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Corrigendum

Corrigendum to "Pore scale modeling of a proton exchange membrane fuel cell catalyst layer: Effects of water vapor and temperature" [J. Power Sources, 196 (6) (2011) 3195–3203]

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Due to an incorrect conversion of units, the thermal conductivity values in Table 2 were given incorrectly and consequently Figs. 6 and 7 were incorrectly scaled on the *y*-axis. The table and figures should have appeared as follows:

Table 2
Thermal conductivities for different regions of the catalyst layer.

Region	λ (W/cm K)
Pore	2.99×10^{-4}
Ionomer	2.0×10^{-3} [28]
Carbon	3.75×10^{-3} [29]

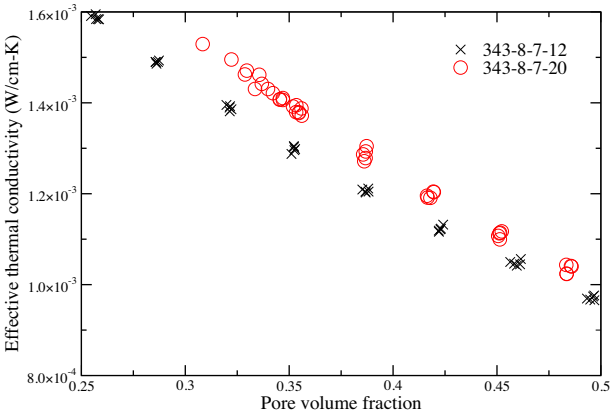

Fig. 6. Effective thermal conductivities as a function of pore volume fraction. Labels: T (K), c_{H_2O} (mol m^{-3}), c_{O_2} (mol m^{-3}), sphere radius (nm).

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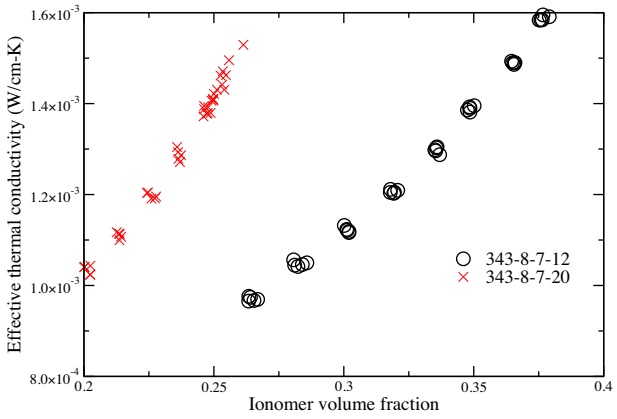

Fig. 7. Effective thermal conductivities as a function of ionomer volume fraction. Labels: T (K), c_{H_2O} (mol m^{-3}), c_{O_2} (mol m^{-3}), sphere radius (nm).