

Reciprocal space mapping and single-crystal scattering rods. Erratum

Detlef-M. Smilgies,^{a*} Daniel R. Blasini,^{a,b} Shu Hotta^c and Hisao Yanagi^d

^aCHESS Center, Cornell University, Ithaca, NY 14853, USA, ^bDepartment of Chemistry and Chemical Biology, Cornell University, Ithaca, NY 14853, USA, ^cDepartment of Polymer Science and Engineering, Kyoto Institute of Technology, Matsugasaki, Sakyo-ku, Kyoto 606-8585, Japan, and ^dFaculty of Engineering, Kobe University, Rokkodai, Nada-ku, Kobe 657-8501, Japan. E-mail: dms79@cornell.edu

While assembling the high-resolution graphics for Fig. 2 of Smilgies *et al.* [*J. Synchrotron Rad.* (2005). **12**, 807–811], the curve from the bottom panel was erroneously repeated in the top panel. Below is the corrected figure.

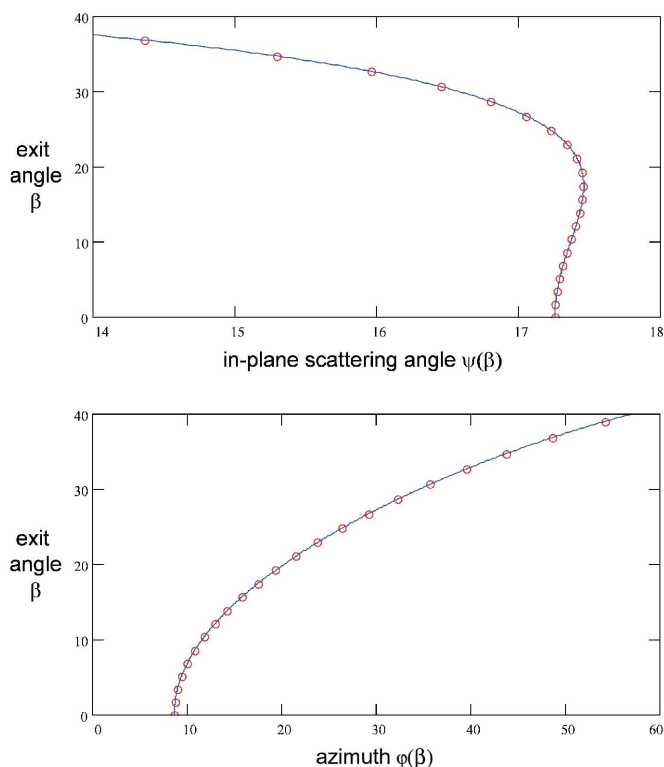


Figure 2

Exit-angle dependence of the scattering angle ψ (top panel) and the sample azimuth ϕ (bottom panel). Solid lines, based on equations (5a) and (5b), respectively, were compared with angular calculations (dots) using the *sixc* diffractometer code (Lohmeier & Vlieg, 1993). Axes were chosen such that they correspond to the experimental scattering intensity maps.

References

Lohmeier, M. & Vlieg, E. (1993). *J. Appl. Cryst.* **26**, 706–716.