

## Analysis output: all-atom contacts and geometry for SLITRK\_31-267H.pdb

## **Summary statistics**

All-Atom Contacts	Clashscore, all atoms:	28.17		17 <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	8	3.72%	Goal: <0.3%
	Favored rotamers	195	90.70%	Goal: >98%
	Ramachandran outliers	4	1.70%	Goal: <0.05%
	Ramachandran favored	208	88.51%	Goal: >98%
	MolProbity score <sup>^</sup>	2.96		25 <sup>th</sup> percentile* (N=27675, 0Å - 99Å)
	Cβ deviations >0.25Å	1	0.45%	Goal: 0
	Bad bonds:	13 / 1902	0.68%	Goal: 0%
	Bad angles:	12 / 2582	0.46%	Goal: <0.1%
Peptide Omegas	Cis Prolines:	1 / 12	8.33%	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 (264 Kb)

<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.

## **Single-criterion visualizations**

- Clash list (4.6 Kb): View
- Ramachandran plot kinemage (419 Kb): View in KiNG | Download
- Ramachandran plot PDF (1.7 Mb): View
- Cβ deviation scatter plot (23 Kb): View in KiNG | Download

Continue >

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