

HADDOCK 2.4

@Bonvinlab

WELCOME TO THE UTRECHT BIOMOLECULAR INTERACTION WEB PORTAL >>

HADDOCK server status for job "cford_TCR-csp-1"

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. Kastiris, E. Karaca, A.S.J. Melquiand, 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:

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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 **195** structures in **11** cluster(s), which represents **97 %** 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 [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)).

Cluster 1

HADDOCK score	-123.7 +/- 6.3
Cluster size	83
RMSD from the overall lowest-energy structure	0.4 +/- 0.3
Van der Waals energy	-88.4 +/- 0.7
Electrostatic energy	-229.4 +/- 35.8
Desolvation energy	-1.5 +/- 1.9
Restraints violation energy	121.5 +/- 39.6
Buried Surface Area	2130.2 +/- 41.6
Z-Score	-1.5

- Nr 1 best structure

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Cluster 6

HADDOCK score	-119.7 +/- 4.8
Cluster size	10
RMSD from the overall lowest-energy structure	0.7 +/- 0.2
Van der Waals energy	-87.7 +/- 5.4
Electrostatic energy	-266.4 +/- 54.5
Desolvation energy	-1.5 +/- 3.1
Restraints violation energy	227.8 +/- 46.2
Buried Surface Area	2230.3 +/- 65.5
Z-Score	-1.2

- Nr 1 best structure

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Cluster 2


HADDOCK score	-114.0 +/- 4.4
Cluster size	37
RMSD from the overall lowest-energy structure	8.5 +/- 0.2
Van der Waals energy	-88.3 +/- 2.2
Electrostatic energy	-193.8 +/- 28.1
Desolvation energy	-7.7 +/- 2.6
Restraints violation energy	206.9 +/- 52.3
Buried Surface Area	2212.7 +/- 22.5


Z-Score -0.9

- Nr 1 best structure




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- Nr 2 best structure



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- Nr 3 best structure



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Cluster 10

HADDOCK score -113.6 +/- 7.2

Cluster size 5

RMSD from the overall lowest-energy structure 8.7 +/- 0.1

Van der Waals energy -82.5 +/- 4.0

Electrostatic energy -231.8 +/- 23.6


Desolvation energy -7.3 +/- 3.4

Restraints violation energy 226.1 +/- 66.0


Buried Surface Area 2217.6 +/- 57.5


Z-Score -0.9

- Nr 1 best structure




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- Nr 2 best structure



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- Nr 3 best structure



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Cluster 3

HADDOCK score -96.5 +/- 3.6

Cluster size 14

RMSD from the overall lowest-energy structure 9.1 +/- 0.2

Van der Waals energy -77.1 +/- 5.3

Electrostatic energy -132.1 +/- 6.0


Desolvation energy -14.7 +/- 1.3

Restraints violation energy 217.0 +/- 31.4


Buried Surface Area 2120.7 +/- 70.2


Z-Score 0.1

- Nr 1 best structure




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- Nr 2 best structure



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- Nr 3 best structure



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Cluster 9



HADDOCK score -92.3 +/- 6.6



Cluster size	6
RMSD from the overall lowest-energy structure	3.6 +/- 0.3
Van der Waals energy	-65.0 +/- 4.1
Electrostatic energy	-245.8 +/- 21.4
Desolvation energy	-4.5 +/- 1.3
Restraints violation energy	263.7 +/- 58.7
Buried Surface Area	1878.0 +/- 23.4
Z-Score	0.3


- Nr 1 best structure 

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Cluster 5



HADDOCK score	-86.5 +/- 6.5
Cluster size	10
RMSD from the overall lowest-energy structure	12.0 +/- 0.1
Van der Waals energy	-61.7 +/- 9.6
Electrostatic energy	-230.8 +/- 29.5
Desolvation energy	-4.7 +/- 2.5
Restraints violation energy	260.1 +/- 34.4
Buried Surface Area	1873.5 +/- 111.0
Z-Score	0.7


- Nr 1 best structure 

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Cluster 7

HADDOCK score	-84.2 +/- 1.7
Cluster size	7
RMSD from the overall lowest-energy structure	18.9 +/- 1.2
Van der Waals energy	-60.4 +/- 3.0
Electrostatic energy	-173.3 +/- 9.4
Desolvation energy	-6.9 +/- 2.5
Restraints violation energy	177.2 +/- 28.2
Buried Surface Area	1720.7 +/- 95.2
Z-Score	0.8

