

Supplementary material

State-to-state vacuum ultraviolet photodissociation study of CO₂ on the formation of state-correlated CO(*X*¹Σ⁺; *v*) with O(¹D) and O(¹S) photoproducts at 11.95-12.22 eV

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Figure S1: One-dimensional evolution of the MRCI/aug-cc-pV5Z potential energy surfaces of the singlet electronic states of CO_2 along the R_{CO} for several bending angles where the other CO distance is kept fixed to 2.0-2.9 Bohr. The energies are given with respect to the energy of CO_2 ground state at equilibrium.



