

# Correction to Mechanisms and Transition States of 1,3-Dipolar Cycloadditions of Phenyl Azide with Enamines: A Computational Analysis

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## Supporting Information

The Supporting Information incorrectly states the forces on the atoms of the stationary points instead of the  $(x,y,z)$  coordinates. The correct coordinates are available here. This change does not alter any conclusions made in the manuscript.

## ASSOCIATED CONTENT

### Supporting Information

Correct  $(x,y,z)$  coordinates. This material is available free of charge via the Internet at <http://pubs.acs.org>.