

# Parametric X-ray radiation in the backward geometry under interaction of relativistic electrons with crystalline and polycrystalline structures

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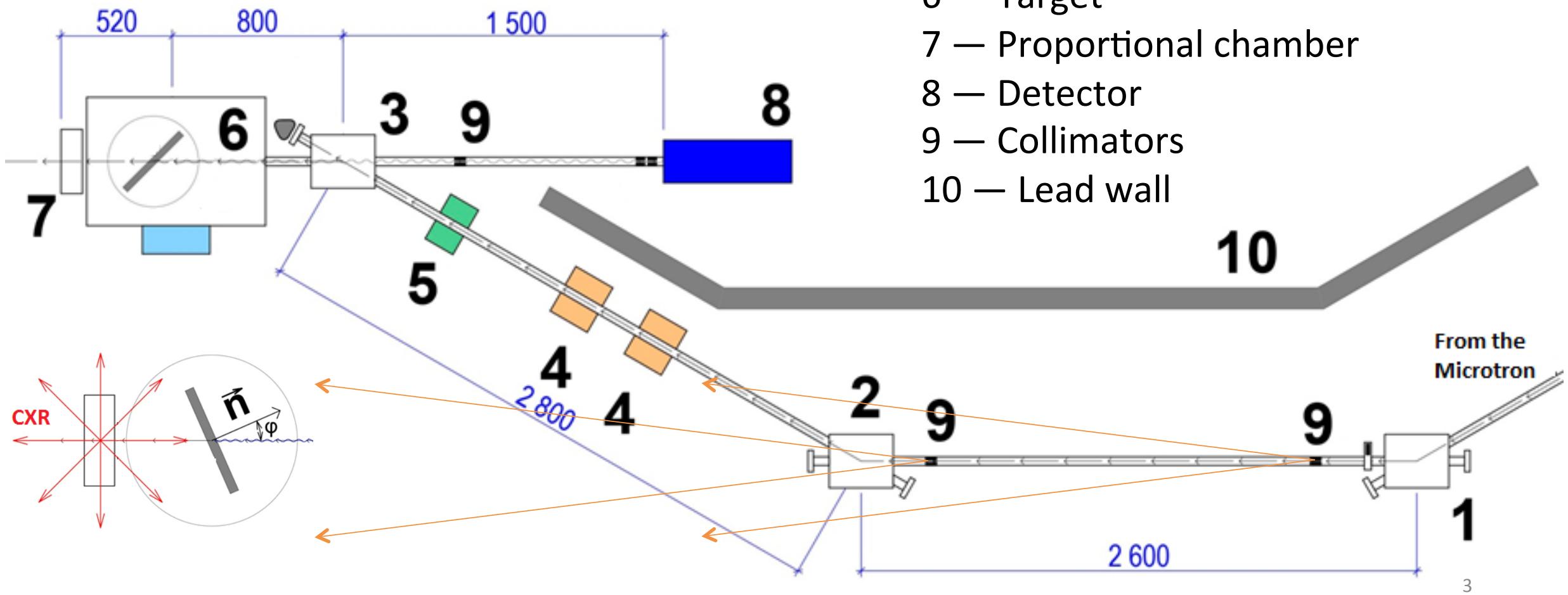
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# Plan

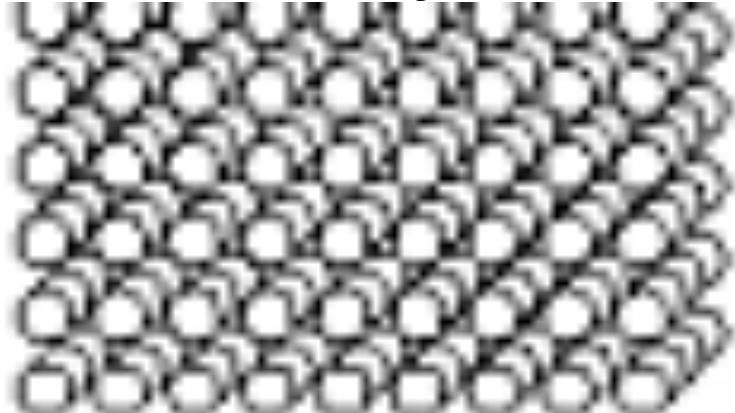
- Experimental setup
- PXR in:
  - Polycrystals (Nickel and Molybdenum)
  - Textured polycrystals (Nickel and Tungsten)
- Orientation dependence and the observation angle
- Conclusion

