

Weronika Puchała^{1,*}, Michał Burdukiewicz², Michał Kistowski¹, Katarzyna Dabrowa-Dabrowska¹, Aleksandra E. Badaczewska-Dawid³,
Dominik Cysewski¹, Michał Dadlez¹
*puchala.weronika@gmail.com

Introduction

Main functionality


Start

Woods plot

Coverage

Sequence data

Report



Select parameters for the plot.

☐ Theoretical calculations

Choose values type:

☒ relative
 ☐ absolute

Comparison plot parameters:

Choose time parameters:

IN

0.001

CHOSEN

1

OUT

1440

Choose states for comparison:

☒ gg_Nucb2_CaCl2
 ☒ gg_Nucb2_EDTA

Adjust colors

gg_Nucb2_CaCl2 color

#ED90A4

gg_Nucb2_EDTA color

#00C1B2

Woods plot parameters:

State 1

gg_Nucb2_CaCl2

State 2

gg_Nucb2_EDTA

Confidence limit 1:

98%

Confidence limit 2:

99%

Adjust plot:

Choose y range for comparison plot:

2

-2

-1.6

-1.2

-0.8

-0.4

0

0.4

0.8

1.2

1.6

2

Choose y range for Woods plot:

2

-2

-1.6

-1.2

-0.8

-0.4

0

0.4

0.8

1.2

1.6

2

Choose x range for both plots:

1

407

Adjust labels

Comparison plot

Data

Fraction exchanged in state comparison in 1 min

gg_Nucb2_CaCl2

gg_Nucb2_EDTA

Save chart (.svg)

Woods plot

Data

Delta Fraction exchanged in 1 min between gg_Nucb2_CaCl2 and gg_Nucb2_EDTA

Confidence interval 98% - 0.0343

Confidence interval 99% - 0.0489

Save chart (.svg)

Calculations and uncertainty

$$D = \frac{D_t - D_0}{D_{100} - D_0} \quad (1)$$
$$diff = D_1 - D_2 \quad (2)$$
$$u_c(y) = \sqrt{\sum_k \left[\frac{\partial y}{\partial x_k} u(x_k) \right]^2} \quad (3)$$

Availability



Protein name
db_Nucb2

Reconstructed sequence

```

XXXXXXXXXXXXXXXXXXVPIDDKTKYKGEGHVEGEKINPDTGLYYDEYLQGVIDVLEDKHFKREKLQDTAIEEKISOKLSRELDLVSHVIRTLDELKRGVARI.RMLIKAKMDSVQDTIDHQA.LKQFEHLNHQNPOTFEPKDLMLTKAATSDLENYOKTRHEEFKYYKXXXX
XXXXXXXXXXXXLDEKRRKEESFKGEKXXXXXXXXXXXXXXXXXXXXKVEEADGLDPNEFDPKFFFKLHVNWRDFLDELEAxFKTELEKYDPKNEEDDYMVEEERLXXXXHVNVEVDINKDRLVLEELFRAKTEKFEFLPSDWLDDQQLFTEDELKFESHSIQDEDLR
KKAELQKQKEELGRHDLQADEDELQQVYKQNEKRLQANPPAPGAPGLK

```

Name	Value
Length	407
Coverage	97.22%
Cys	0

Correct sequence length:

If C-terminal of is not covered by peptides, enter to correct position.

Sequence length from the file is 427

Hydro-

- ☒ Hydrophobic
- ☒ Hydrophilic

Amino acid composition

Amino Acid	Count	Charge
P	10	Neutral
A	12	Neutral
G	15	Neutral
S	10	Neutral
T	10	Neutral
C	0	Neutral
I	38	Neutral
L	12	Neutral
V	10	Neutral
M	8	Neutral
F	8	Neutral
Y	5	Neutral
N	5	Neutral
D	15	Neutral
E	30	Negative
K	55	Positive
R	30	Positive
Q	15	Neutral
H	10	Neutral

[Save chart \(.svg\)](#)

Choose state:

- gg_Nutrio_CoQ2
- gg_Nutrio_EDTA

Choose range:

Peptide Coverage
Data

Peptide coverage

Relative intensity

Position

[Save chart \(.svg\)](#)

Choose state:

- gg_Nutrio_CoQ2
- gg_Nutrio_EDTA

Choose range:

Position Frequency
Data

Position frequency

Relative frequency

Position

Average frequency: 0.25

[Save chart \(.svg\)](#)

log(time) [min]	Average error (first state)	Average error (second state)
1	0.030	0.078
10	0.008	0.010
100	0.007	0.008
1000	0.007	0.008

Choose items for report:

☒ Region Frequency
 ☒ Region Coverage
 ☒ Comparison Rate
 ☒ Theoretical Comparison Plot
 ☒ Woods Plot
 ☒ Theoretical Woods Plot

☐ Region Frequency Date
 ☐ Region Coverage Date
 ☐ Comparison Plot Date
 ☐ Theoretical Comparison Plot Date
 ☐ Woods Plot Date
 ☐ Theoretical Woods Plot Date

Elements chosen for report have the same parameters as chosen in panel **4**, with cell **1** and cell **2**. Adjust parameters for plots as needed in this report.

Create report



- 1 Start2Fold: a database of hydrogen/deuterium exchange data on protein folding and stability. Pancsa R, Varadi M, Tompa P, Vranken WF. *Nucleic Acids Res.* 2016 Jan 4;44(D1):D429-34. doi: 10.1093/nar/gkv1185. Epub 2015 Nov 17
- 2 HDX workbook: software for the analysis of H/D exchange MS data. Pascal BD, Willis S, Lauer JL, Landgraf RR, West GM, Marciano D, Novick S, Goswami D, Chalmers MJ, Griffin PR. *J Am Soc Mass Spectrom.* 2012 Sep;23(9):1512-21. doi: 10.1007/s13361-012-0419-6. Epub 2012 Jun 13.
- 3 MEMHDX: an interactive tool to expedite the statistical validation and visualization of large HDX-MS datasets. Hourdel V, Volant S, O'Brien DP, Chenal A, Chamot-Rooke J, Dillies MA, Brier S. *Bioinformatics.* 2016 Nov 15;32(22):3413-3419. Epub 2016 Jul 13
- 4 Deuterios: software for rapid analysis and visualization of data from differential hydrogen deuterium exchange-mass spectrometry. Andy M C Lau, Zainab Ahdash, Chloe Martens, Argyris Politis, *Bioinformatics*, btz022, <https://doi.org/10.1093/bioinformatics/btz022>
- 5 ISO, Guide to the Expression of Uncertainty in Measurement (International Organization for Standardization, Geneva, Switzerland, 1993)
- 6 The Utility of Hydrogen/Deuterium Exchange Mass Spectrometry in Biopharmaceutical Comparability Studies. Houde, D.et al.(2011). *Journal of pharmaceutical sciences*,100(6), 2071–2086