- Nr 1 best structure 

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Cluster 4

HADDOCK score	-81.2 +/- 5.3
Cluster size	13
RMSD from the overall lowest-energy structure	9.8 +/- 0.4
Van der Waals energy	-58.6 +/- 4.4
Electrostatic energy	-179.3 +/- 3.0
Desolvation energy	-5.9 +/- 3.9
Restraints violation energy	192.2 +/- 37.6
Buried Surface Area	1918.3 +/- 58.7
Z-Score	1.0

- Nr 1 best structure

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Cluster 8

HADDOCK score	-69.9 +/- 12.4
Cluster size	6
RMSD from the overall lowest-energy structure	7.2 +/- 1.2
Van der Waals energy	-55.1 +/- 7.7
Electrostatic energy	-162.9 +/- 59.5
Desolvation energy	-11.8 +/- 3.3
Restraints violation energy	295.7 +/- 65.1
Buried Surface Area	1832.1 +/- 59.7
Z-Score	1.6

- Nr 1 best structure

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- Nr 2 best structure

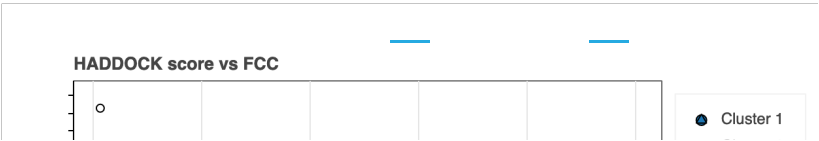
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- Nr 3 best structure

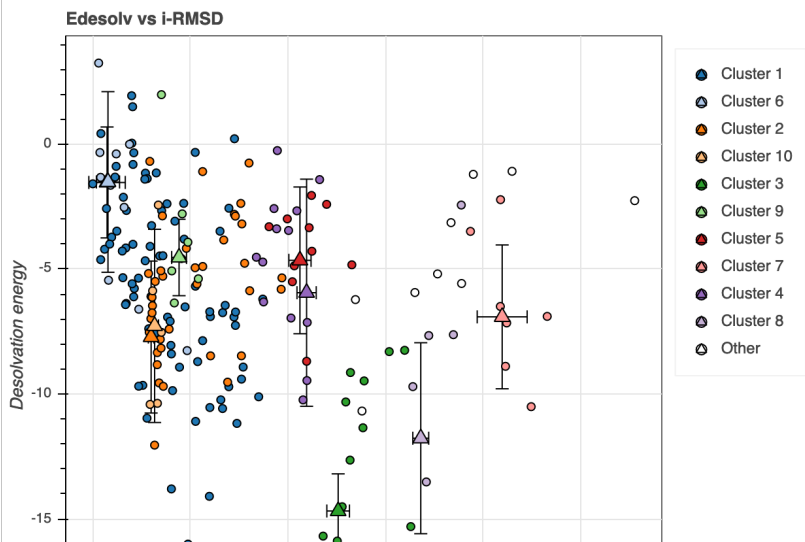
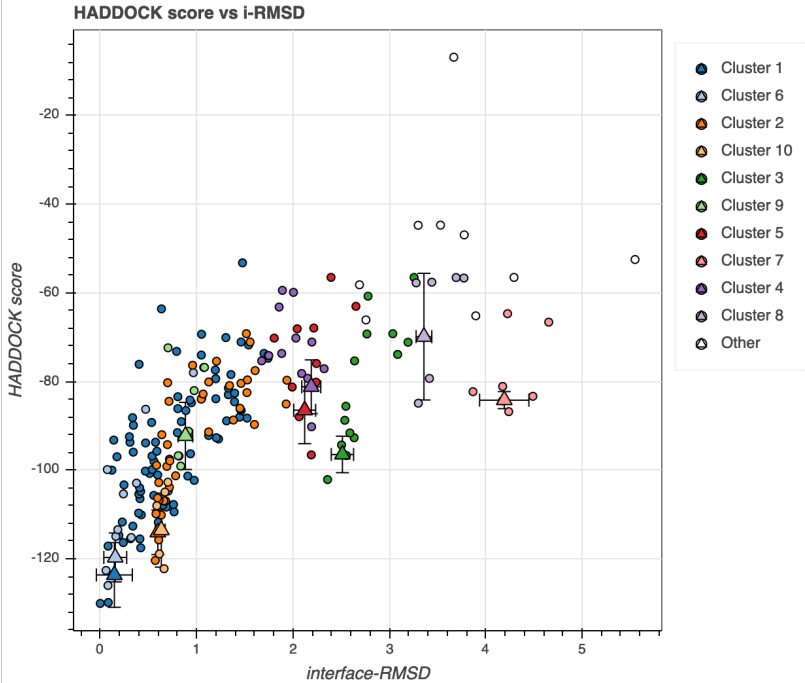
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- Nr 4 best structure

