

Analysis output: all-atom contacts and geometry for MITF_324-402H.pdb

Summary statistics

All-Atom Contacts	Clashscore, all atoms:	4.75		94 th percentile* (N=1784, all resolutions)
	Clashscore is the number of serious steric overlaps (> 0.4 Å) per 1000 atoms.			
Protein Geometry	Poor rotamers	0	0.00%	Goal: <0.3%
	Favored rotamers	140	97.22%	Goal: >98%
	Ramachandran outliers	0	0.00%	Goal: <0.05%
	Ramachandran favored	154	100.00%	Goal: >98%
	MolProbity score [^]	1.24		99 th percentile* (N=27675, 0Å - 99Å)
	Cβ deviations >0.25Å	0	0.00%	Goal: 0
	Bad bonds:	11 / 1358	0.81%	Goal: 0%
	Bad angles:	1 / 1812	0.06%	Goal: <0.1%
Peptide Omegas	Cis Prolines:	0 / 4	0.00%	Expected: ≤1 per chain, or ≤5%

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

Multi-criterion visualizations



View (170 Kb)

Single-criterion visualizations

^{* 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.

- Clash list (721 bytes): View
- Ramachandran plot kinemage (414 Kb): View in KiNG | Download
 Ramachandran plot PDF (1.7 Mb): View
- Cβ deviation scatter plot (19 Kb): View in KiNG | Download

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