

## Analysis output: all-atom contacts and geometry for TJP2\_21129\_FFX1H.pdb

## **Summary statistics**

All-Atom Contacts	Clashscore, all atoms:	0		100 <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	3	3.37%	Goal: <0.3%
	Favored rotamers	82	92.13%	Goal: >98%
	Ramachandran outliers	3	2.80%	Goal: <0.05%
	Ramachandran favored	92	85.98%	Goal: >98%
	MolProbity score <sup>^</sup>	1.54		94 <sup>th</sup> percentile* (N=27675, 0Å - 99Å)
	Cβ deviations >0.25Å	0	0.00%	Goal: 0
	Bad bonds:	0 / 840	0.00%	Goal: 0%
	Bad angles:	10 / 1137	0.88%	Goal: <0.1%
Peptide Omegas	Cis Prolines:	0 / 6	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 (116 Kb)

## Single-criterion visualizations

<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.

- Clash list (175 bytes): View
- Ramachandran plot kinemage (411 Kb): View in KiNG | Download
  Ramachandran plot PDF (1.7 Mb): View
- **C**β deviation scatter plot (16 Kb): View in KiNG | Download

Continue >

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