wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	1ZR9
Title	:	Solution Structure of a Human C2H2-type Zinc Finger Protein
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This is a wwPDB NMR Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/NMRValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (i)) were used in the production of this report:

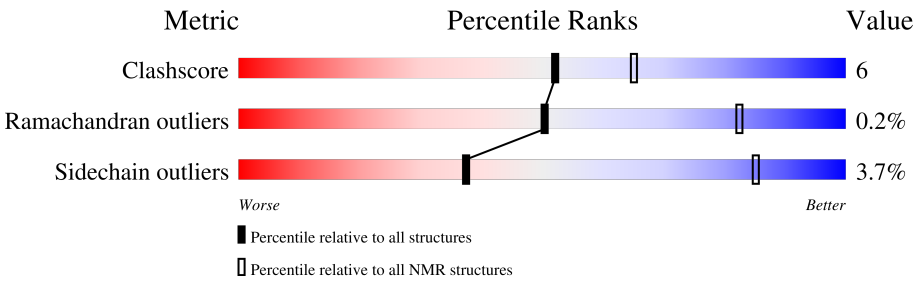
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.26
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.26

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and a grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Mol	Chain	Length	Quality of chain
1	A	124	

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 1 is the overall representative, medoid model (most similar to other models).

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	**Residue range (total)**	**Backbone RMSD (Å)**	**Medoid model**
1	A:38-A:68 (31)	0.27	1

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 4 clusters and 3 single-model clusters were found.

Cluster number	Models
1	1, 2, 3, 4, 5, 10, 14, 18, 20
2	6, 7, 19
3	9, 11, 17
4	12, 16
Single-model clusters	8; 13; 15

3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1039 atoms, of which 508 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Zinc finger protein 593.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	67	1038	329	508	97	101	3	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	cloning artifact	UNP O00488
A	-6	HIS	-	expression tag	UNP O00488
A	-5	HIS	-	expression tag	UNP O00488
A	-4	HIS	-	expression tag	UNP O00488
A	-3	HIS	-	expression tag	UNP O00488
A	-2	HIS	-	expression tag	UNP O00488
A	-1	HIS	-	expression tag	UNP O00488
A	0	LEU	-	cloning artifact	UNP O00488
A	1	GLU	-	cloning artifact	UNP O00488

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	
			Total	Zn
2	A	1	1	1

4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: Zinc finger protein 593

Chain A: 20% • • 29% 46%

GLY HIS HIS HIS HIS HIS HIS LEU GLU LYS ALA LYS ARG ARG ARG PRO ASP LEU ASP GLU ILE HIS ARG GLU LEU ARG PRO GLN GLY SER ALA ARG PRO GLN PRO D28 P29 N30 A31 E32 F33 D34 P35 D36 L37 C45 L58 H61 F62 K65 D66 H67 K68 K69 R70 L71

K72 Q73 L74 S75 V76 E77 P78 Y79 S80 Q81 E82 E83 A84 E85 R86 A87 A88 G89 M90 G91 S92 Y93 V94 PRO PRO ARG ARG LEU ALA VAL PRO THR GLU VAL SER THR GLU VAL PRO GLU MET ASP THR SER THR

4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 1. Colouring as in section 4.1 above.

- Molecule 1: Zinc finger protein 593

Chain A: 19% • • 29% 46%

GLY HIS HIS HIS HIS HIS HIS LEU GLU LYS ALA LYS ARG ARG ARG PRO ASP LEU ASP GLU ILE HIS ARG GLU LEU ARG PRO GLN GLY SER ALA ARG PRO GLN PRO D28 P29 N30 A31 E32 F33 D34 P35 D36 L37 C45 L58 H61 F62 K65 D66 H67 K68 K69 R70 L71

K72 Q73 L74 S75 V76 E77 P78 Y79 S80 Q81 E82 E83 A84 E85 R86 A87 A88 G89 M90 G91 S92 Y93 V94 PRO PRO ARG ARG LEU ALA VAL PRO THR GLU VAL SER THR GLU VAL PRO GLU MET ASP THR SER THR

5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *Automated methods were used for backbone chemical shift assignment and iterative NOE refinement. Final structures were obtained by molecular dynamics in explicit solvent.*.

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *target function.*

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
CYANA	structure solution	1.0.6
XPLOR-NIH	refinement	2.0.6

No chemical shift data was provided.

6 Model quality (i)

6.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	0.90±0.04	0±0/254 (0.1± 0.2%)	0.64±0.02	0±0/339 (0.0± 0.0%)
All	All	0.90	6/5080 (0.1%)	0.64	0/6780 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	0.9±0.2
All	All	0	19

All unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	48	CYS	CB-SG	-5.76	1.72	1.81	17	3
1	A	67	HIS	CG-CD2	-5.51	1.26	1.35	15	3

There are no bond-angle outliers.

There are no chirality outliers.

All unique planar outliers are listed below.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	67	HIS	Sidechain	19

6.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	247	243	242	3±1
All	All	4960	4860	4840	56

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

5 of 18 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:62:PHE:O	1:A:68:LYS:HD2	0.56	2.00	10	2
1:A:58:LEU:O	1:A:61:HIS:HB3	0.55	2.02	20	15
1:A:45:CYS:SG	1:A:62:PHE:HE1	0.54	2.25	7	13
1:A:43:HIS:CE1	1:A:55:SER:HB3	0.53	2.38	18	1
1:A:62:PHE:HB3	1:A:68:LYS:HD3	0.52	1.82	8	2

6.3 Torsion angles (i)

6.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the backbone conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	31/124 (25%)	30±1 (98±2%)	0±1 (2±2%)	0±0 (0±1%)	50	82
All	All	620/2480 (25%)	609 (98%)	10 (2%)	1 (0%)	50	82

All 1 unique Ramachandran outliers are listed below.

Mol	Chain	Res	Type	Models (Total)
1	A	38	PRO	1

6.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the sidechain conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	26/108 (24%)	25±1 (96±3%)	1±1 (4±3%)	37 85
All	All	520/2160 (24%)	501 (96%)	19 (4%)	37 85

All 4 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	65	LYS	9
1	A	44	ARG	7
1	A	53	ILE	2
1	A	57	ASN	1

6.3.3 RNA (i)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates (i)

There are no monosaccharides in this entry.

6.6 Ligand geometry (i)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

6.7 Other polymers (i)

There are no such molecules in this entry.

6.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

7 Chemical shift validation (i)

No chemical shift data were provided