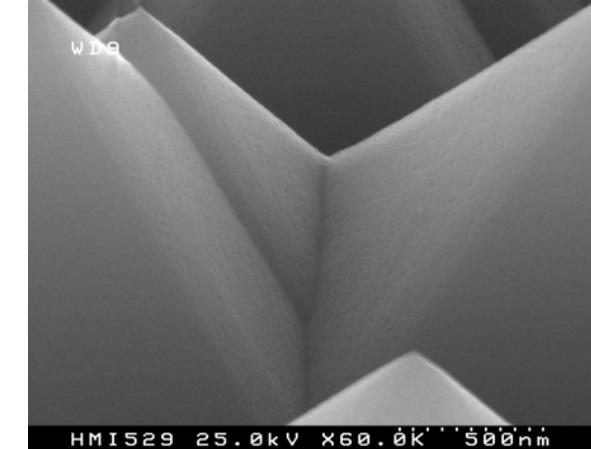
0.5 μm

HMI475 25.0kV x60.0K 500nm

optimized smoothing



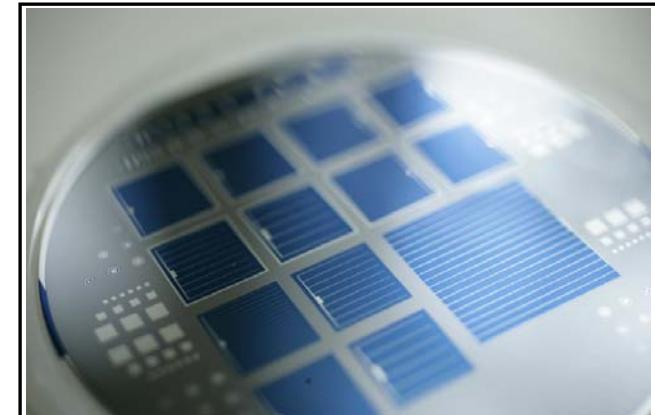
HMI531 25.0kV x10.0K 3.00μm



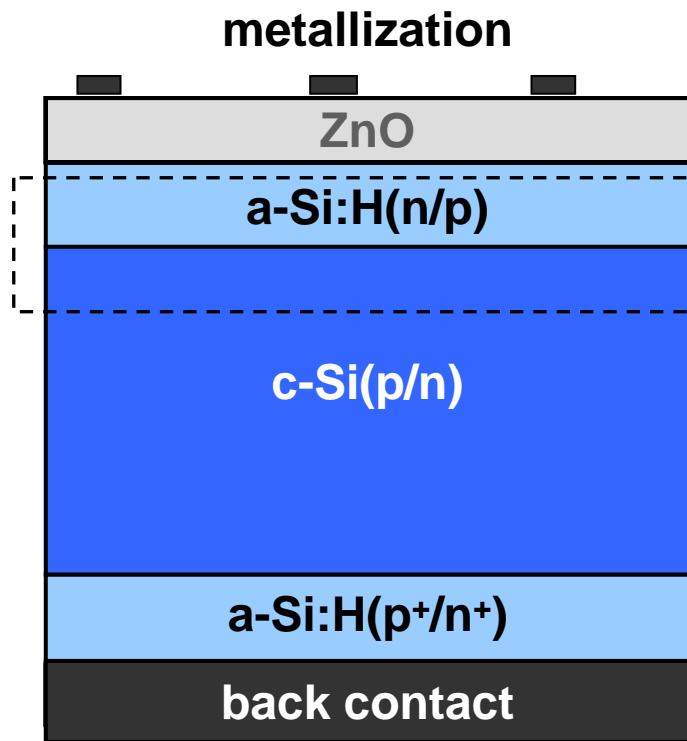
HMI529 25.0kV x60.0K 500nm

Outline

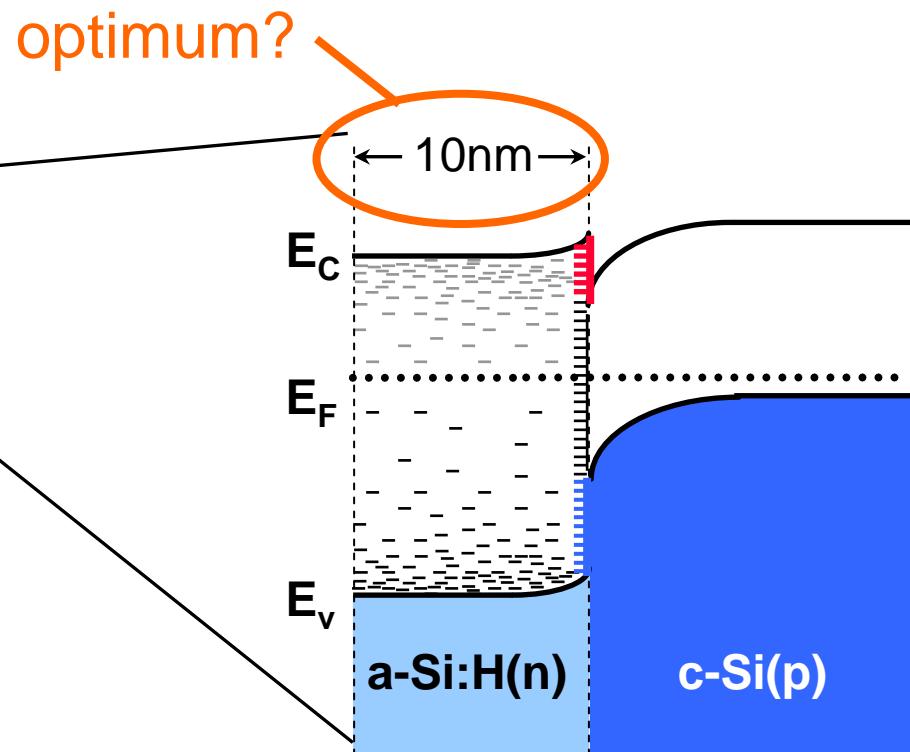
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cell structure

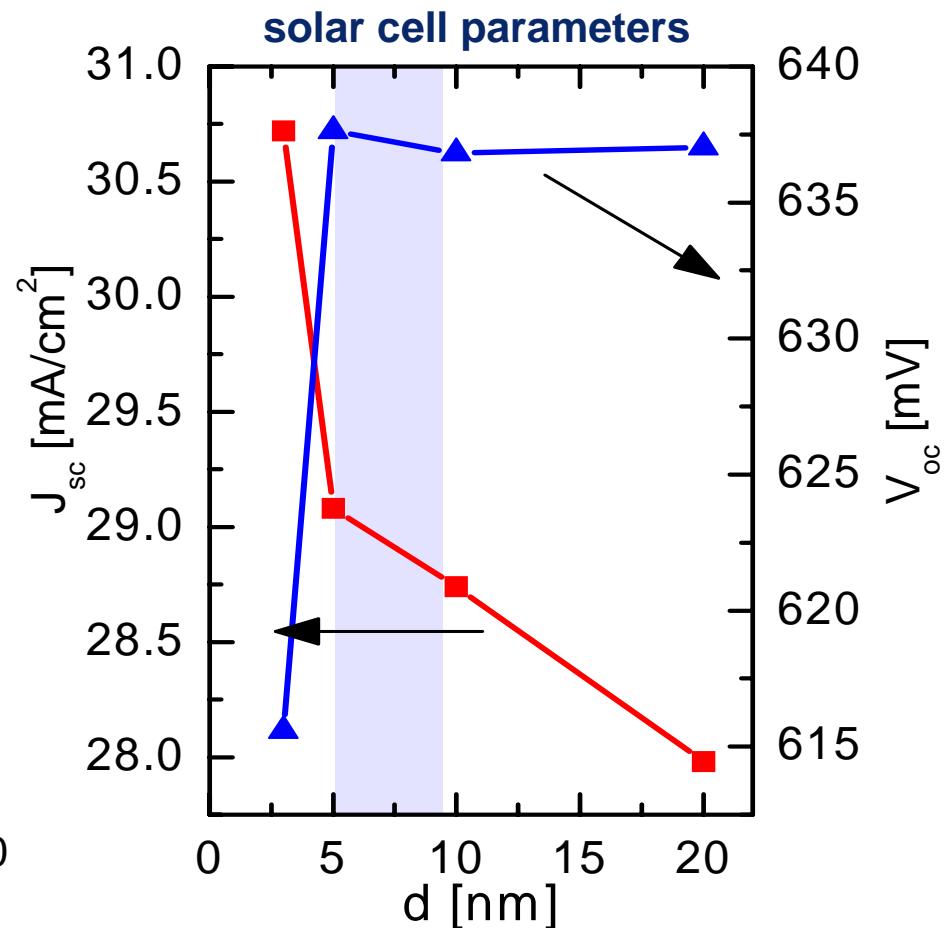
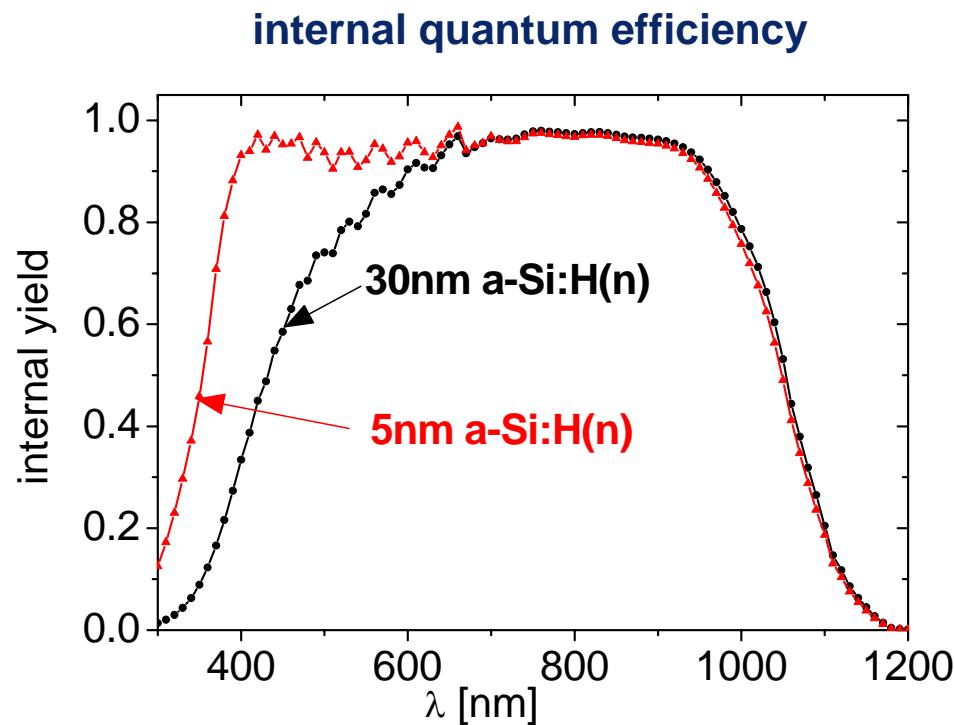


