



# Analysis output: all-atom contacts and geometry for PDZD7\_82- 170\_FFX1H.pdb

## Summary statistics

All-Atom Contacts	Clashscore, all atoms:	0.75	99 <sup>th</sup> percentile* (N=1784, all resolutions)	
	Clashscore is the number of serious steric overlaps (> 0.4 Å) per 1000 atoms.			
Protein Geometry	Poor rotamers	2	2.74%	Goal: <0.3%
	Favored rotamers	61	83.56%	Goal: >98%
	Ramachandran outliers	1	1.15%	Goal: <0.05%
	Ramachandran favored	75	86.21%	Goal: >98%
	MolProbity score <sup>^</sup>	1.71	89 <sup>th</sup> percentile* (N=27675, 0Å - 99Å)	
	Cβ deviations >0.25Å	0	0.00%	Goal: 0
	Bad bonds:	0 / 661	0.00%	Goal: 0%
	Bad angles:	2 / 884	0.23%	Goal: <0.1%
Peptide Omegas	Cis Prolines:	0 / 2	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.

\* 100<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.

## Multi-criterion visualizations



[View](#) (95 Kb)

## Single-criterion visualizations

- **Clash list** (217 bytes): [View](#)
- **Ramachandran plot kinemage** (410 Kb): [View in KiNG](#) | [Download](#)
- **Ramachandran plot PDF** (1.7 Mb): [View](#)
- **C $\beta$  deviation scatter plot** (15 Kb): [View in KiNG](#) | [Download](#)