# Experimental setup



# Crystals

## Parametric X-ray Radiation (PXR)



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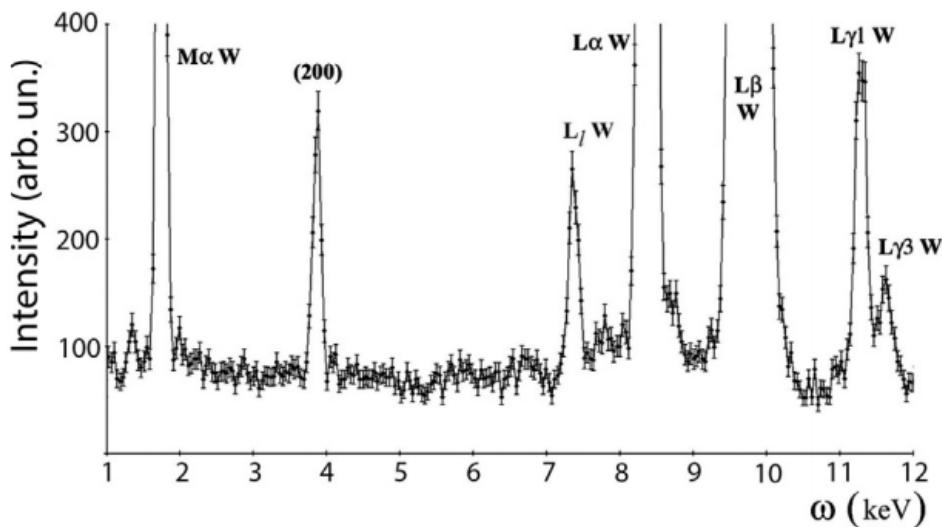
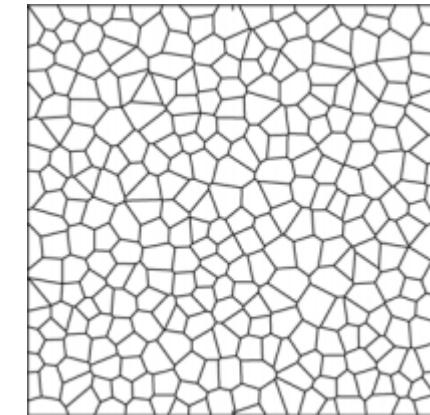


Fig. 9. The PBBS spectrum measured from W polycrystalline foil.

