

***Interactive comment on “Application of random  
forest regression to the calculation of gas-phase  
chemistry within the GEOS-Chem chemistry  
model v10” by Christoph A. Keller and  
Mat J. Evans***

**Christoph A. Keller and Mat J. Evans**

[christoph.a.keller@nasa.gov](mailto:christoph.a.keller@nasa.gov)

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<https://www.geosci-model-dev-discuss.net/gmd-2018-229/gmd-2018-229-AC1-supplement.pdf>

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Interactive comment on Geosci. Model Dev. Discuss., <https://doi.org/10.5194/gmd-2018-229>,  
2018.