

Correction to Mitigating Heterocycle Metabolism in Drug Discovery

David J. St. Jean Jr. and Christopher Fotsch*

Journal of Medicinal Chemistry 2012, 55, 6002–6020. DOI: 10.1021/jm300343m

Page 6017. Reference 24 should read as follows: Unless noted otherwise, cLogP values were calculated using the PCModels software, version 4.9.2, from Daylight, which is developed and supported by BioByte, Inc., Claremont, CA. Also, unless noted otherwise, the term “cLogD_{7.4}” refers to log *D*_{7.4} values calculated using ACD/Labs software, version 10.