

SPECIAL COLLECTION: BUILDING PLANETS: THE DYNAMICS AND GEOCHEMISTRY OF CORE FORMATION

The W-WO<sub>2</sub> oxygen fugacity buffer (WWO) at high pressure and temperature:  
Implications for  $f_{O_2}$  buffering and metal-silicate partitioning

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ABSTRACT

Synchrotron X-ray diffraction data were obtained to simultaneously measure unit-cell volumes of W and WO<sub>2</sub> at pressures and temperatures up to 70 GPa and 2300 K. Both W and WO<sub>2</sub> unit-cell volume data were fit to Mie-Grüneisen equations of state; parameters for W are  $K_T = 307 (\pm 0.4)$  GPa,  $K'_T = 4.05 (\pm 0.04)$ ,  $\gamma_0 = 1.61 (\pm 0.03)$ , and  $q = 1.54 (\pm 0.13)$ . Three phases were observed in WO<sub>2</sub> with structures in the  $P2_1/c$ ,  $Pnma$ , and  $C2/c$  space groups. The transition pressures are 4 and 32 GPa for the  $P2_1/c$ - $Pnma$  and  $Pnma$ - $C2/c$  phase changes, respectively. The  $P2_1/c$  and  $Pnma$  phases have previously been described, whereas the  $C2/c$  phase is newly described here. Equations of state were fitted for these phases over their respective pressure ranges yielding the parameters  $K_T = 238 (\pm 7)$ ,  $230 (\pm 5)$ ,  $304 (\pm 3)$  GPa,  $K'_T = 4$  (fixed),  $4$  (fixed),  $4$  (fixed) GPa,  $\gamma_0 = 1.45 (\pm 0.18)$ ,  $1.22 (\pm 0.07)$ ,  $1.21 (\pm 0.12)$ , and  $q = 1$  (fixed),  $2.90 (\pm 1.5)$ ,  $1$  (fixed) for the  $P2_1/c$ ,  $Pnma$ , and  $C2/c$  phases, respectively. The W-WO<sub>2</sub> buffer (WWO) was extended to high pressure using these W and WO<sub>2</sub> equations of state. The  $T$ - $f_{O_2}$  slope of the WWO buffer along isobars is positive from 1000 to 2500 K with increasing pressure up to at least 60 GPa. The WWO buffer is at a higher  $f_{O_2}$  than the iron-wüstite (IW) buffer at pressures lower than 40 GPa, and the magnitude of this difference decreases at higher pressures. This implies an increasingly lithophile character for W at higher pressures. The WWO buffer was quantitatively applied to W metal-silicate partitioning by using the WWO-IW buffer difference in combination with literature data on W metal-silicate partitioning to model the exchange coefficient ( $K_D$ ) for the Fe-W exchange reaction. This approach captures the non-linear pressure dependence of W metal-silicate partitioning using the WWO-IW buffer difference. Calculation of  $K_D$  along a peridotite liquidus predicts a decrease in W siderophilicity at higher pressures that supports the qualitative behavior predicted by the WWO-IW buffer difference, and agrees with findings of others. Comparing the competing effects of temperature and pressure the results here indicate that pressure exerts a greater effect on W metal-silicate partitioning.

**Keywords:** High pressure, tungsten, oxygen fugacity buffer, equation of state, metal-silicate partitioning