

Supporting Information for “Design of robust 2,2′-bipyridine ligand linkers for the stable
immobilization of molecular catalysts on silicon(111) surfaces”

Samantha I. Johnson, James D. Blakemore, Bruce S. Brunshwig, Nathan S. Lewis, Harry B.
Gray, William A. Goddard III*, Petter Persson*

Geometries

1

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Cartesian coordinates for molecules calculated in this manuscript can be found in the accompanying XYZ file. These can be opened with any computational chemistry graphical viewer, for example Mercury, which can be obtained free of charge at www.ccdc.cam.ac.uk.