

# ***Interactive comment on “Accelerating simulations using Direct Relation Graphs for atmospheric chemistry mechanism reduction” by Zacharias M. Nikolaou et al.***

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The code used to reduce the chemistry is now attached as supplementary material. Also, the name of the code is now included in the title of the revised manuscript.

Please also note the supplement to this comment:

<https://www.geosci-model-dev-discuss.net/gmd-2018-106/gmd-2018-106-AC3-supplement.zip>

Interactive comment on Geosci. Model Dev. Discuss., <https://doi.org/10.5194/gmd-2018-106, 2018>.

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Discussion paper

