Supplementary Material (ESI) for Dalton Transactions

KZn_4SbO_7 and $KZn_4Sb_3O_{12}$: Syntheses, Structures and Photophysics of Sb^{5+} Control Materials

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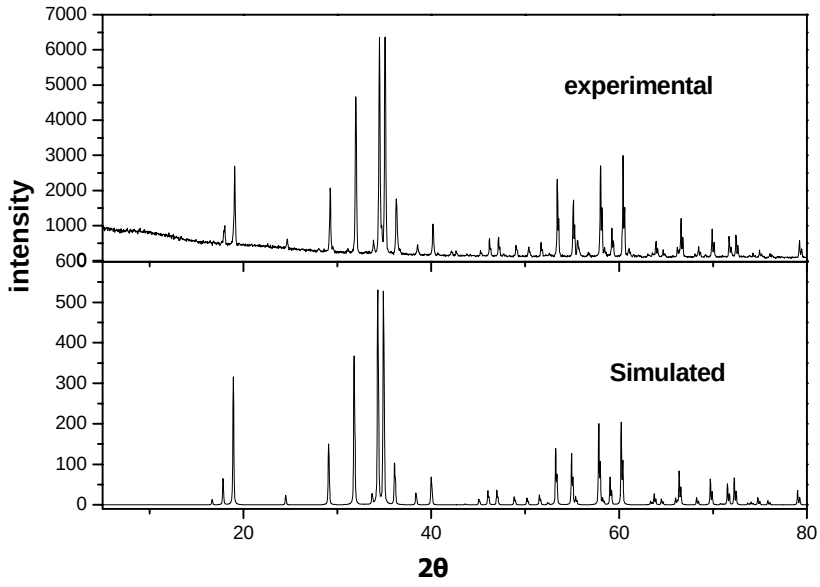
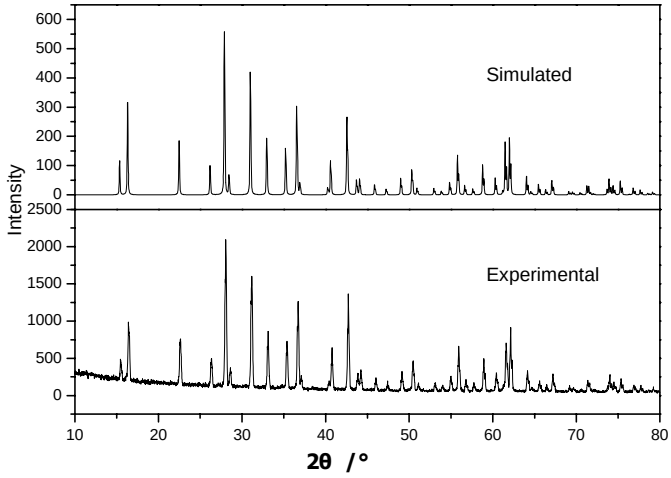
Supporting Information



a: KZn_4SbO_7



b: $KZn_4Sb_3O_{12}$

Figure S1. Simulated and experimental XRD powder patterns for **1** (a) and **2** (b).

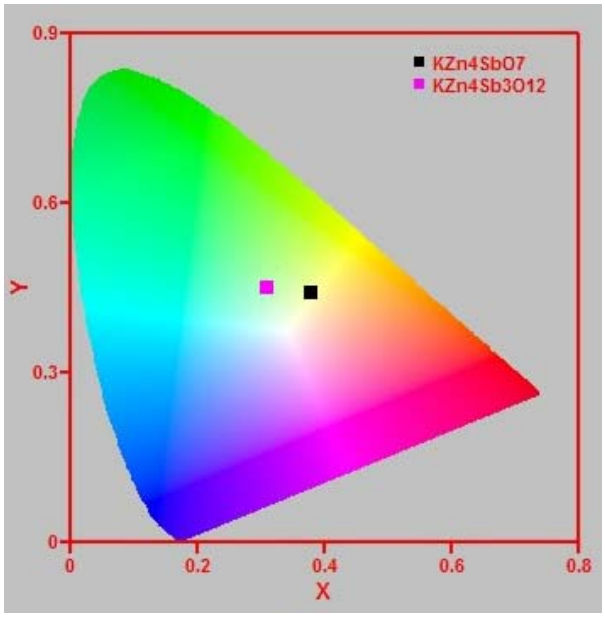
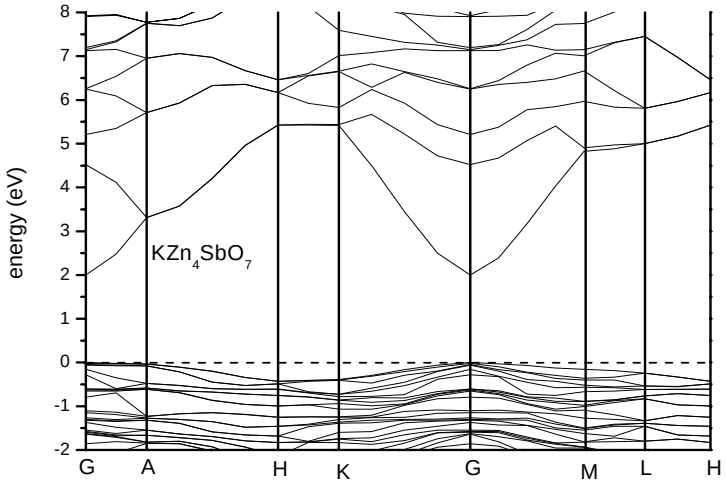
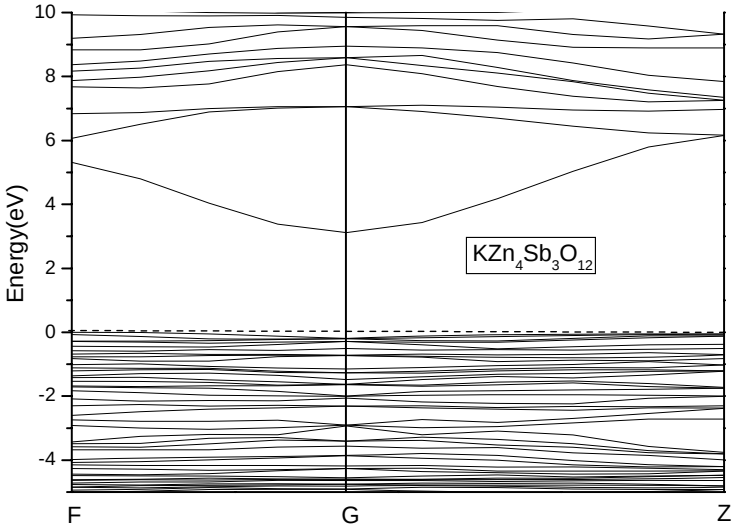


