HADDOCK 2.4

@Bonvinlab

WELCOME TO THE UTRECHT BIOMOLECULAR INTERACTION WEB PORTAL >>

HADDOCK server status for job "cford_HLAcsp-16"

Status: FINISHED

Your HADDOCK run has successfully completed. The complete run can be downloaded as a gzipped tar file here. The file containing your docking parameters is here.

Please cite the following paper in your work:

G.C.P van Zundert, J.P.G.L.M. Rodrigues, M. Trellet, C. Schmitz, P.L. Kastritis, E. Karaca, A.S.J. Melquiond, M. van Dijk, S.J. de Vries and A.M.J.J. Bonvin (2016). "The HADDOCK2.2 webserver: User-friendly integrative modeling of biomolecular complexes

J. Mol. Biol., 428, 720-725 (2015).

and add the following acknowledgment:

The FP7 WeNMR (project# 261572), H2020 West-Life (project# 675858), the EOSC-hub (project# 777536) and the EGI-ACE (project# 101017567) European e-Infrastructure projects are acknowledged for the use of their web portals, which make use of the EGI infrastructure with the dedicated support of CESNET-MCC, INFN-PADOVA-STACK, INFN-LNL-2, NCG-INGRID-PT, TW-NCHC, CESGA, IFCA-LCG2, UA-BITP, SURFsara and NIKHEF, and the additional support of the national GRID Initiatives of Belgium, France, Italy, Germany, the Netherlands, Poland, Portugal, Spain, UK, Taiwan and the US Open Science Grid.

How would you rate your experience with our portal?









Questions / feedback ? ask.bioexcel.eu

Do check up the HADDOCK best practice guide! There you can learn more about which settings are best used in which scenario and use HADDOCK in its full potential!

In the aim to improve our new web portal, we would really appreciate 2 min of your time to complete a short survey here! Thanks!

Post-processing: SUCCESS

Summary

HADDOCK clustered 171 structures in 14 cluster(s), which represents 85 % of the water-refined models HADDOCK generated. Note that currently the maximum number of models considered for clustering is 200.

The statistics of the top 10 clusters are shown below. The top cluster is the most reliable according to HADDOCK. Its Z-score indicates how many standard deviations from the average this cluster is located in terms of score (the more negative the better).

A $\underline{\text{graphical representation}}$ of the results is also provided at the bottom of the page.

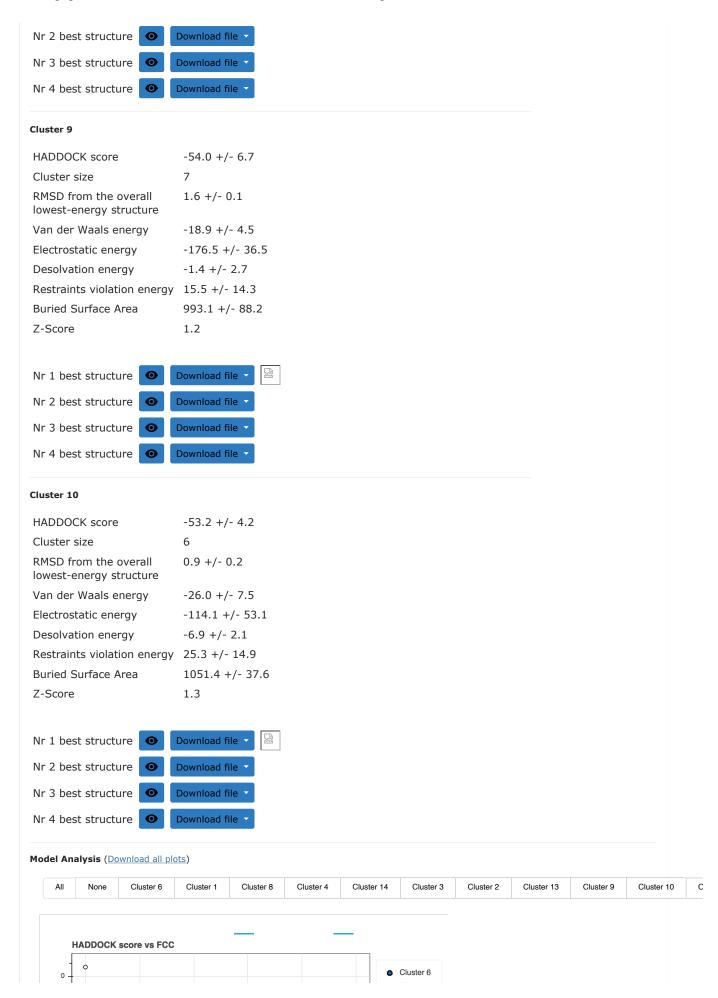
You can also download all cluster files (best 4 of the top 10 cluster(s)).

