

Analysis output: all-atom contacts and geometry for CDH23_1628-1742H.pdb

Summary statistics

Contacts	Clashscore, all atoms:	39.36		8 th percentile* (N=1784, all resolutions)
	Clashscore is the number of serious steric overlaps (> 0.4 Å) per 1000 atoms.			
Protein Geometry	Poor rotamers	2	1.96%	Goal: <1%
	Ramachandran outliers	5	4.42%	Goal: <0.05%
	Ramachandran favored	90	79.65%	Goal: >98%
	MolProbity score [^]	3.04		21 st percentile* (N=27675, 0Å - 99Å)
	Cβ deviations >0.25Å	0	0.00%	Goal: 0
	Bad backbone bonds:	8 / 903	0.89%	Goal: 0%
	Bad backbone angles:	10 / 1240	0.81%	Goal: <0.1%

In the two column results, the left column gives the raw count, right column gives the percentage.

By adding H to this model and allowing Asn/Gln/His flips, we could *automatically* improve your clashscore by 1.71 points.

Multi-criterion visualizations

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^{* 100&}lt;sup>th</sup> percentile is the best among structures of comparable resolution; 0th percentile is the worst. For clashscore the comparative set of structures was selected in 2004, for MolProbity score in 2006.

[^] MolProbity score combines the clashscore, rotamer, and Ramachandran evaluations into a single score, normalized to be on the same scale as X-ray resolution.



View (119 Kb)

Single-criterion visualizations

- Clash list (2.2 Kb): View
- Ramachandran plot kinemage (412 Kb): View in KiNG | Download
- Ramachandran plot PDF (1.7 Mb): View
- Cβ deviation scatter plot (17 Kb): View in KiNG | Download

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