

Analysis output: all-atom contacts and geometry for MYO6_1175-1277.pdb_4_1H.pdb

http://molprobity.biochem.duke.edu/index.php?MolProbSID=qv...

Summary statistics

All-Atom Contacts	Clashscore, all atoms:	2.92		98 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: <1%
	Ramachandran outliers	0	0.00%	Goal: <0.05%
	Ramachandran favored	100	99.01%	Goal: >98%
	MolProbity score [^]	1.08		100 th percentile* (N=27675, 0Å - 99Å)
	Cβ deviations >0.25Å	0	0.00%	Goal: 0
	Bad backbone bonds:	4 / 892	0.45%	Goal: 0%
	Bad backbone angles:	1 / 1203	0.08%	Goal: <0.1%

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

Multi-criterion visualizations



View (100 Kb)

1 of 2 6/18/14, 10:11 AM

^{* 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.

Single-criterion visualizations

- Clash list (385 bytes): View
- Ramachandran plot kinemage (410 Kb): View in KiNG | Download
- Ramachandran plot PDF (1.7 Mb): View
- Cβ deviation scatter plot (17 Kb): View in KiNG | Download



About MolProbity | Website for the Richardson Lab | Using ecloud x-H | Internal reference 4.1-537

2 of 2