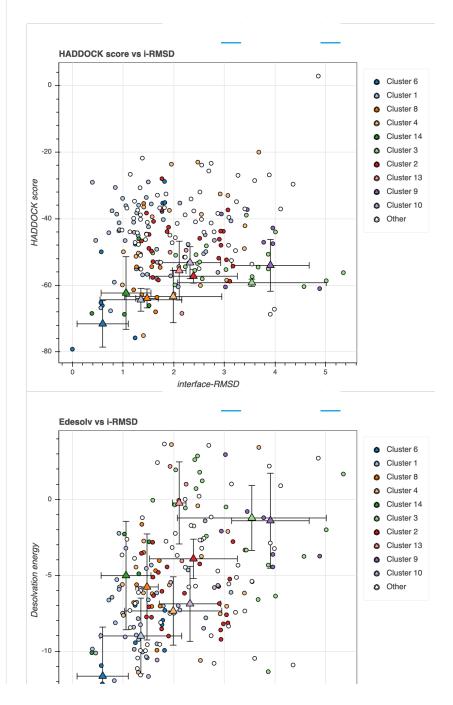
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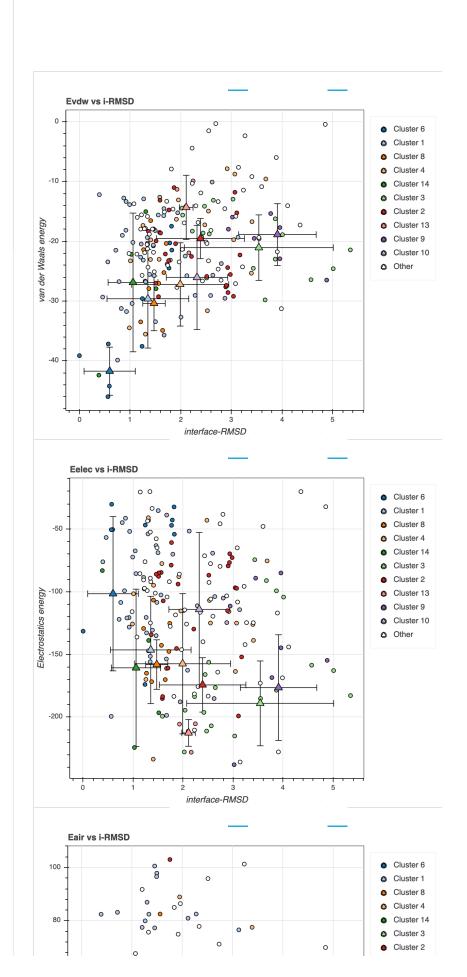
Cluster 6		
HADDOCK score	-71.6 +/- 6.0	
Cluster size	10	
RMSD from the overall owest-energy structure	0.4 +/- 0.3	
Van der Waals energy	-41.7 +/- 3.5	
Electrostatic energy	-101.6 +/- 53.2	
Desolvation energy	-11.7 +/- 2.8	
Restraints violation energy	21.4 +/- 16.6	
Buried Surface Area	1237.7 +/- 30.2	
Z-Score	-2.1	
Nr 1 best structure	Download file 🔻	
Nr 2 best structure	Download file 🔻	
Nr 3 best structure	Download file 🔻	
Nr 4 best structure	Download file 🔻	
Cluster 1		
HADDOCK score	-64.4 +/- 3.0	
Cluster size	39	
RMSD from the overall lowest-energy structure	0.6 +/- 0.3	
Van der Waals energy	-29.6 +/- 7.1	
Electrostatic energy	-146.4 +/- 37.0	
Desolvation energy	-9.0 +/- 2.2	
Restraints violation energy	35.5 +/- 25.1	
Buried Surface Area	1149.8 +/- 105.3	
Z-Score	-0.7	
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Nr 2 best structure	Download file 🔻	
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Nr 4 best structure	Download file 🔻	
Cluster 8		
HADDOCK score	-63.9 +/- 2.5	
Cluster size	9	
RMSD from the overall lowest-energy structure	0.7 +/- 0.1	
Van der Waals energy	-30.4 +/- 4.0	
Electrostatic energy	-158.1 +/- 17.1	
Desolvation energy	-5.8 +/- 3.0	
Restraints violation energy	38.3 +/- 17.3	
Buried Surface Area	1159.3 +/- 89.9	

