




ACEInhibPKPD: An open-source MATLAB app for a pharmacokinetic/pharmacodynamic model of ACE inhibition

Ashlee N. Ford Versypt¹, Grace K. Harrell¹, and Alexandra N. McPeak¹

DOI: [10.21105/joss.00340](https://doi.org/10.21105/joss.00340)

¹ School of Chemical Engineering, Oklahoma State University

Software

- [Review](#) 
- [Repository](#) 
- [Archive](#) 

Licence

Authors of JOSS papers retain copyright and release the work under a Creative Commons Attribution 4.0 International License ([CC-BY](#)).

Summary

This software features an MATLAB app install file and the source code for an interactive computer simulation packaged as a MATLAB app that can be used to design the best dosage of pharmaceuticals (ACE inhibitors) for reducing high blood pressure. The app uses concepts from chemical engineering and tools from mathematics, pharmacology, and computational science to describe and solve the dynamics of the chemical reactions in the human body involved in the absorption, metabolism, and excretion of the drug and how the blood pressure-regulating hormone angiotensin II is affected by the drug concentration as a function of time. The scientific significance is that the model solved in the software fills the gap in pharmacokinetic/pharmacodynamic models for ACE inhibitors in impaired renal function, given the prevalence of using this class of pharmaceuticals for treating chronic kidney disease in addition to high blood pressure. The model used in this software is described in detail in (Ford Versypt, Harrell, and McPeak 2017). Use of the software in the context of engineering education is described in (Harrell, McPeak, and Ford Versypt 2017).

References

- Ford Versypt, A. N., G. K. Harrell, and A. N. McPeak. 2017. “A pharmacokinetic/pharmacodynamic model of ACE inhibition of the renin-angiotensin system for normal and impaired renal function.” *Computers & Chemical Engineering* 104: 311–22. doi:10.1016/j.compchemeng.2017.03.027.
- Harrell, G. K., A. N. McPeak, and A. N. Ford Versypt. 2017. “A pharmacokinetic simulation-based module to introduce mass balances and chemical engineering design concepts to engineering freshmen.” *Proceedings of the ASEE Annual Conference* Columbus, OH.

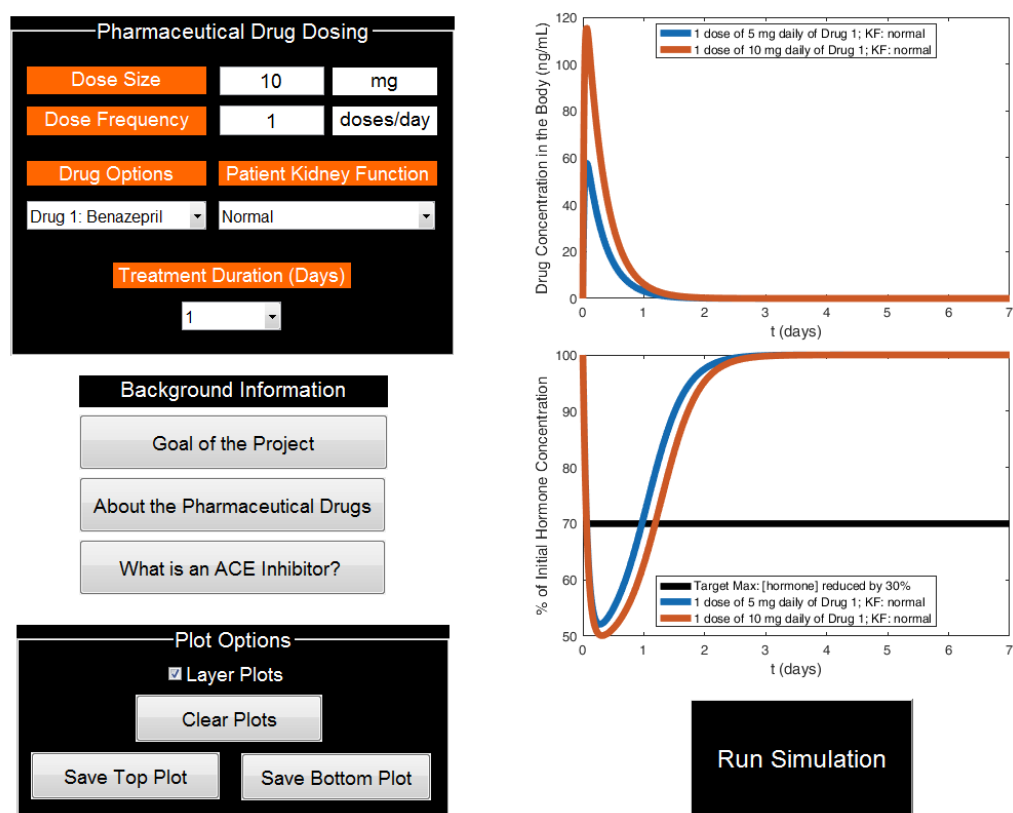


Figure 1: GUI screenshot