Tungsten Polycrystalline foil

# Polycrystals

## Polarization Bremsstrahlung (PB)



Mosaic  
crystals

Textured  
Polycrystals

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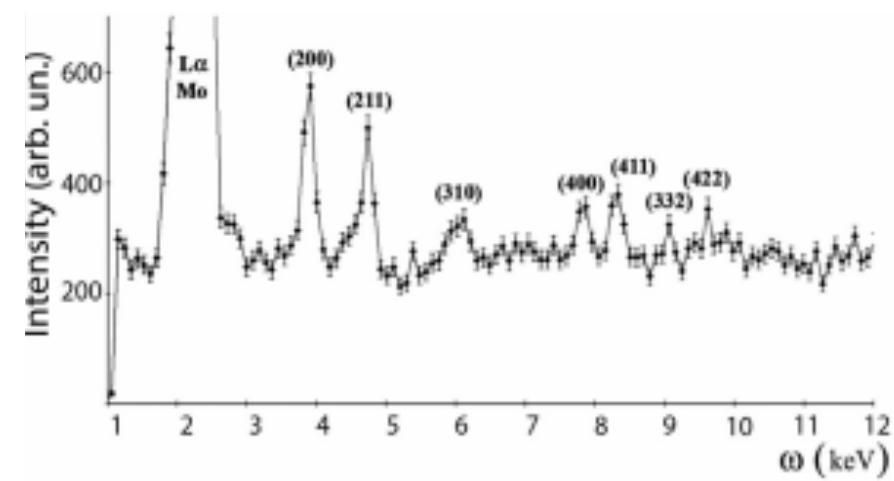
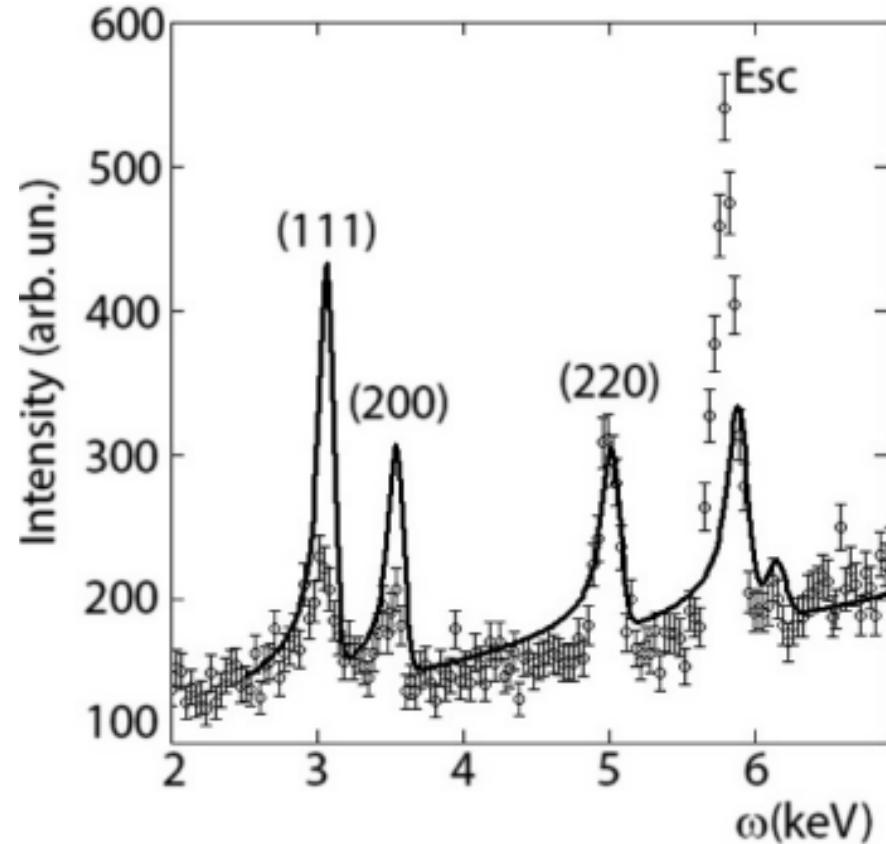


Fig. 8. The PBBS spectrum measured from Mo polycrystalline foil.