Figure S2. The calculated CIE chromaticity coordinates of (0.380, 0.440) and (0.310, 0.450) close to the coordinates of white-light emission for KZn_4SbO_7 and $KZn_4Sb_3O_{12}$.





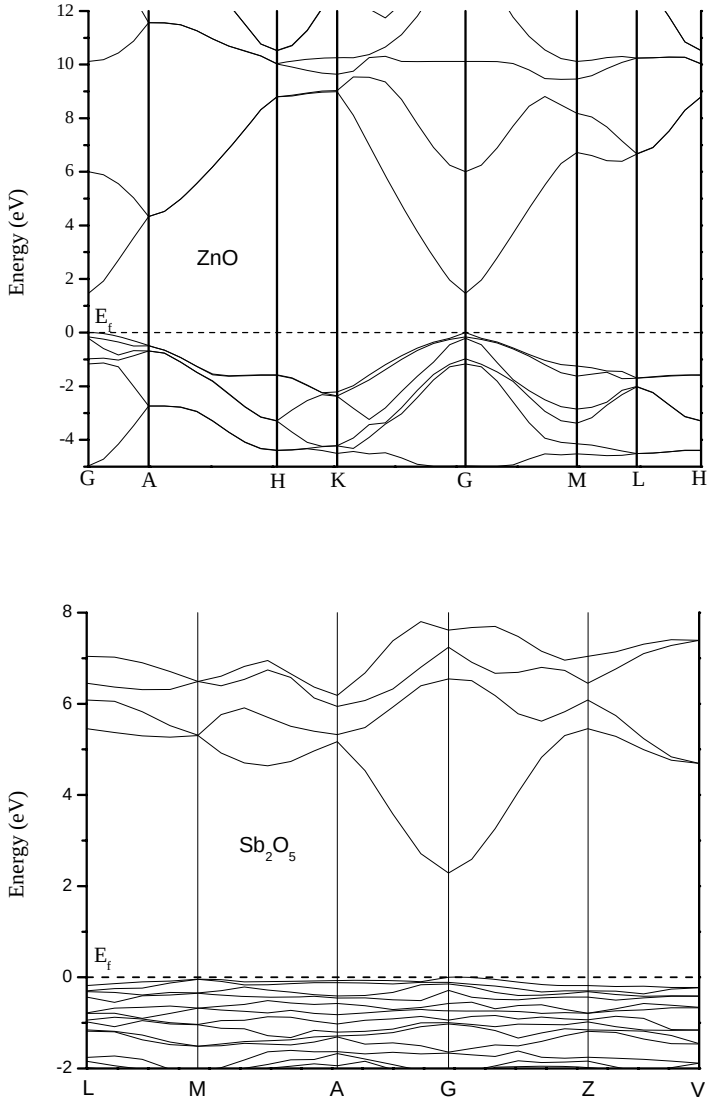


Figure S3. The band structures of plots. The lowest conduction band is at G-point and the highest valence band at F-point for $KZn_4Sb_3O_{12}$ crystal; both the lowest conduction band and highest valence band are at G-point for KZn_4SbO_7, ZnO and Sb_2O_5 crystals.