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Addition/Correction

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Correction to Orientation of Phenylphosphonic Acid Self-Assembled Monolayers on a Transparent Conductive Oxide: A Combined NEXAFS, PM-IRRAS, and DFT Study

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In the Results and Discussion section, there are several instances where α and φ have been inadvertently interchanged in the text. The data were processed using the correct definitions, eq 4 and Figure 1 are correct, and the conclusions remain unaffected. However, for consistency throughout the Article, α should always be the orientation of the molecular dipole moment relative to the surface normal, and φ is the angle that the incoming X-rays make with the surface normal. The following typographical errors may confuse the reader:

On page 2168, α is given as the incident X-ray angle. φ is the incident X-ray angle.

On page 2169, (1) the angle that the transition dipole moment makes with the surface normal is given erroneously as ϕ. α is the angle between the molecular dipole moment and the surface normal. (2) α is again referenced incorrectly as the incident X-ray angle. (3) The sample curves given by dotted lines in Figure 1b are plots of I_v (ϕ) (eq 4) for values of α = 0−90° in 10° steps. The orientation of the phenyl ring normal is correct but is incorrectly referred to in the text as ϕ (although the figure notation is correct). The best fit of I_v (ϕ) to the data in Figure 1b yields the result α = 77 ± 5°.

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