

## Analysis output: all-atom contacts and geometry for Model\_3H.pdb

## **Summary statistics**

| Contacts            | Clashscore, all atoms:  | 31.11        |        | 14 <sup>th</sup> percentile* (N=1784, all resolutions) |
|---------------------|---|--------------|--------|--|
|                     | Clashscore is the number of serious steric overlaps (> 0.4 Å) per 1000 atoms. |              |        |  |
| Protein<br>Geometry | Poor rotamers   | 10           | 5.13%  | Goal: <1%  |
|                     | Ramachandran<br>outliers  | 10           | 4.67%  | Goal: <0.05%   |
|                     | Ramachandran<br>favored   | 176          | 82.24% | Goal: >98%   |
|                     | MolProbity score <sup>^</sup>   | 3.22         |        | 16 <sup>th</sup> percentile* (N=27675, 0Å - 99Å)       |
|                     | Cβ deviations >0.25Å  | 2            | 0.97%  | Goal: 0  |
|                     | Bad backbone bonds:   | 14 /<br>1823 | 0.77%  | Goal: 0%   |
|                     | Bad backbone angles:  | 13 /<br>2473 | 0.53%  | Goal: <0.1%  |

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

## **Multi-criterion visualizations**



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<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.

View (218 Kb)

## Single-criterion visualizations

- Clash list (3.5 Kb): View
- Ramachandran plot kinemage (418 Kb): View in KiNG | Download
- Ramachandran plot PDF (1.7 Mb): View
- Cβ deviation scatter plot (21 Kb): View in KiNG | Download

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