

Analysis output: all-atom contacts and geometry for KITLG_33-507_FFX1H.pdb

Summary statistics

All-Atom Contacts	Clashscore, all atoms:	0.13		99 th percentile* (N=1784, all resolutions)
	Clashscore is the number of serious steric overlaps (> 0.4 Å) per 1000 atoms.			
Protein Geometry	Poor rotamers	7	1.64%	Goal: <0.3%
	Favored rotamers	383	89.91%	Goal: >98%
	Ramachandran outliers	8	1.69%	Goal: <0.05%
	Ramachandran favored	402	84.99%	Goal: >98%
	MolProbity score [^]	1.38		97 th percentile* (N=27675, 0Å - 99Å)
	Cβ deviations >0.25Å	1	0.22%	Goal: 0
	Bad bonds:	0/3858	0.00%	Goal: 0%
	Bad angles:	28 / 5245	0.53%	Goal: <0.1%
Peptide Omegas	Cis Prolines:	3 / 25	12.00%	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 (497 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 (217 bytes): View
- Ramachandran plot kinemage (433 Kb): View in KiNG | Download
 Ramachandran plot PDF (1.7 Mb): View
- **C**β deviation scatter plot (38 Kb): View in KiNG | Download

Continue >

About MolProbity | Website for the Richardson Lab | Using ecloud x-H | Internal reference 4.3