Molybdenum Polycrystalline foil

# Nickel polycrystalline foil

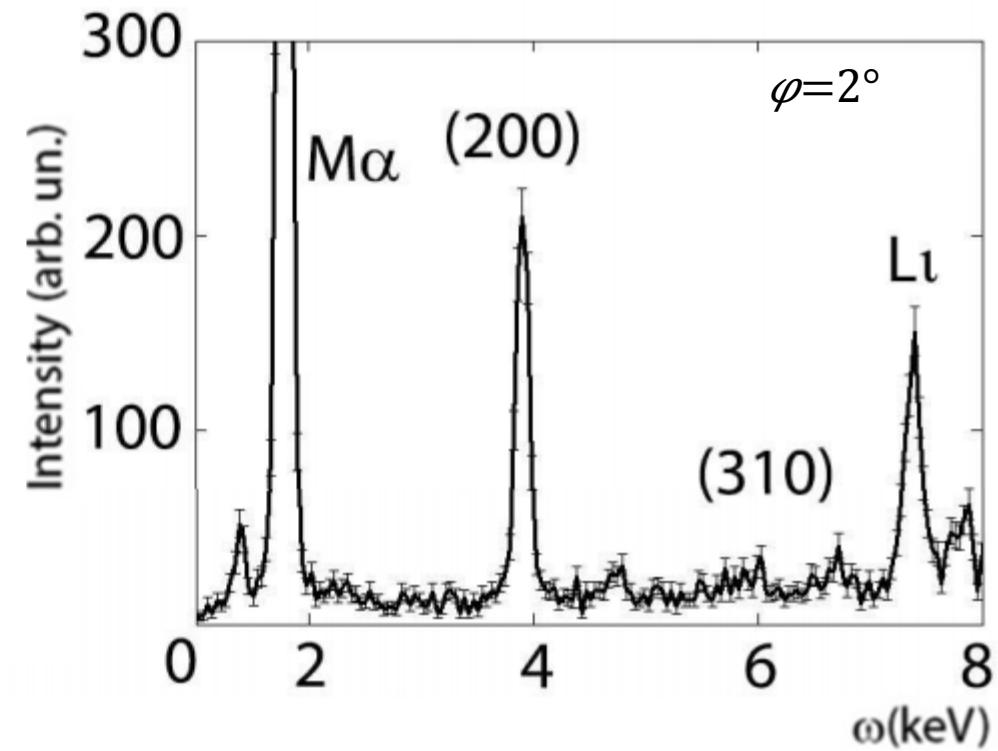
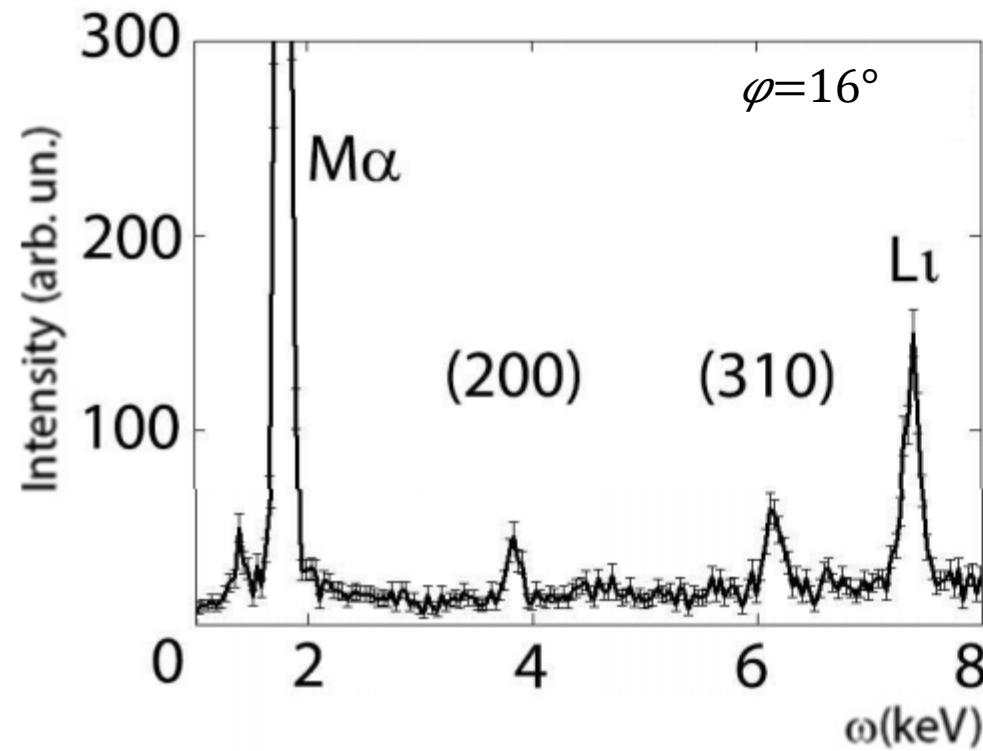
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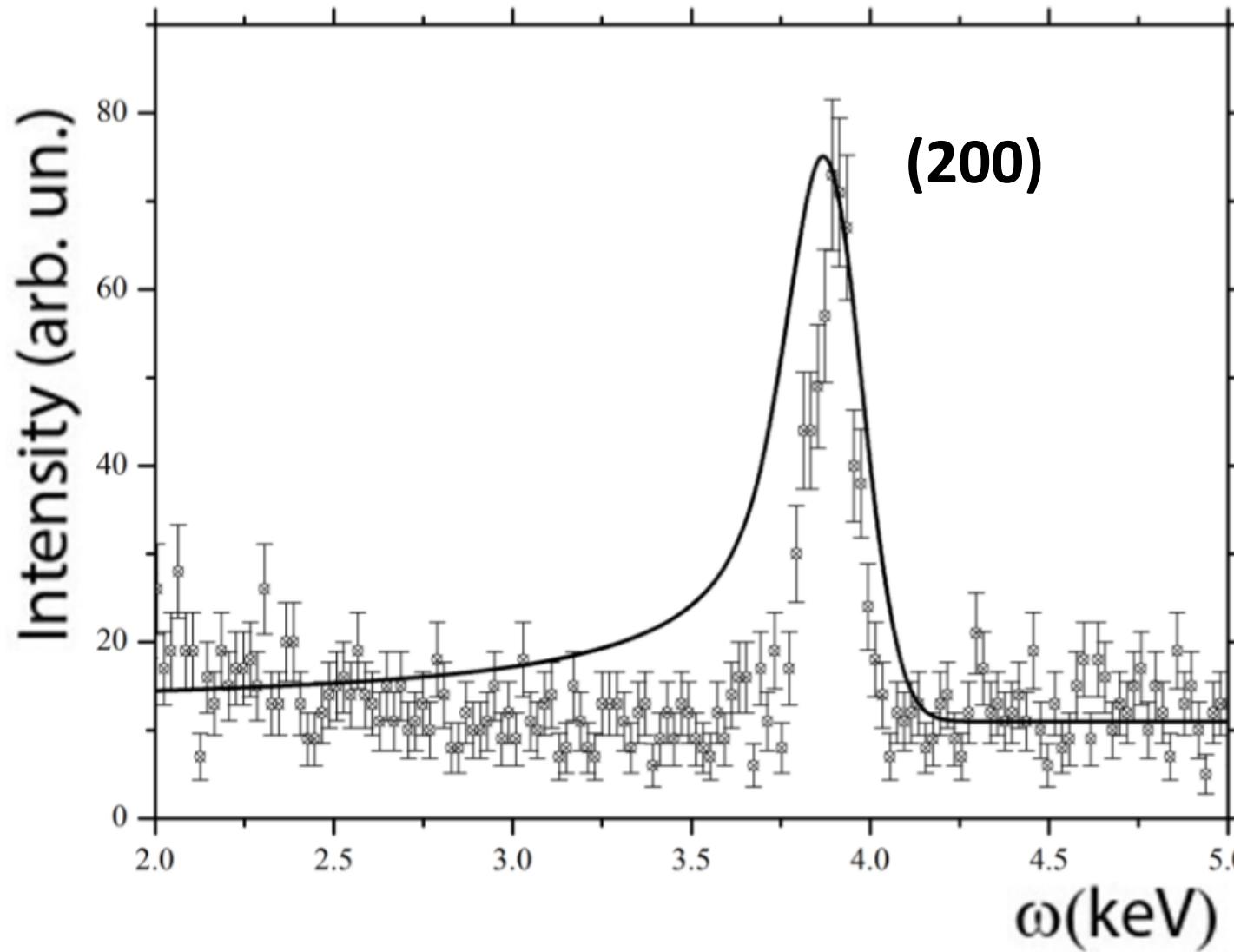
Theoretical peak position (220)	4.97 keV
Experimental peak position (220)	$4.940 \text{ keV} \pm 7 \text{ eV}$
Full Width Half Maximum (FWHM)	$159 \text{ eV} \pm 18 \text{ eV}$

