PANASQUEIRAITE, A NEW MINERAL: THE OH-EQUIVALENT OF ISOKITE. ERRATUM

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As reported in the paper "Panasqueiraite, the OH-equivalent of isokite" by A.M. Isaacs and D.R. Peacor, which appeared in volume 19, pages 389-392, errors occur in the section on "Crystallography and Structure". The space group of panasqueiraite is incorrectly described as $C2/c$ for the unit cell as reported (a 6.535, b 8.753, c 6.191 Å, β 112.33°) and should be $C2/n$. Furthermore, the results obtained by Deans & McConnell (1955) for isokite (which is isostructural with panasqueiraite) were incorrectly reported to be in error. The unit cell of panasqueiraite as given above may be transformed to one having a 6.535, b 8.753, c 7.498 Å, β 121.40°. With this setting the space group is $C2/c$ and all data are consistent with those reported by Deans & McConnell (1955) for isokite.