band diagram



finding the optimal a-Si:H emitter thickness

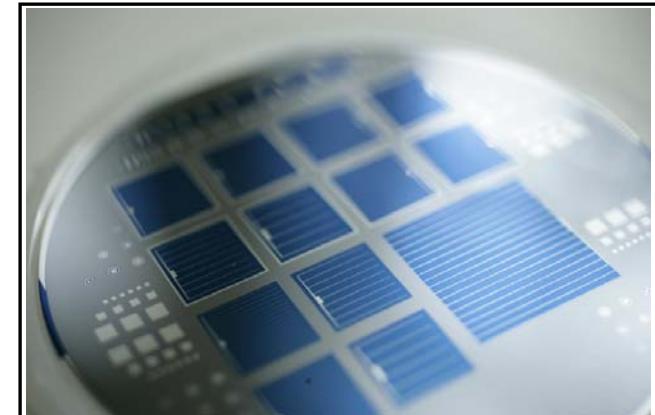
cell structure: TCO/(n)a-Si:H/(p)c-Si/Al



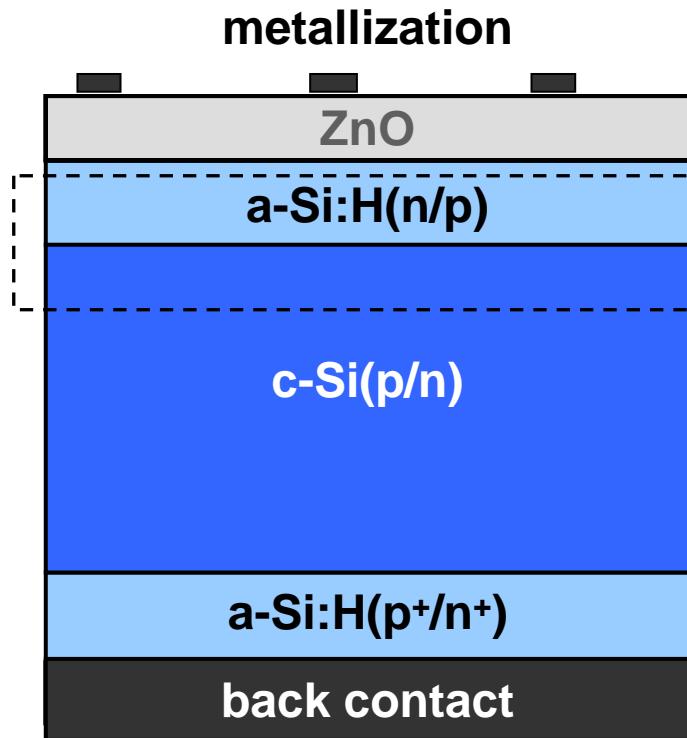
absorption and recombination in the a-Si:H emitter lead to a decrease in J_{SC}
➤ optimum a-Si:H thickness: 5-10nm – how to analyze its electronic properties?

Outline

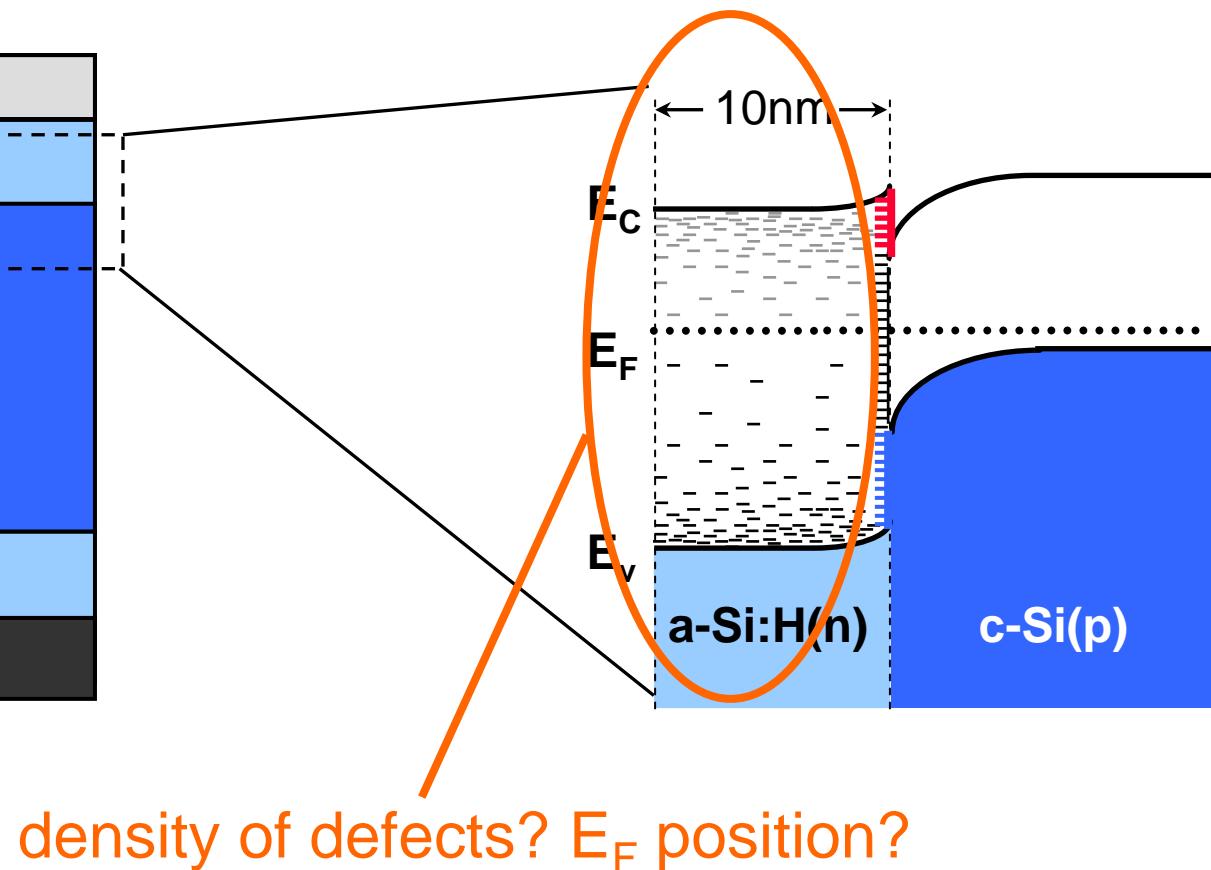
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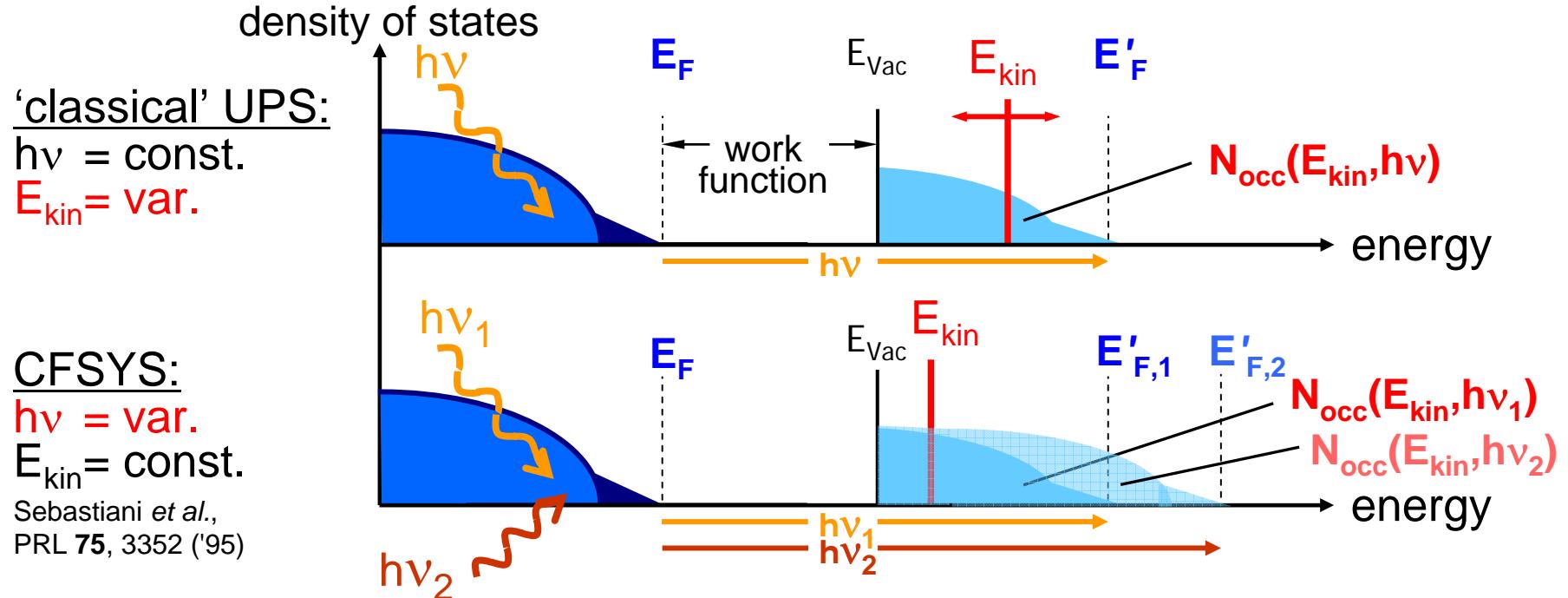
cell structure



