

Open up a new world of opportunities for your chromatographic processes





Ypso-lonic[®] is a software solution for designing chromatographic processes. It maximizes the value of experimental data and makes predictive simulation a game changer.

WHY "IONIC"?

Does pH matter in your separation process? Do you work with mineral ions? With organic buffers? With proteins, peptides, or oligonucleotides whose retention behaviour is dramatically impacted by their charge?

If yes, properly dealing with ions and thus using IONIC is a must!

If no, that's also fine, this will just make the calculations easier!

IONIC KEY STRENGTH

A unique feature of lonic is the way it deals with ions: it offers the possibility to account for solution equilibria (e.g., acid-base or complexation) in combination with a rigorous description of interaction mechanisms with the solid media.

A WIDE RANGE OF APPLICATIONS

Ionic can be used for various types of chromatography:

Ion exchange, adsorption, affinity, hydrophobic interactions, size exclusion, ion exclusion, chelation, etc

... from the simplest to the most complex molecules:

mineral ions, sugars, organic acids, amino acids, oligonucleotides, proteins, monoclonal antibodies, etc

• ... with any kind of process configuration:

single-column or multicolumn, isocratic or gradient, backflush, with or without recycling, etc



Two complementary means to address your challenges: experimental data (X) and predictive simulations (P)

MAXIMIZE THE VALUE OF YOUR DATA

 collect and structure experimental data to ensure data completeness and reusability: our YIF Data format associates the experimental protocol (column dimensions, raw materials composition, flowrates, etc) with clearly labelled measured results (UV, pH, conductivity, species concentration, etc)

 visualize experimental results thanks to flexible graph tools – possibly overlaying online measurements with offline fraction analyses

 easily extract relevant information from experimental data: check mass balance, calculate purity of mock pools, estimate productivity and eluent consumption, etc.



EMBRACE THE POWER OF MECHANISTIC SIMULATION

 enjoy the predictive capabilities of the most advanced mechanistic models in a user-friendly interface

• simulate outlet concentration profiles as well as signals like pH and conductivity

 apply automatic post-treatments to calculate fractograms, mock pools, performance indicators (purity, yield, productivity, eluent consumption)

easily compare experiments with simulations

 investigate the impact of key operating parameters (loading, flowrate, pH, salt concentration, gradient slope, column dimensions, etc) and predict behaviours in non-explored regions

 use our Buffer preparation tool to get recipes based on specified targets (e.g., target pH and concentration)





GPX[®] Concept - Leveraging 3 complementary means to address your challenges.

Guess – the ability of the experts to propose, anticipate, imagine... some routes or performances

Predictive Simulation – for a better understanding of your process, to anticipate issues and make rational decisions

eXperimental data – the hard foundation of the project, which needs to be well targeted, structured and analyzed.

Enjoy working in a virtual lab

Select your raw materials, prepare buffers, pack columns, define separation protocols... as if you were in your own lab! Then, either store your experimental measurements or click Run and predict the results.

Save time thanks to automatic post-treatments

Focus on your field of expertise and let the software perform the tedious work of data treatment, be it for experimental or simulated data. Boost your creativity

Simulation significantly broadens the range of possibilities – explore new process configurations in a few clicks.

Get the most of your experimental data

Reduce experimental workload

Maximize process performances

EXAMPLES OF USE

- Transform data into knowledge
- Quantify the impact of key operating parameters
- Train your team on chromatographic processes
- Optimize technical performances
- Identify the design space
- Use simulation as a troubleshooting tool



WHAT'S MORE?

Ypso-lonic[®] is the chromatography unit operation App within **Ypso-Proxima[®]**, our collaborative software suite for chemical and bioprocess development, optimization, and evaluation. Ionic is compatible with the other Apps of Proxima – making Ionic results directly exploitable to perform an economic evaluation.

Contact us for a demo!

contact@ypso-facto.com | www.ypsofacto.com | +33 3 55 96 16 50

© 2021 YPSO-FACTO, YFI SAS. The content of this brochure is owned and copyrighted by YFI SAS and its subsidiaries. All rights reserved. Except as stated herein, none of the material included in this brochure may be copied, reproduced, distributed, republished, downloaded, displayed, posted or transmitted in any form or by any means including, but not limited to, electronic, mechanical, photocopying, recording or otherwise, except with the prior written consent of Ypso-Facto. Disclaimers: The information contained in this brochure are provided "as is", for informational purposes only, without any representation or warranty of accuracy or completeness. Ypso-Facto will not be liable for any damage or injury from access or reliance upon any information contained in this brochure.

