

Supporting Material:

Tracking Site-Specific C–C Coupling of Formaldehyde Molecules on Rutile TiO₂(110)

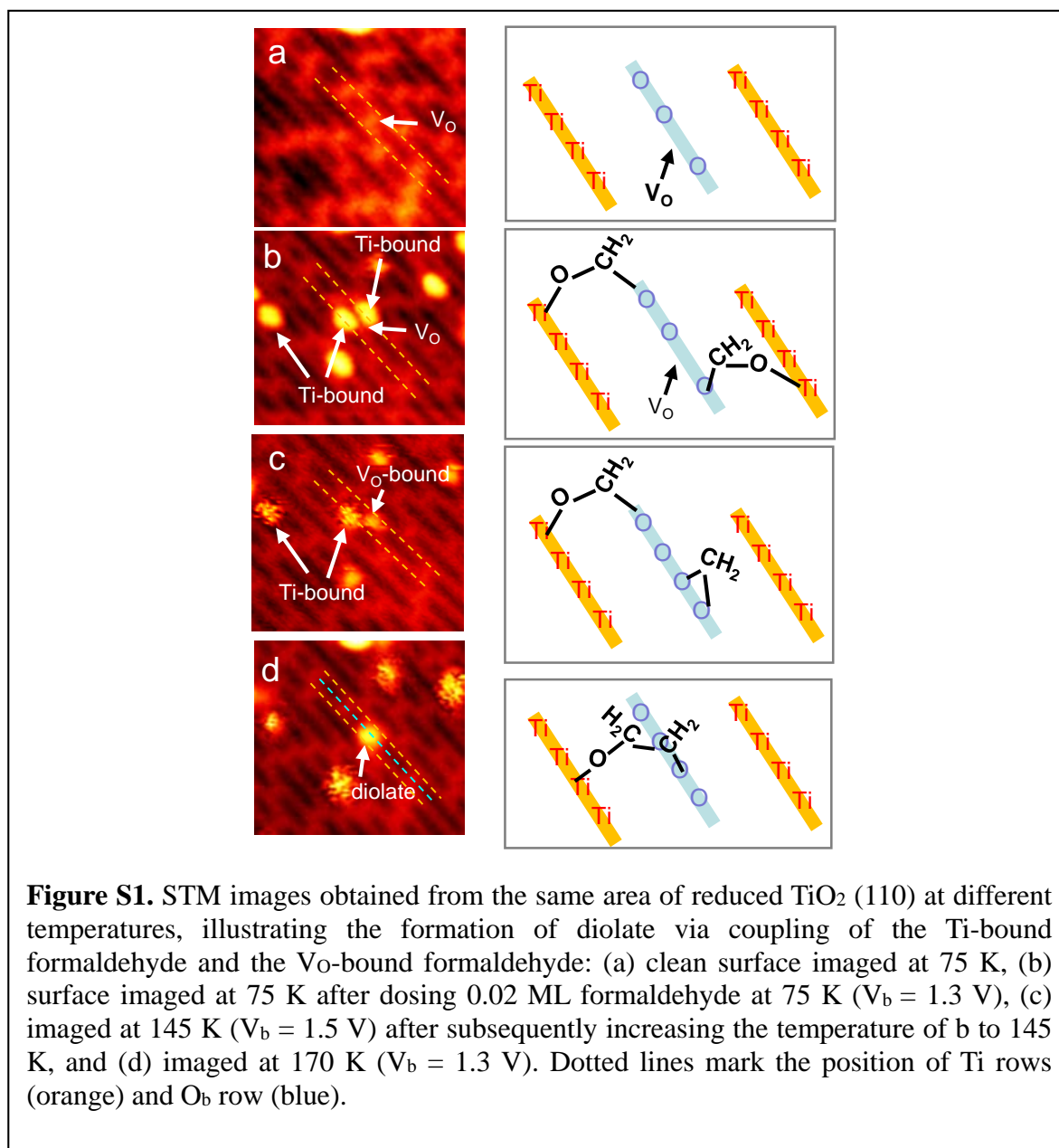
Ke Zhu,[†] Yaobiao Xia,[†] Miru Tang,[‡] Zhi-Tao Wang,[§] Bryan Jan,[†] Igor Lyubinetsky,[§] Qingfeng Ge,[‡] Zdenek Dohnálek,^{||} Kenneth T. Park,^{*,†} and Zhenrong Zhang^{*,†}

[†]Department of Physics, Baylor University, Waco, Texas 76798, United States

[‡]Department of Chemistry and Biochemistry, Southern Illinois University, Carbondale, Illinois 62901, United States

[§]Environmental Molecular Sciences Laboratory and ^{||}Fundamental and Computational Sciences Directorate, Institute for Interfacial Catalysis, Pacific Northwest National Laboratory, Richland, Washington 99352, United States

S1 Additional figure illustrating the formation of diolate



S2 The STM images showing the initial and the final status of sequence A.