band diagram



Constant Final State Yield Photoelectron Spectroscopy



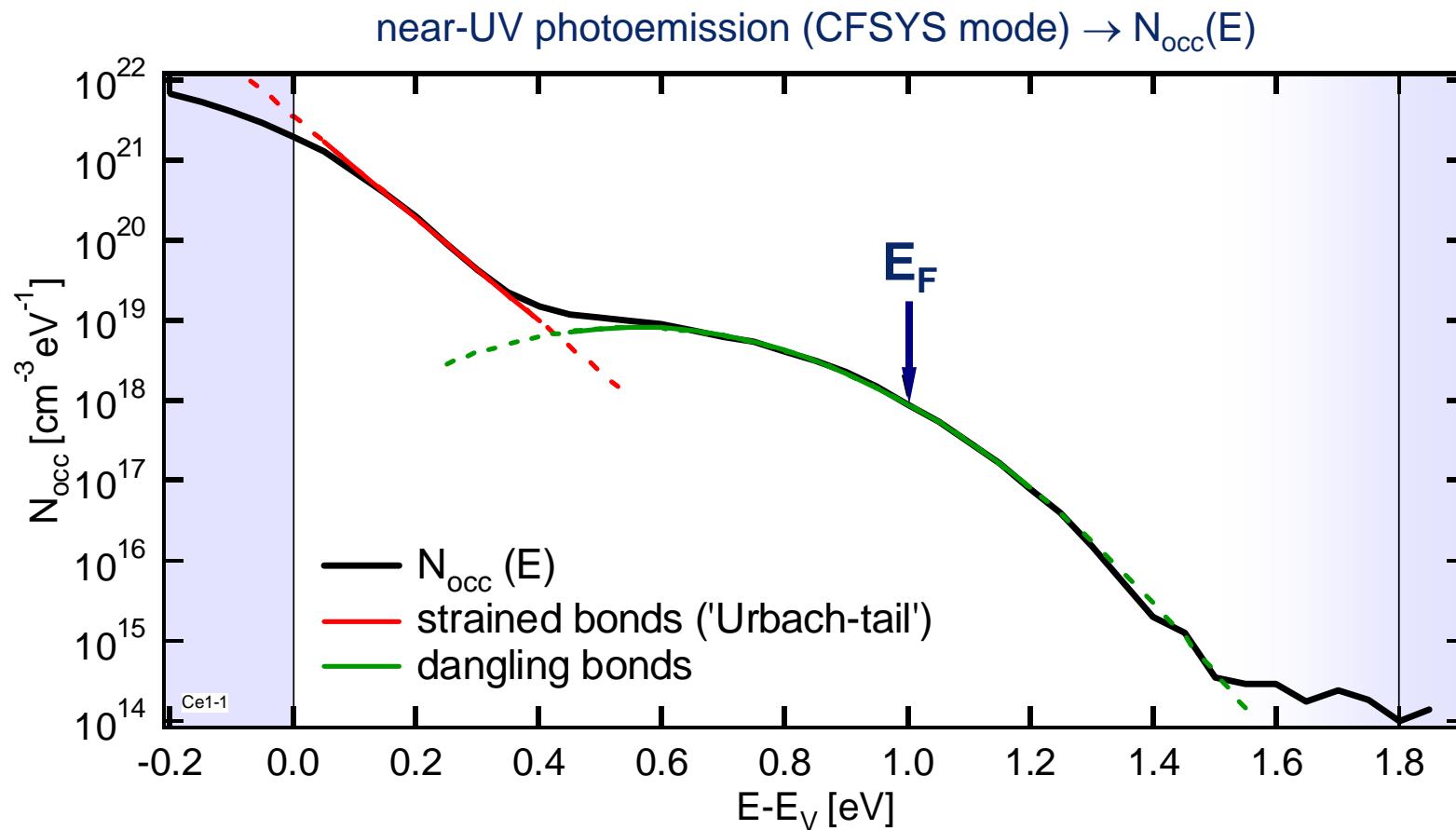
low excitation energy (4...7 eV)

- high electron escape depth (6-10 nm)
- high excitation cross section (at 4eV: 10³ times higher than at 21.2eV!)

