

**Summary of integrative structure determination of Structure of N4BP1
CUE domain from NMR chemical shifts (PDB ID: 9A14, PDB-Dev ID:
PDBDEV_00000076)**

1. Model Composition	
Entry composition	N4BP1 CUE domain: chain(s) A (44 residues)
Datasets used for modeling	- NMR data, BMRB: 50688
2. Representation	
Number of representations	1
Scale	Atomic
Number of rigid and flexible segments	0, 1
3. Restraints	
Physical principles	Information about physical principles was not provided
Experimental data	
4. Validation	
Number of ensembles	0
Number of models in ensembles	Not applicable
Number of deposited models	10
Model precision (uncertainty of models)	Not available
Data quality	Data quality has not been assessed
Model quality: assessment of atomic segments	<ul style="list-style-type: none"> - Clashscore: 0.00-2.80 - Ramachandran outliers: 0-0 - Sidechain outliers: 0-0
Fit to data used for modeling	Fit of model to information used to compute it has not been determined
Fit to data used for validation	Fit of model to information not used to compute it has not been determined
5. Methodology and Software	
1. Name	CS-ROSETTA modeling
Method	None

<i>Number of computed models</i>	20000
<i>Software</i>	CS-ROSETTA (version Not available)