

## Analysis output: all-atom contacts and geometry for MYO7A\_60-686H.pdb

## **Summary statistics**

Contacts	Clashscore, all atoms:	71.86		1 <sup>st</sup> percentile* (N=1784, all resolutions)
	Clashscore is the number of serious steric overlaps (> 0.4 Å) per 1000 atoms.			
Protein Geometry	Poor rotamers	1	0.18%	Goal: <1%
	Ramachandran outliers	13	2.08%	Goal: <0.05%
	Ramachandran favored	590	94.40%	Goal: >98%
	MolProbity score <sup>^</sup>	2.71		36 <sup>th</sup> percentile* (N=27675, 0Å - 99Å)
	Cβ deviations >0.25Å	9	1.52%	Goal: 0
	Bad backbone bonds:	53 / 5181	1.02%	Goal: 0%
	Bad backbone angles:	108 / 6990	1.55%	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 1.09 points.

## Multi-criterion visualizations

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

<sup>^</sup> 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 (639 Kb)

## Single-criterion visualizations

- Clash list (14 Kb): View
- Ramachandran plot kinemage (442 Kb): View in KiNG | Download
- Ramachandran plot PDF (1.7 Mb): View
- Cβ deviation scatter plot (44 Kb): View in KiNG | Download

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