we obtain: photoelectron yield

- density of occupied states $N_{occ}(E)$
- Fermi level position E_F

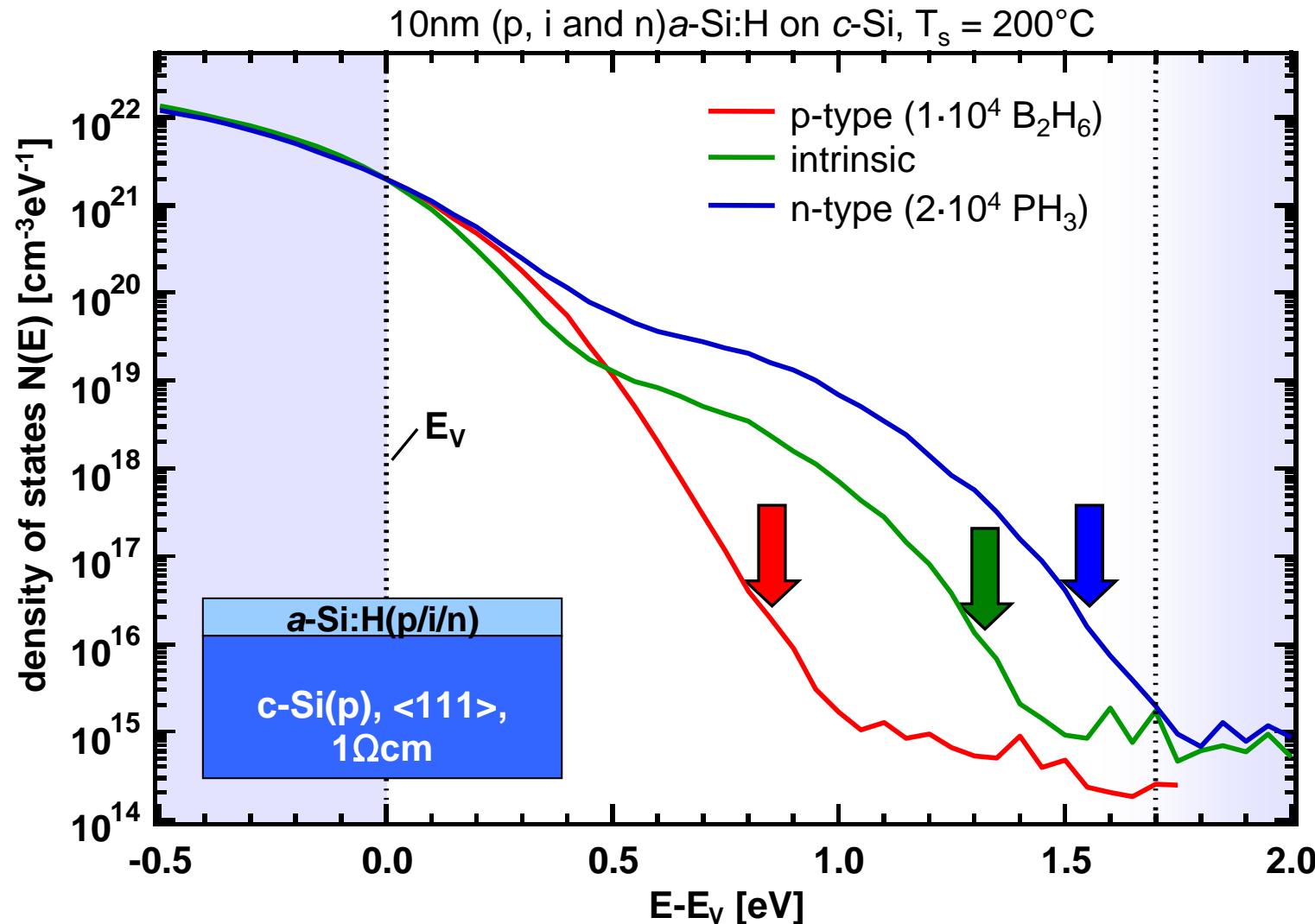
density of a-Si:H gap states



$N_{\text{occ}}(E) = \text{exp. tail ('strained bonds')} + \text{Gaussian distribution ('dangling bonds')}$

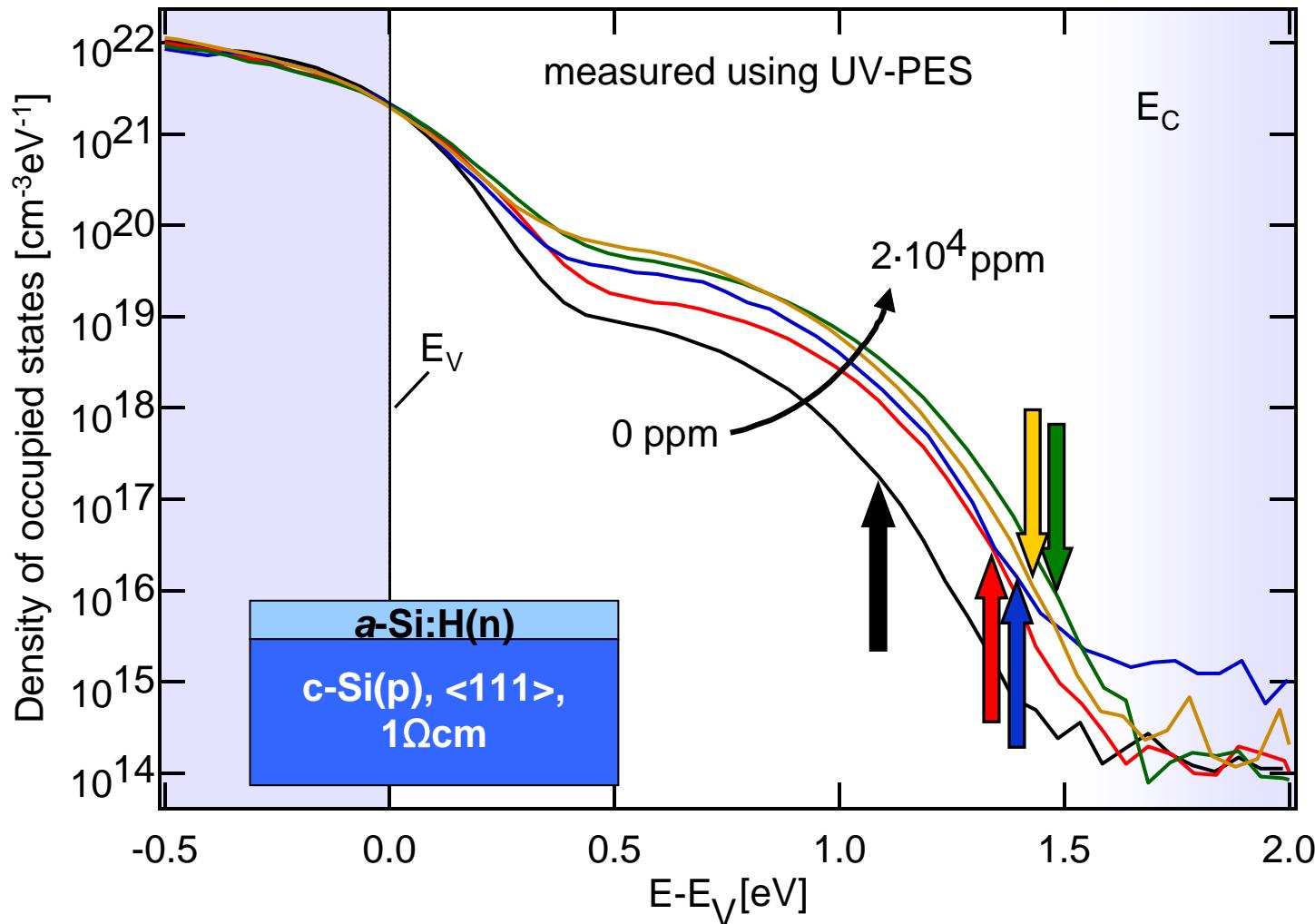
Urbach energy E_{0v} integrated density of deep defects N_D

doping dependence of $N(E)$, E_F



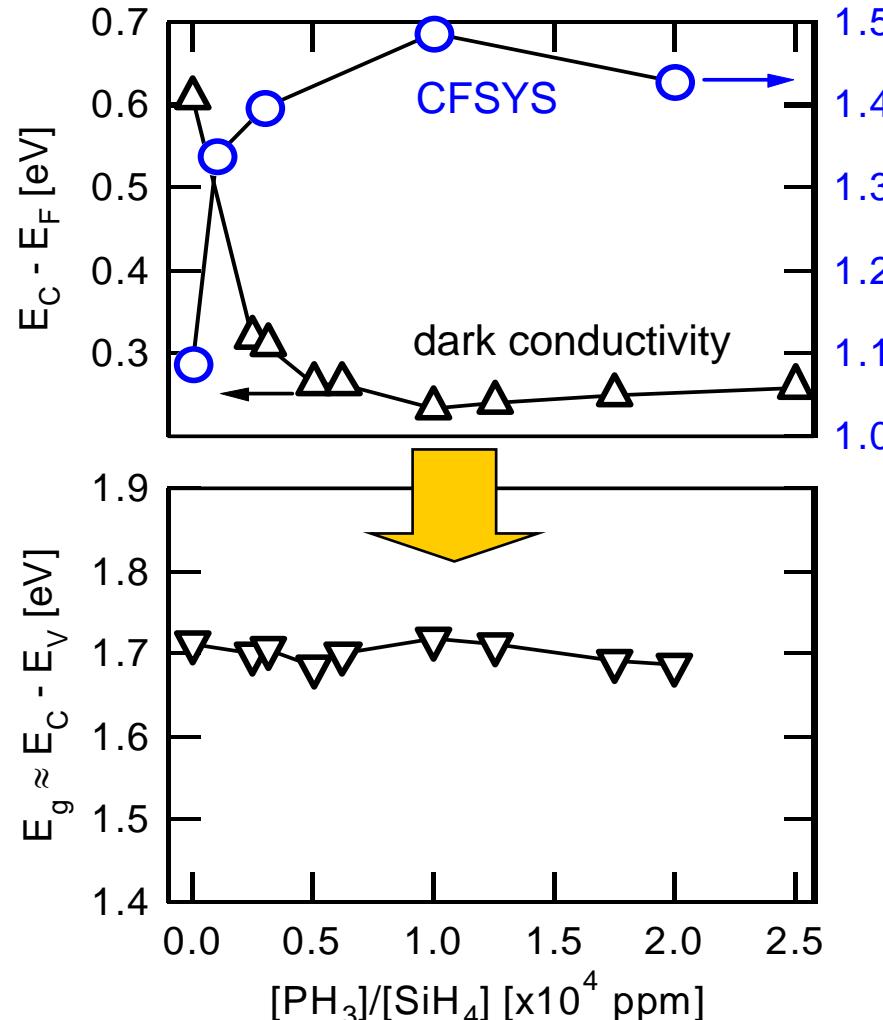
n-type doping series

samples: ~10 nm a-Si:H(n) on c-Si(p), $[\text{PH}_3]/[\text{SiH}_4] = 0 - 2 \times 10^4$ ppm, $T_s = 170^\circ\text{C}$