**Fig. 6.** The PBBS spectrum measured from Ni polycrystalline foil with average grain size about 300 nm.

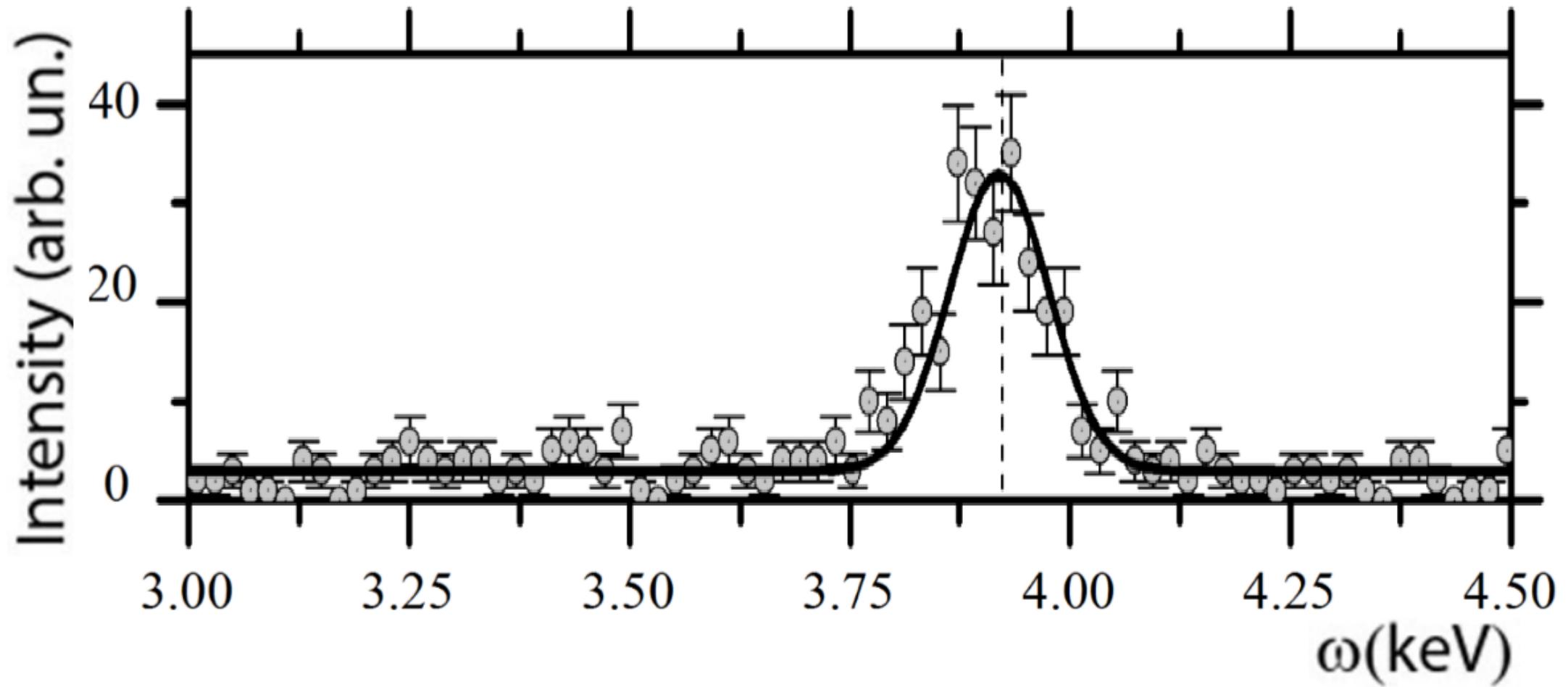
# Textured polycrystalline foil of Tungsten