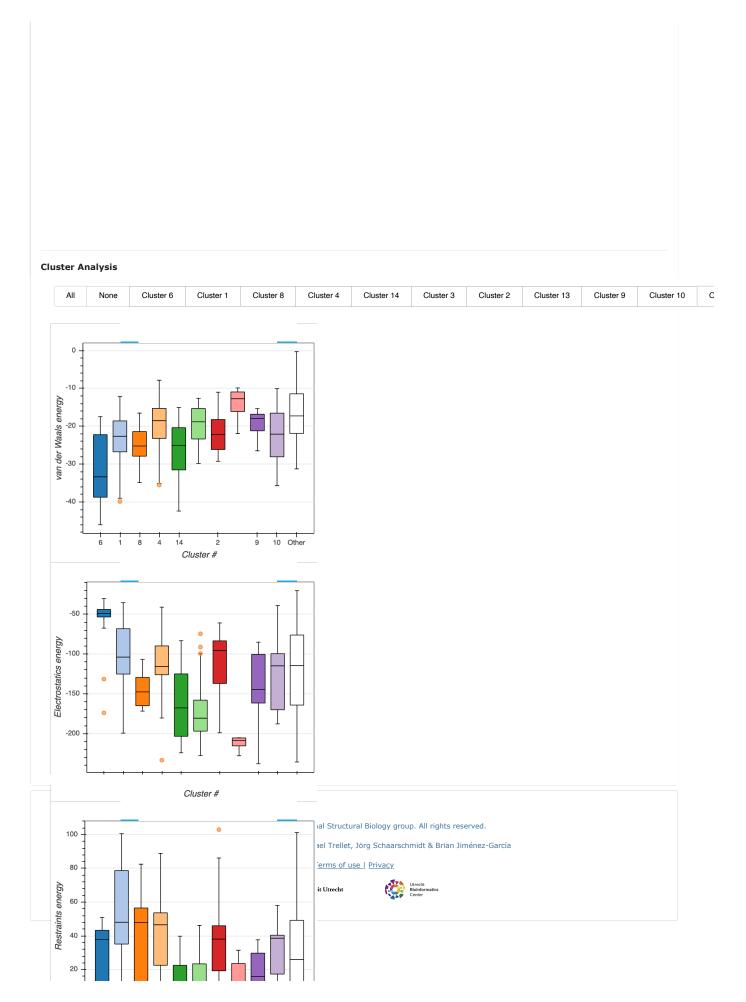
Z-Score	-0.6	
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Nr 3 best structure	Download file 🔻	
Nr 4 best structure	Download file 🔻	
Cluster 4		
HADDOCK score	-63.4 +/- 6.8	
Cluster size	18	
RMSD from the overall lowest-energy structure	0.7 +/- 0.2	
Van der Waals energy	-27.2 +/- 6.0	
Electrostatic energy	-157.5 +/- 48.4	
Desolvation energy	-7.4 +/- 2.0	
Restraints violation ene Buried Surface Area	-	
Z-Score	1096.2 +/- 102.2 -0.5	
2 30010	0.5	
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Nr 4 best structure	Download file •	
Cluster 14		
HADDOCK score	-62.3 +/- 9.5	
Cluster size	4	
RMSD from the overall lowest-energy structure	0.6 +/- 0.1	
Van der Waals energy	-26.9 +/- 10.1	
Electrostatic energy	-160.7 +/- 54.3	
Desolvation energy	-5.0 +/- 3.1	
Restraints violation ene Buried Surface Area	rgy 17.4 +/- 13.6 1169.3 +/- 77.1	
Z-Score	-0.3	
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Nr 2 best structure	Download file 🔻	
Nr 3 best structure	Download file •	
Nr 4 best structure		
Cluster 3		
	E0.2 . / 0.0	
HADDOCK score	-59.2 +/- 0.9	

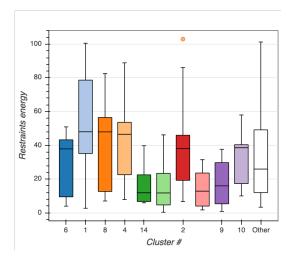
Cluster size	20
RMSD from the overall lowest-energy structure	1.4 +/- 0.2
Van der Waals energy	-21.1 +/- 4.8
Electrostatic energy	-189.1 +/- 29.3
Desolvation energy	-1.2 +/- 1.9
Restraints violation energy	8.5 +/- 4.1
Buried Surface Area	979.5 +/- 49.5
Z-Score	0.2
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Cluster 2	
HADDOCK score	-57.3 +/- 1.8
Cluster size	24
RMSD from the overall lowest-energy structure	1.5 +/- 0.0
Van der Waals energy	-19.5 +/- 2.9
Electrostatic energy	-174.3 +/- 18.8
Desolvation energy	-3.9 +/- 1.1
Restraints violation energy	
Buried Surface Area	1037.6 +/- 22.5
Z-Score	0.6
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Nr 2 best structure	Download file 🔻
Nr 3 best structure	Download file •
Nr 4 best structure	Download file •
Cluster 13	
	EE 6 1 / 7 7
HADDOCK score	-55.6 +/- 7.7
Cluster size	4
RMSD from the overall lowest-energy structure	1.0 +/- 0.1
Van der Waals energy	-14.3 +/- 4.7
Electrostatic energy	-212.6 +/- 9.2
Desolvation energy	-0.2 +/- 2.3
Restraints violation energy	
Buried Surface Area	937.5 +/- 66.3
Z-Score	0.9
Nr 1 best structure	Download file











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