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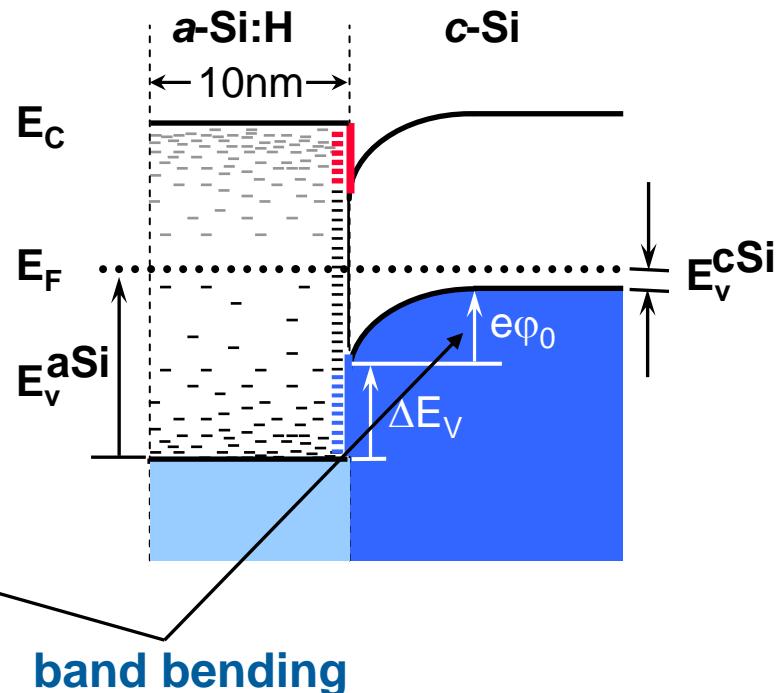
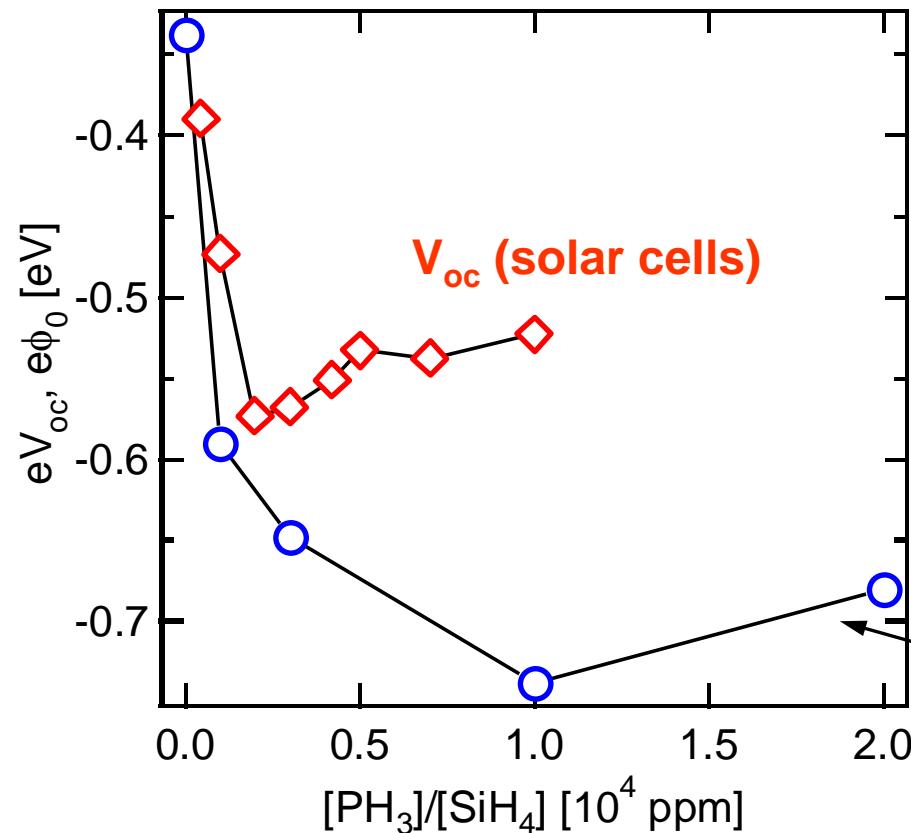
optimum doping:
 $[\text{PH}_3]/[\text{SiH}_4] = 10^4 \text{ ppm}?$

no change in band gap
with doping ($E_g \sim 1.7 \text{ eV}$)

same as for thick films
(e.g. V. Chacorn, D. Haneman,
Solid State Comm. 65 (1988) 609)

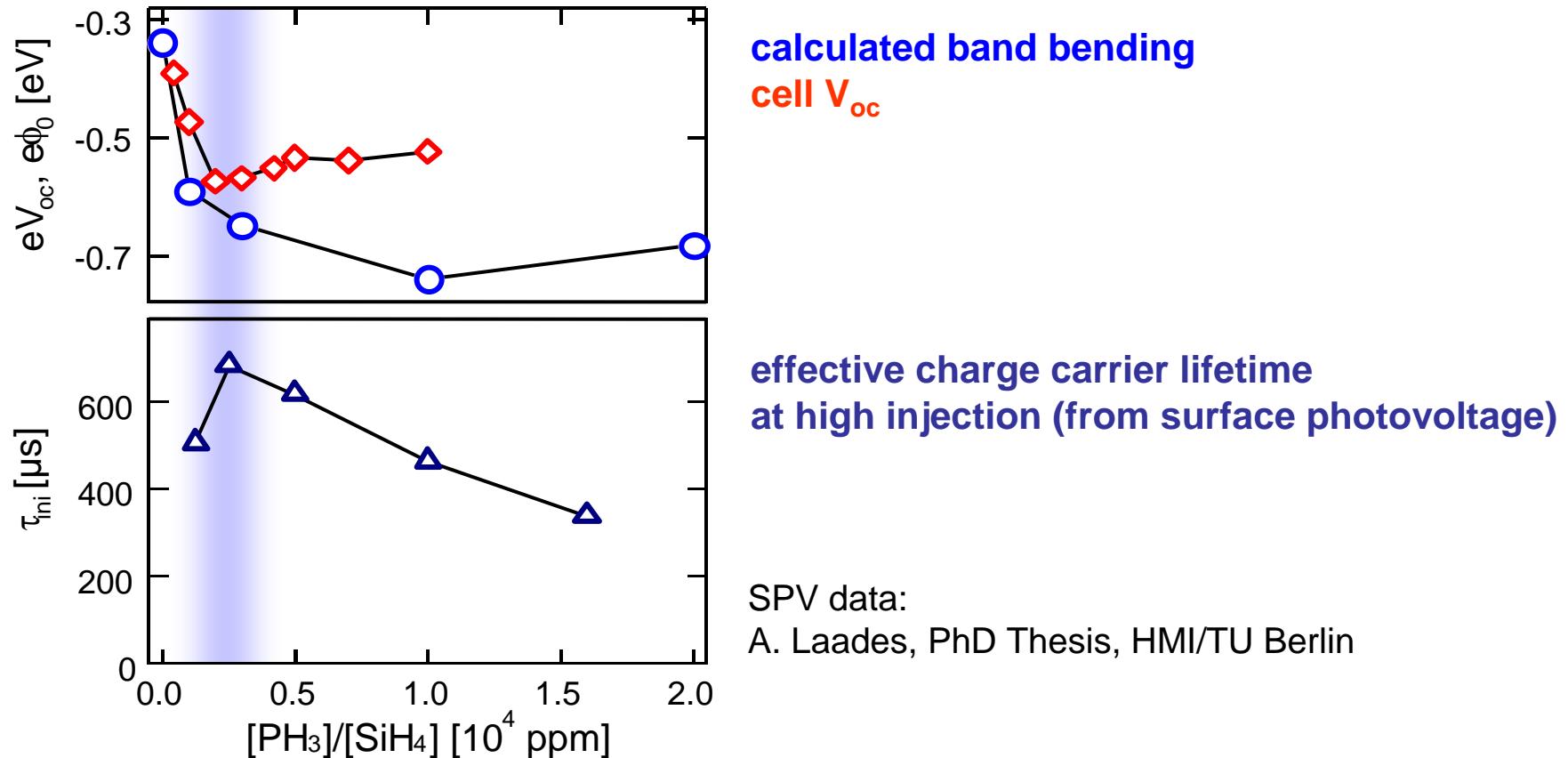
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→ optimum doping ~ 2000 ppm - higher doping: enhanced recombination?