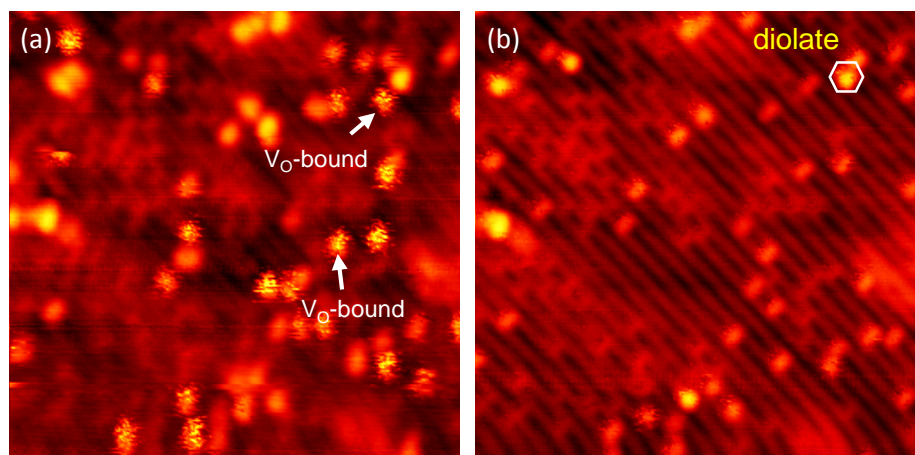


Figure S2. The isothermal STM images showing (a) the initial and (b) the final status of sequence A taken at 190 K.

S3 The images showing the initial and the final status of sequence B.

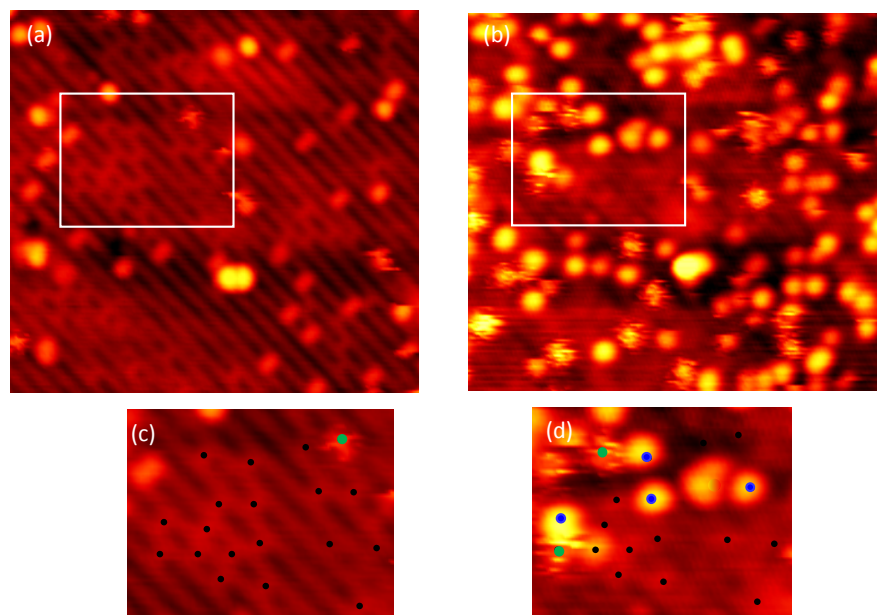


Figure S3. The isothermal STM images showing (a) the initial and (b) the final status of sequence B taken at 190 K. (c) and (d): Magnified areas marked in (a) and (b). Black dots mark the position of V_O 's. Blue dots and green dots mark the position of diolates and V_O -bound CH_2O molecules.

S4 Another set of statistical analysis performed on the images of a surface at 180 K with less V_o -concentration (0.075 ML) to compliment Figure 3A.