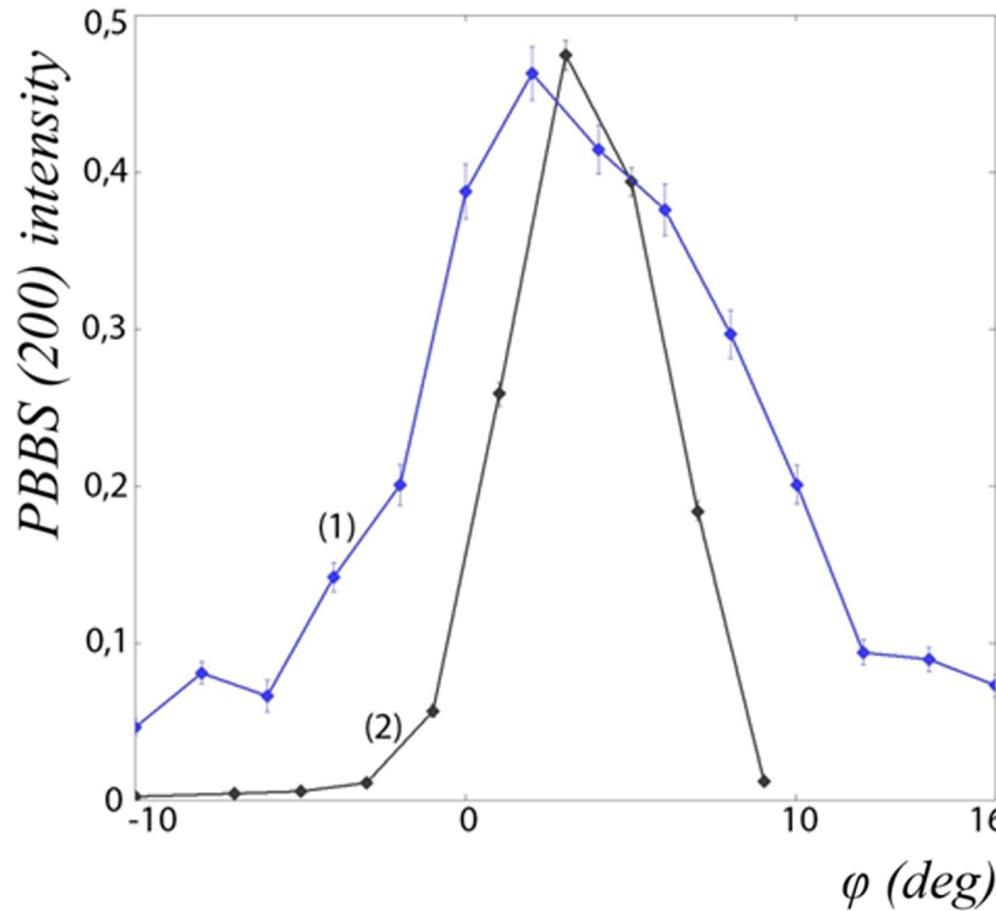


# Tungsten polycrystalline foil



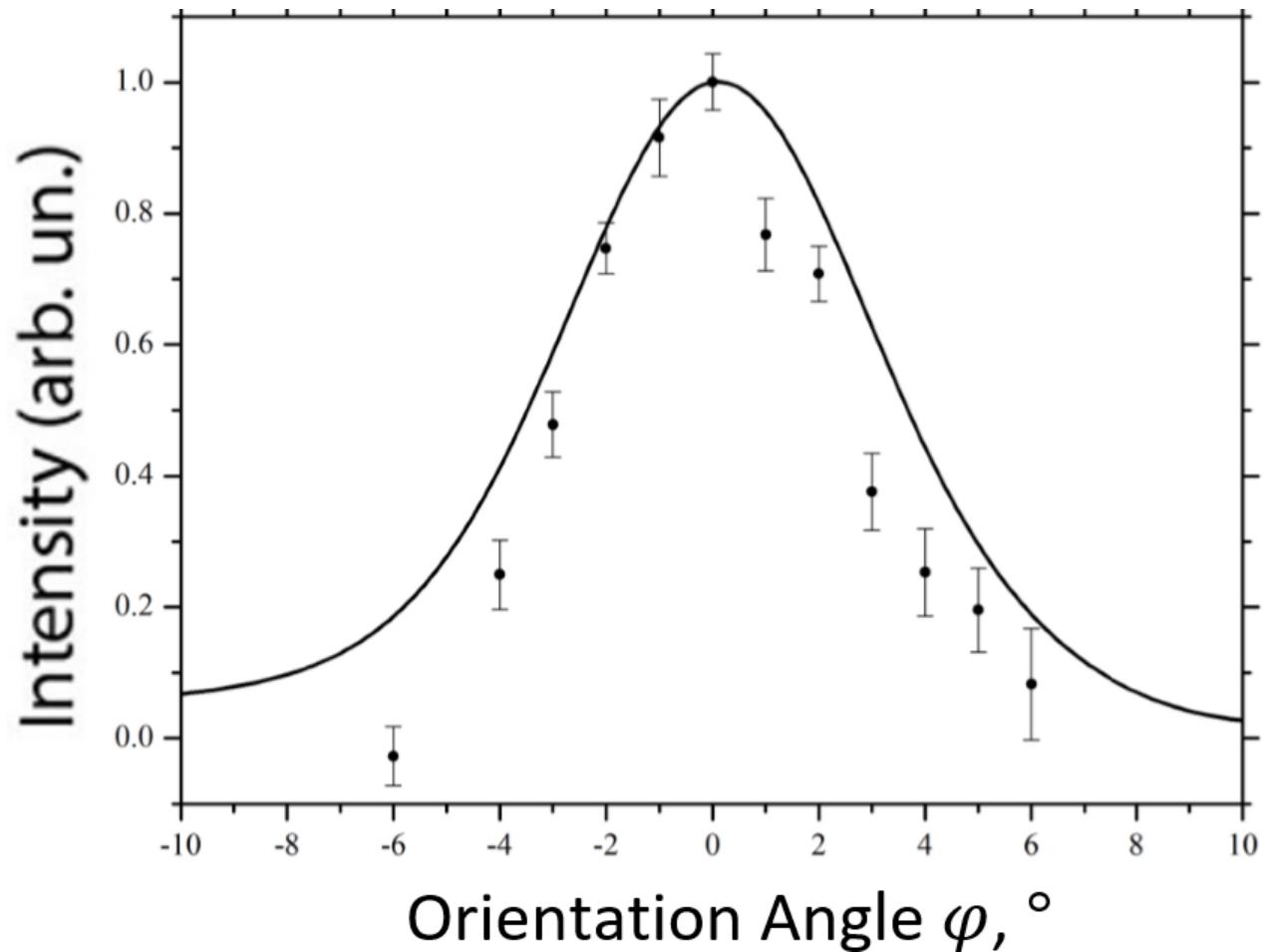
# Tungsten polycrystalline foil





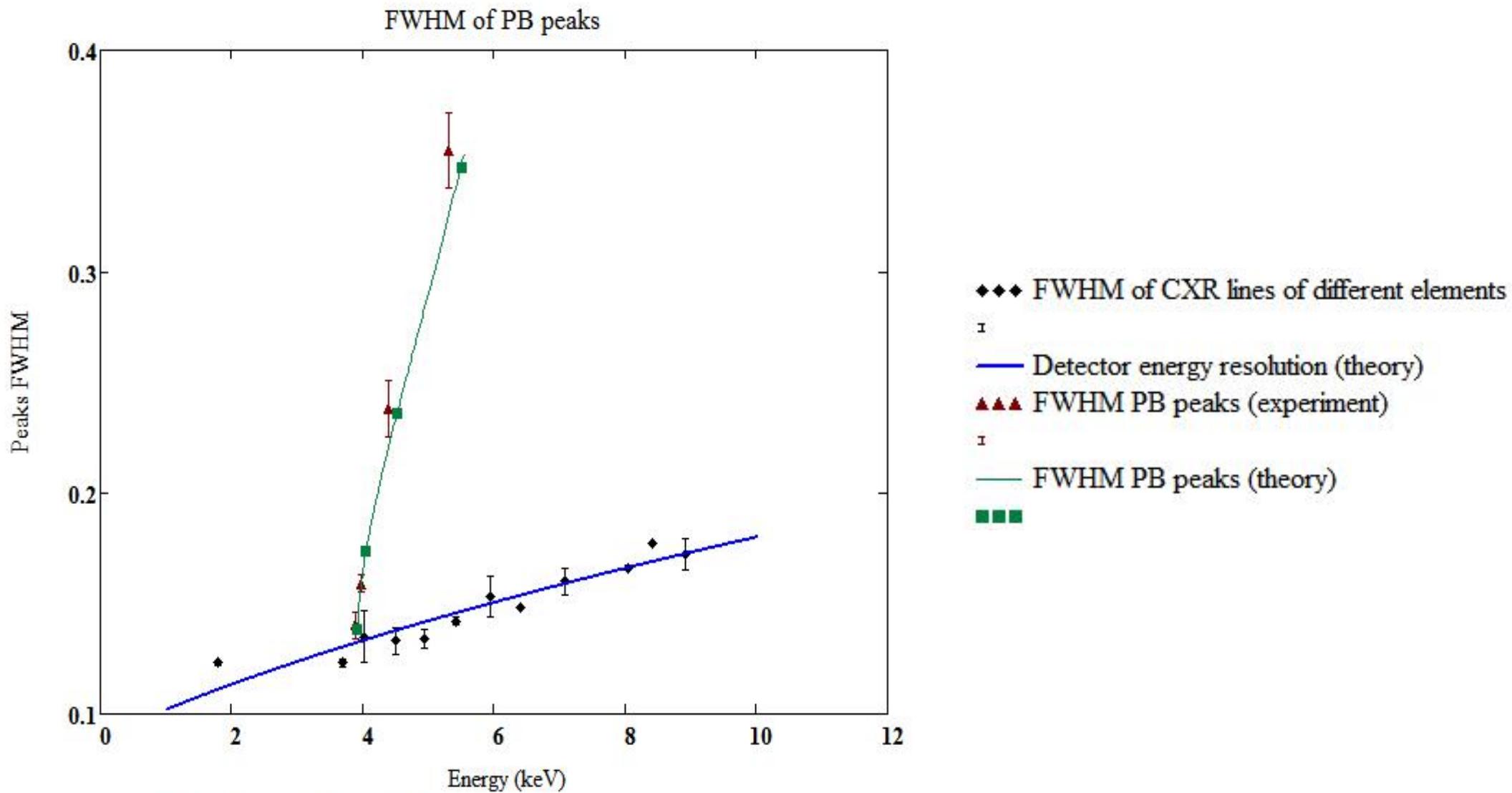
The orientation dependence of PBBS peak (200)  
from W textured polycrystalline foil (1). The comparison with XRD (2).

PBBS (220) distribution width	XRD distribution width
<b>10.2°</b>	<b>5.4°</b>



Orientation dependence of the (200) plane of the tungsten polycrystalline foil for an observation angle of  $90.8^{\circ}$

PXR (220) distribution width (Backward geometry)	PXR (220) distribution width (Backward geometry)	XRD distribution width
$10.2^{\circ}$	$5.7^{\circ}$	$5.4^{\circ}$



# Conclusions

- A deviation between the existing theories and measurements is observed for textured polycrystals which can be considered as a transition structure between crystals and polycrystals.
- The FWHM of the orientation dependence changes when the observation angle changes.
- Is the backward geometry the most suitable to diagnostics polycrystalline materials using PXR?

Thank you for your attention!