Supplementary Information

Oxide-Oxide Nanojunctions in coaxial SnO_2/TiO_2 , SnO_2/V_2O_3 and $SnO_2/(Ti_{0.5}V_{0.5})_2O_3$ Nanowire Heterostructures

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S1. Atomic simulation of the interfaces between rutile $SnO_2 NW$ core and different shells: (a) Ti_2O_3 , (b) Ti_2O_3 , (c) TiO2 rutile and (d) TiO₂ anatase.



Figure S1. Atomic simulation of the interfaces between rutile SnO_2 NW core and different shells: (a) Ti_2O_3 , (b) Ti_2O_3 , (c) TiO2 rutile and (d) TiO₂ anatase. Top views show the $[10-10]_{Ti2O}$ // $[0-11]_{SnO2}$ in the case of Ti_2O on SnO_2 , $[11-20]_{Ti2O3}$ // $[0-11]_{SnO2}$ zone axis in the case of Ti_2O_3 on SnO_2 and $[0-11]_{TiO2}$ // $[0-11]_{SnO2}$ in the case of TiO_2 rutile and anatase on SnO_2 . Front views are visualized along the $[0001]_{Ti2O/Ti2O3}$ // $[100]_{SnO2}$ zone axis in the case of Ti_2O_3 and Ti_2O_3 on SnO_2 , and along the $[100]_{TiO2}$ // $[100]_{SnO2}$ zone axis in the case of TiO_2 rutile and anatase on SnO_2 . Finally, the side view corresponds to the $[-12-10]_{Ti2O}$ // $[011]_{SnO2}$, $[-1100]_{Ti2O3}$ // $[011]_{SnO2}$ and $[011]_{TiO2}$ anat/rutile // $[011]_{SnO2}$ directions. In each case mismatch percentage is indicated under the model.

S2. Formation mechanism of SnO2- based heterostructures.



Figure S2. Formation mechanism of SnO2- based heterostructures. While titania shells degrade, vanadia overlayers show improved epitaxial relationships and thermochemical stability.



S3. In-situ mass spectrometical analysis of the gas phase during the deposition process.

Figure S3. *In-situ* mass spectrometical analysis of the gas phase during the deposition process of $Ti(OPr^{i})_{4}$: m/z=2 (H₂⁺), 18 (H₂O⁺), 15 (CH₃⁺), 42 (C₃H₅⁺), 45 (C₂H₆O⁺), 59 (C₃H₇O⁺).