Accepted Manuscript

Title: Prediction tool for loading, isocratic elution, gradient elution and scaling up of ion exchange chromatography of proteins

Authors: Wojciech Kazimierz Marek, Dominik Sauer, Astrid Dürauer, Alois Jungbauer, Wojciech Piątkowski, Dorota Antos

PII: S0021-9673(18)30812-4

DOI: https://doi.org/10.1016/j.chroma.2018.06.057

Reference: CHROMA 359501

To appear in: Journal of Chromatography A

Received date: 6-3-2018 Revised date: 20-6-2018 Accepted date: 22-6-2018

Please cite this article as: Marek WK, Sauer D, Dürauer A, Jungbauer A, Piatkowski W, Antos D, Prediction tool for loading, isocratic elution, gradient elution and scaling up of ion exchange chromatography of proteins, *Journal of Chromatography A* (2018), https://doi.org/10.1016/j.chroma.2018.06.057

This is a PDF file of an unedited manuscript that has been accepted for publication. As a service to our customers we are providing this early version of the manuscript. The manuscript will undergo copyediting, typesetting, and review of the resulting proof before it is published in its final form. Please note that during the production process errors may be discovered which could affect the content, and all legal disclaimers that apply to the journal pertain.



ACCEPTED MANUSCRIPT

Prediction tool for loading, isocratic elution, gradient elution and scaling up of ion exchange chromatography of proteins

Wojciech Kazimierz Marek a ,*, Dominik Sauer b , Astrid Dürauer b,c , Alois Jungbauer b ,c, Wojciech Piątkowski a , Dorota Antos a

^aDepartment of Chemical and Process Engineering, Powstańców Warszawy Ave. 6, 35-959 Rzeszów, Poland

^bAustrian Centre of Industrial Biotechnology, Muthgasse 11, 1190 Vienna, Austria

*Corresponding author; E-mail address: wmarek@prz.edu.pl (W.K. Marek), tel.: +48 177432020; fax: +48 178543655

Highlights

- Mathematical tool for design and scaling up of protein chromatography is suggested
- Moment analysis was used for determining kinetic parameters of mass transfer
- Two model ion-exchange chromatography systems were selected for the study
- The design procedure was successfully verified over a wide space of the operating conditions

Abstract

An efficient mathematical tool for the design and scaling up of protein chromatography is suggested, in which the model parameters can be determined quickly over a wide operating space without large material investments. The design method is based on mathematical modelling of column dynamics and moment analysis. The accuracy of the dynamic models that are most frequently used for simulations of chromatographic processes is analyzed, and possible errors that can be generated using the moment analysis are indicated. The so-called transport dispersive model was eventually employed for the process simulations. The model was modified to account for the protein dispersion in void volumes of chromatographic systems. The manner of the model calibration was

^cDepartment of Biotechnology, University of Natural Resources and Life Sciences Vienna, Muthgasse 18, 1190 Vienna, Austria

Download English Version:

https://daneshyari.com/en/article/7607504

Download Persian Version:

https://daneshyari.com/article/7607504

<u>Daneshyari.com</u>