Band bending, V_{oc} and recombination - optimized a-Si:H doping



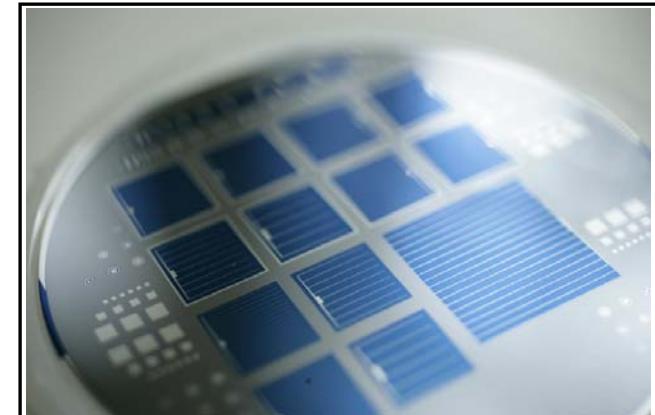
doping > 2000ppm: enhanced recombination

- reduced charge carrier conc.
- V_{oc} decreases

and also: hopping into a-Si:H via a-Si:H tail states possible, additional recombination
(see also: Boehme *et al.*, J. Non-Cryst. Sol. 352 (2006) 1113)

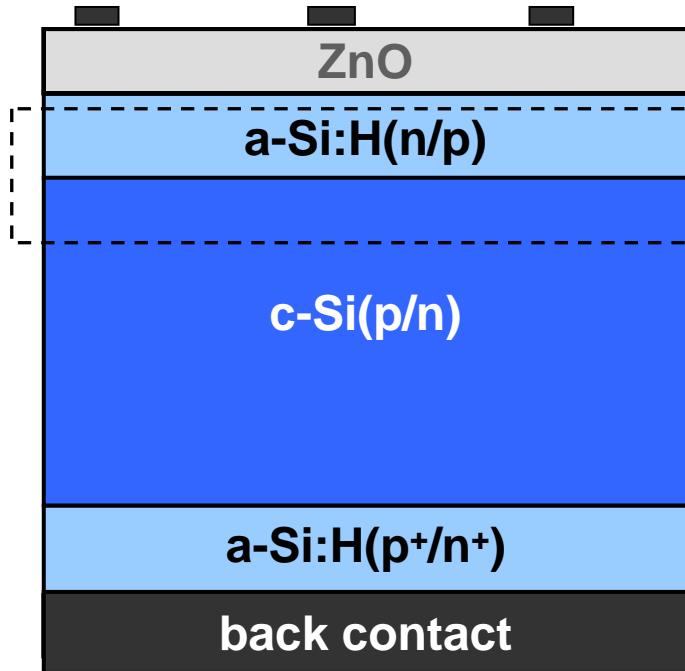
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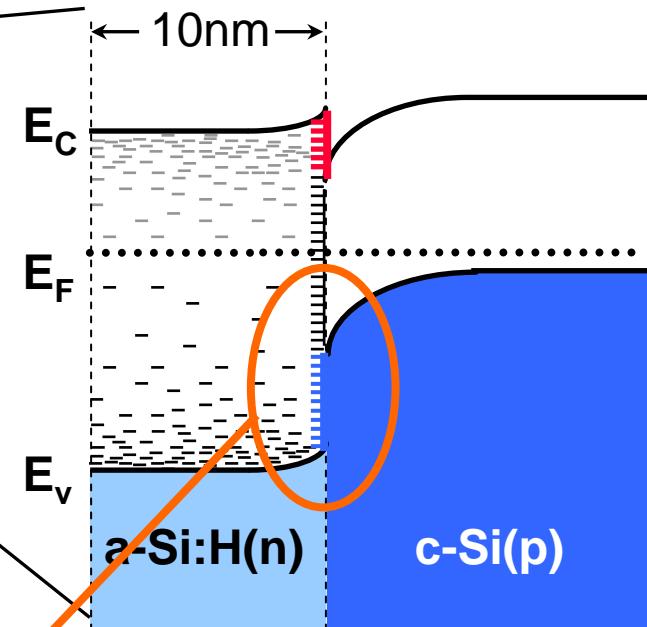


cell structure

metallization



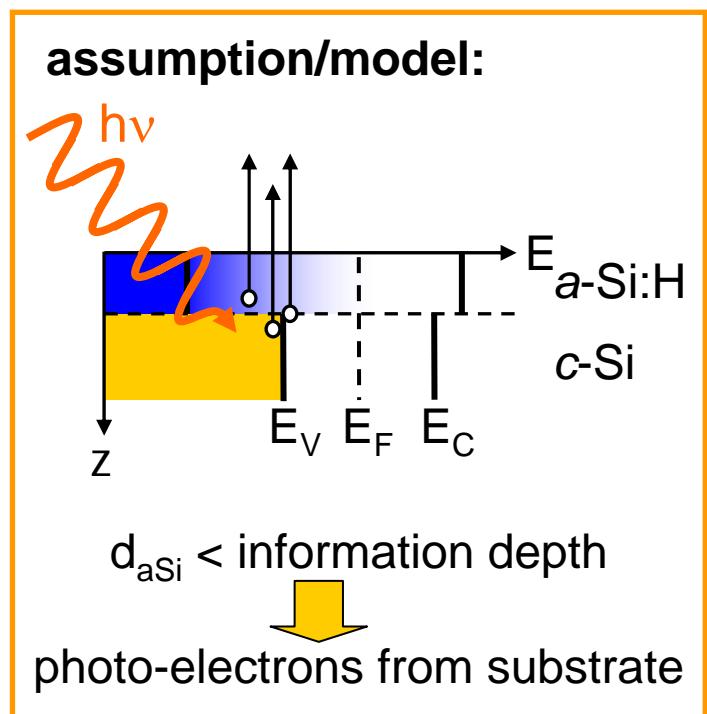
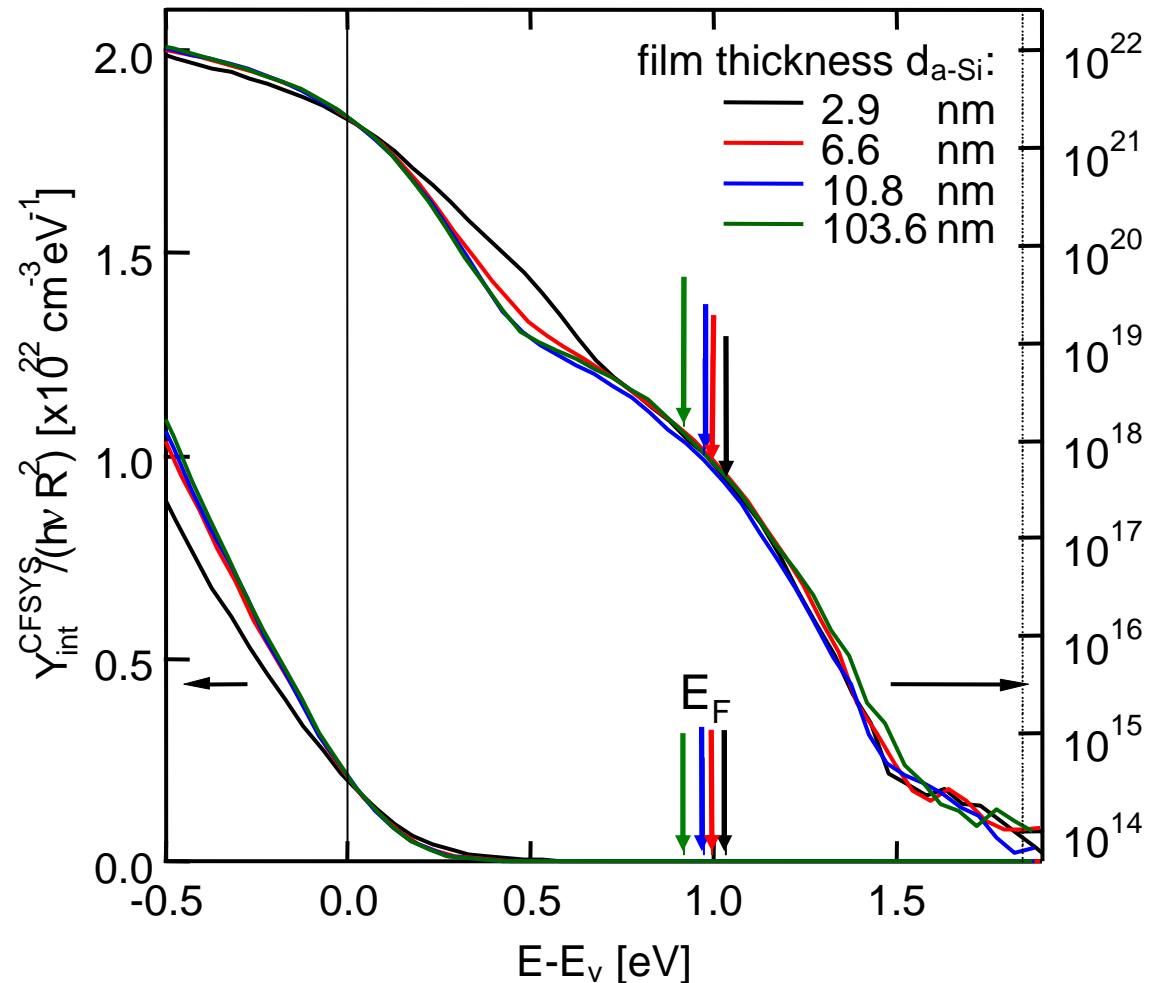
band diagram