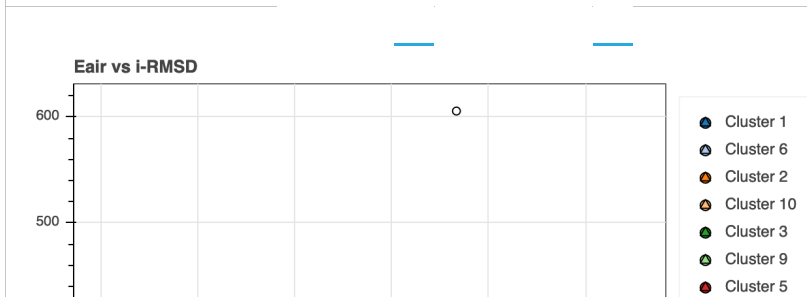
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Model Analysis [\(Download all plots\)](#)

All	None	Cluster 1	Cluster 6	Cluster 2	Cluster 10	Cluster 3	Cluster 9	Cluster 5	Cluster 7	Cluster 4	Cluster 8	C
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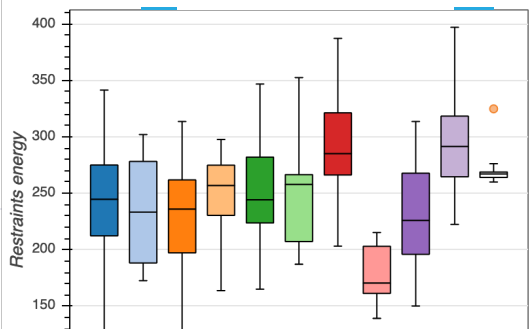
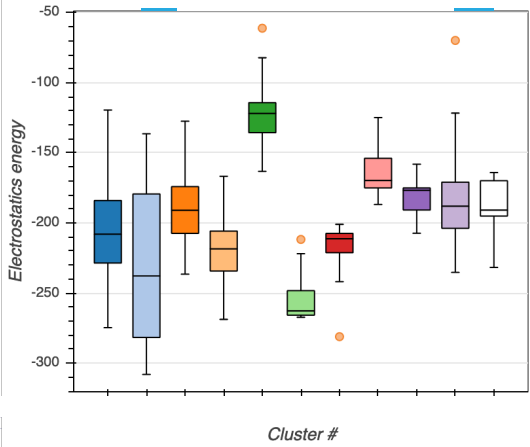
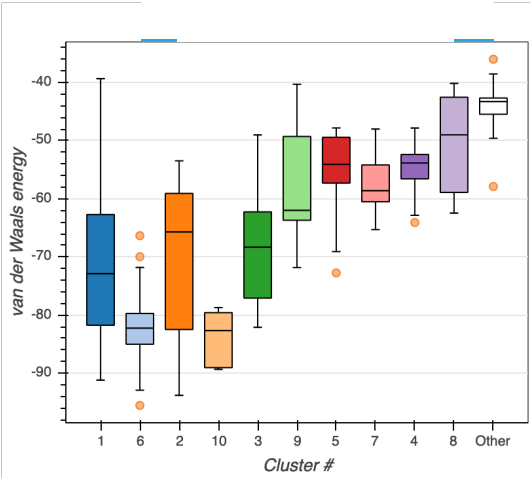






Cluster Analysis

All	None	Cluster 1	Cluster 6	Cluster 2	Cluster 10	Cluster 3	Cluster 9	Cluster 5	Cluster 7	Cluster 4	Cluster 8	O
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