



The atmospheric chemistry box model CAABA/MECCA-4.0

Rolf Sander (1), Andreas Baumgaertner (2), David Cabrera (1), Franziska Frank (4), Sergey Gromov (1,8), Hartwig Harder (1), Vincent Huijnen (3), Patrick Jöckel (4), Vlassis A. Karydis (1), Kyle Niemeyer (5), Andrea Pozzer (1), Martin Schultz (7), Domenico Taraborrelli (7), and Sebastian Tauer (1)

(1) Max Planck Institute for Chemistry, Air Chemistry Department, Mainz, Germany (rolf.sander@mpic.de), (2) Deutsches Zentrum für Luft- und Raumfahrt (DLR), Project Management Agency, 53227 Bonn, Germany, (4) Deutsches Zentrum für Luft- und Raumfahrt (DLR), Institut für Physik der Atmosphäre, Oberpfaffenhofen, 82234 Wessling, Germany, (8) also at: Institute of Global Climate and Ecology (Roshydromet and RAS), Moscow, Russia, (3) Royal Netherlands Meteorological Institute (KNMI), De Bilt, the Netherlands, (5) School of Mechanical, Industrial & Manufacturing Engineering, Oregon State University, USA, (7) IEK-8, Forschungszentrum Jülich, Jülich, Germany

We present version 4.0 of the atmospheric chemistry box model CAABA/MECCA which now includes a number of new features: skeletal mechanism reduction; the MOM chemical mechanism for terpenes and aromatics; an option to include reactions from The Master Chemical Mechanism (MCM) and other chemical mechanisms; updated isotope tagging; new and updated photolysis modules (RADJIMT, JVAL). Another new feature can be used when MECCA is connected to a global model: coexisting multiple chemistry mechanisms (Polymecca/CHEMGLUE). Additional changes have been implemented to make the code more user-friendly and to facilitate the analysis of the model results. Like earlier versions, CAABA/MECCA-4.0 is a community model published under the GNU General Public License.