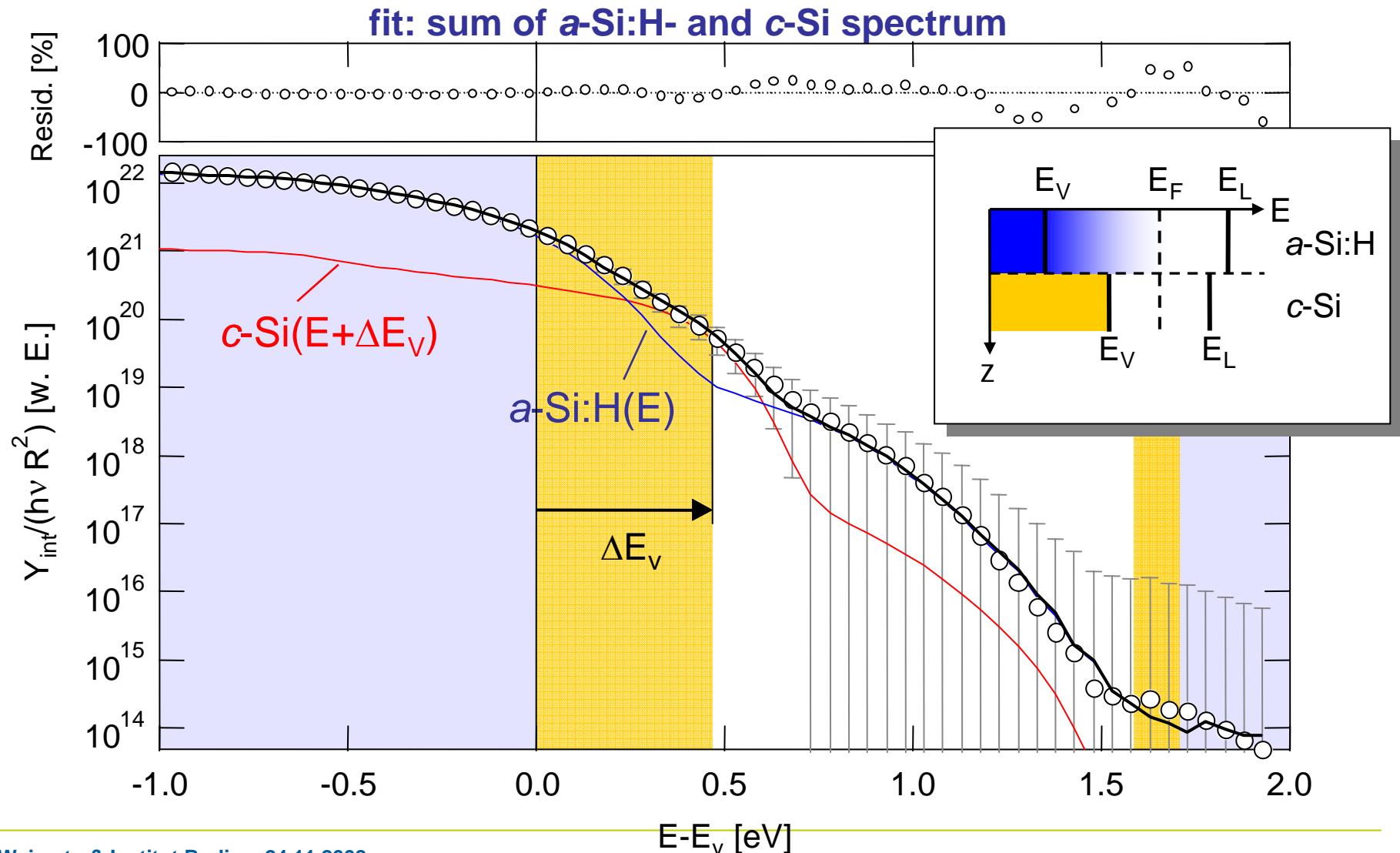
band offset?

a-Si:H/c-Si valence band offset

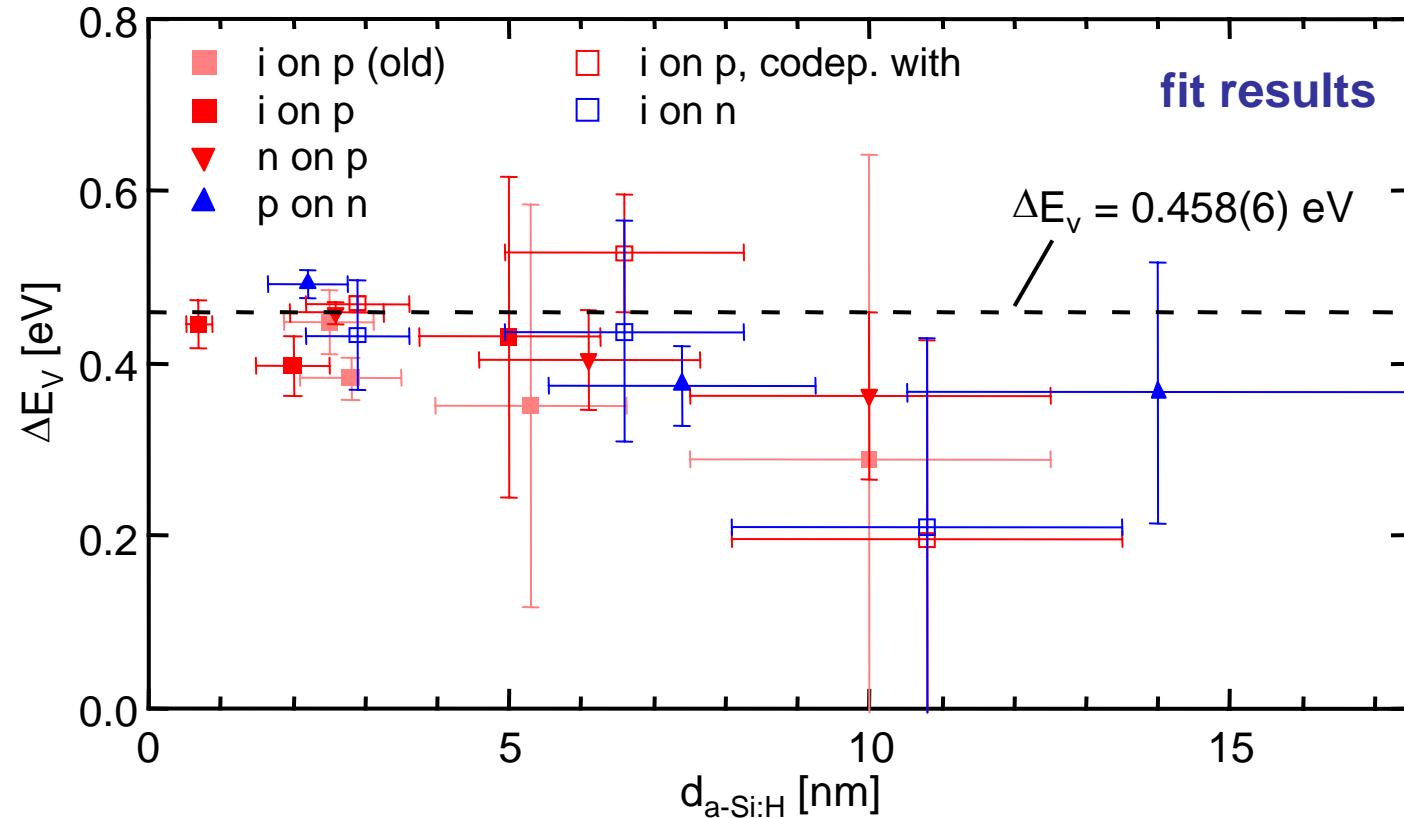
example: a-Si:H(i) / c-Si(p)



