




# PyBox: An automated box-model generator for atmospheric chemistry and aerosol simulations.

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## Software

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## Summary

Air pollution and climate change are two of the biggest multidisciplinary challenges in society today. The need to understand the chemical and physical processes in the atmosphere that dictate the impacts of both has created a wide range of research platforms. These include numerical models. Volatile organic compounds (VOCs), emitted from both natural and anthropogenic sources, are oxidised in the atmosphere to form lower-volatility species that form organic particulate matter in the atmosphere through gas-to-particle partitioning. Chemical mechanisms, such as the Master Chemical Mechanism (MCM) (Jenkin, Saunders, Wagner, & Pilling, 2003), have been built to hold information of all relevant species and reactions in the atmosphere.

However, it is becoming extremely hard to develop models that can respond to our growing body of knowledge since these mechanisms now treat thousands of species and tens of thousands of reactions. Researchers need to simulate the evolution of each individual species over a range of time-scales and ambient conditions. Manually setting up the relevant ordinary differential equations definitions [ODEs] and associated solvers would present a huge challenge given the number of compounds and equations involved. In addition, it is important to test sensitivity to rate coefficients or evaluate the impact of emerging laboratory data that might identify new reaction pathways or new, improved rate coefficients. Added to this complexity is the need to account for gas-to-particle partitioning of each compound. The atmosphere has varying concentrations of particulate matter that might act as a condensational sink for each compound. Predicting that partitioning requires calculations of molecular properties including saturation vapour pressures. Automation is essential, not least to ensure the gas phase chemistry is accounted for alongside the gas-to-particle partitioning and reproducibility is ensured. This is the driver behind PyBox.

PyBox is a 0-D box model, where all species are homogeneously distributed, built around a chemical mechanism file; a file that represents all the individual chemical reactions taking place starting from a wide range of precursors. By parsing a chemical equation file, obtained from the MCM project, PyBox creates files that account for the gas phase chemistry as well as automatically calculating properties that dictate gas-to-particle partitioning through connection with the UManSysProp informatics suite (Topping et al., 2016). Written in Python, PyBox uses Numba (Lam, Pitrou, & Seibert, 2015) or the Fortran-to-Python-Generator `f2py` (Peterson, 2009) to perform calculations currently within a library of ODE solvers provided by the `Assimulo` package (Andersson, Führer, & Åkesson, 2015).

## Acknowledgements

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## References

Andersson, C., Führer, C., & Åkesson, J. (2015). Assimulo: A unified framework for {ode} solvers. *Mathematics and Computers in Simulation*, 116(0), 26–43. doi:<http://dx.doi.org/10.1016/j.matcom.2015.04.007>

Jenkin, M. E., Saunders, S. M., Wagner, V., & Pilling, M. J. (2003). Protocol for the development of the master chemical mechanism, mcm v3 (part b): Tropospheric degradation of aromatic volatile organic compounds. *Atmospheric Chemistry and Physics*, 3(1), 181–193. doi:[10.5194/acp-3-181-2003](https://doi.org/10.5194/acp-3-181-2003)

Lam, S. K., Pitrou, A., & Seibert, S. (2015). Numba: A llvm-based python jit compiler. In *Proceedings of the second workshop on the llvm compiler infrastructure in hpc, LLVM '15* (pp. 7:1–7:6). New York, NY, USA: ACM. doi:[10.1145/2833157.2833162](https://doi.org/10.1145/2833157.2833162)

Peterson, P. (2009). F2PY: A tool for connecting fortran and python programs. *Int. J. of Computational Science and Engineering*, 4(4), 296–305. doi:<http://dx.doi.org/10.1504/IJCSE.2009.0>

Topping, D., Barley, M., Bane, M. K., Higham, N., Aumont, B., Dingle, N., & McFiggans, G. (2016). UManSysProp v1.0: An online and open-source facility for molecular property prediction and atmospheric aerosol calculations. *Geoscientific Model Development*, 9(2), 899–914. doi:[10.5194/gmd-9-899-2016](https://doi.org/10.5194/gmd-9-899-2016)