a-Si:H/c-Si valence band offset



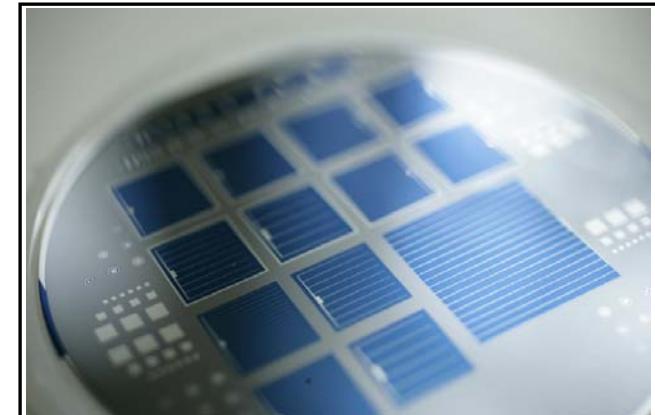
a-Si:H/c-Si valence band offset



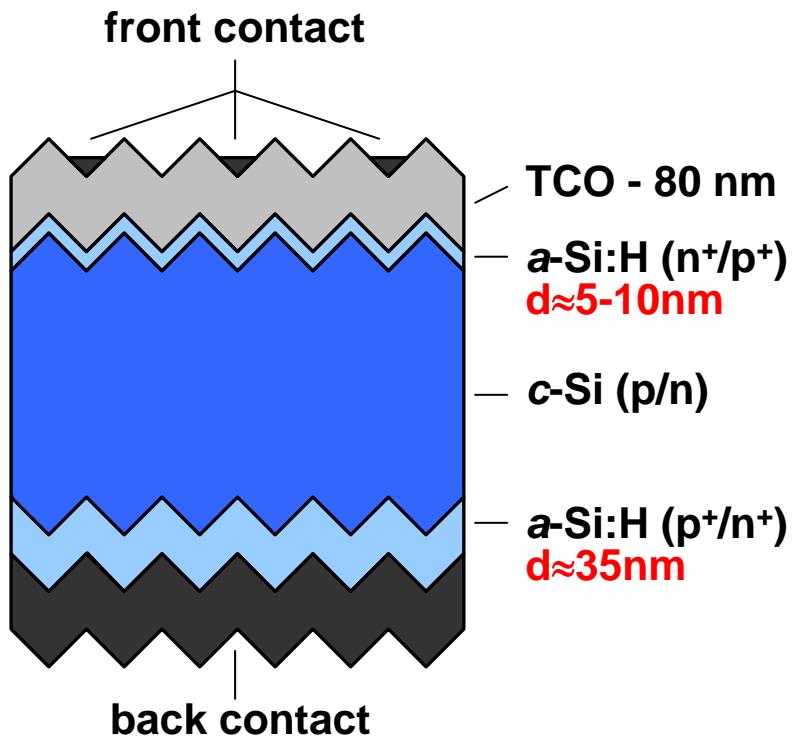
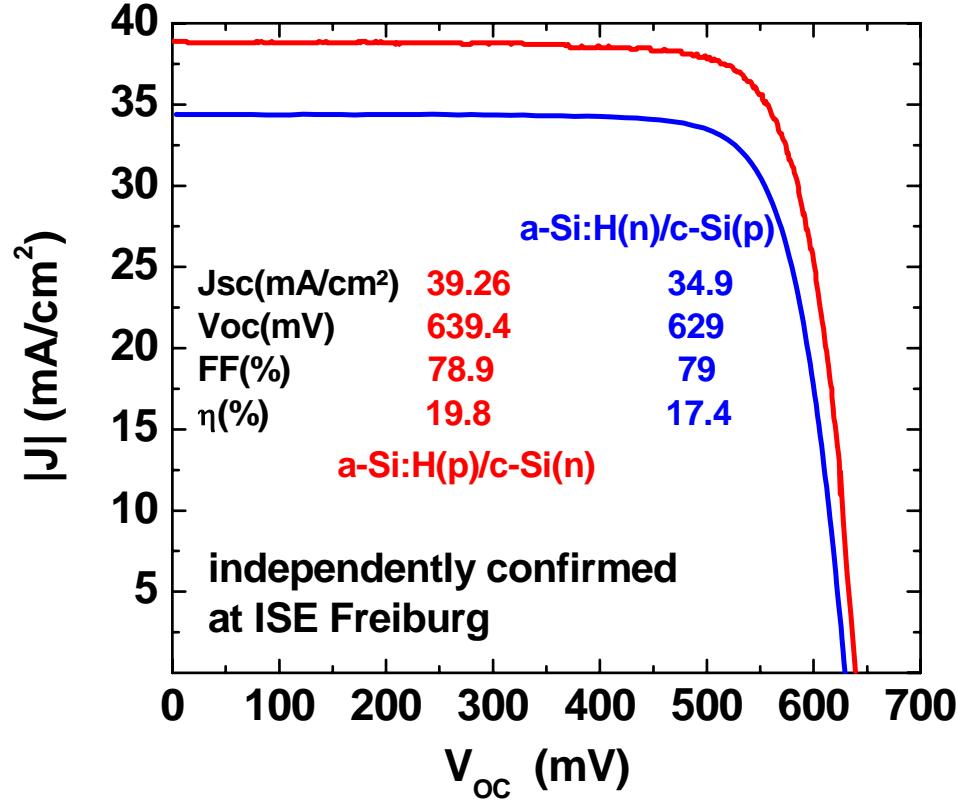
- mean: $\Delta E_V=0.458(6)$ eV (systematic error: ~50 meV)
- no dependence on substrate- or film doping
- (weak) trend: decreasing ΔE_V with film thickness - possible explanation: decreasing Si-H interface dipole, because Si-H bonds are substituted with Si-Si

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Solar cell results



doping sequence	texture	η	V _{oc}	J _{sc}	area
a-Si(n)/c-Si(p)/a-Si(p)	pyramids	17.4 %	629 mV	34.9 mA/cm ²	1 cm ²
	pyramids	18.4 %	629 mV	34.9 mA/cm ²	1 cm ²
← uncertified					
a-Si(p)/c-Si(n)/a-Si(n)	pyramids	19.8 %	639 mV	39.3 mA/cm ²	1 cm ²

Summary

a-Si:H growth

- growth mode on c-Si: islands – coalescence – thickening
 - high hydrogen content at a-Si/c-Si interface \leftrightarrow enhanced defect density
 - **optimum growth conditions:**
 - high c-Si surface quality prior to a-Si:H deposition
 - low plasma damage during growth
 - low defect density in a-Si:H “bulk”
- reduced recombination**

a-Si:H/c-Si interface

- device simulation: (effective) **interface DOS $< 10^{10} \text{ cm}^{-2}$**
 \rightarrow **not detrimental to cell** parameters
- asymmetric band offset: $\Delta E_V \sim 460 \text{ meV}$ ($\rightarrow \Delta E_C \sim 150\text{-}200 \text{ meV}$)

n-doped a-Si:H emitters

- Urbach-Energy and N_D increase with doping, comparable to thick films
- **optimum doping for device $\sim 2000 \text{ ppm}$** , not at minimum of $E_F\text{-}E_C$ –
trade-off: doping \leftrightarrow defect generation

Thank you

People

Walther Fuhs

Bernd Rech

Thomas Lußky

Matthias Schulz

Aziz Laades

Karsten von Maydell

Andreas Schöpke

Kerstin Jacob

Brunhilde Rabe

Dagmar Patzek

Funding



BMBF Netzwerk-Projekt Nr. 01SF0012

„Grundlagen und Technologie von Solarzellen auf der Basis von a-Si/c-Si Heterostrukturen“



EU FP7 project no. 211821

“Heterojunction Solar Cells based on a-Si c-Si”