



Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 06:41 PM BST

PDB ID : 1ZR9
Title : Solution Structure of a Human C2H2-type Zinc Finger Protein
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Deposited on : 2005-05-19

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/NMRValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

Cyrange : Kirchner and Güntert (2011)
NmrClust : Kelley et al. (1996)
MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : rb-20027457
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027457

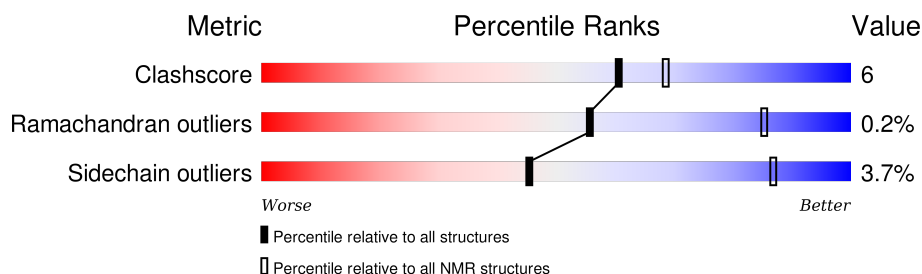
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 86%.

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	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	124	 20% . . 29% 46%

2 Ensemble composition and analysis

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

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.

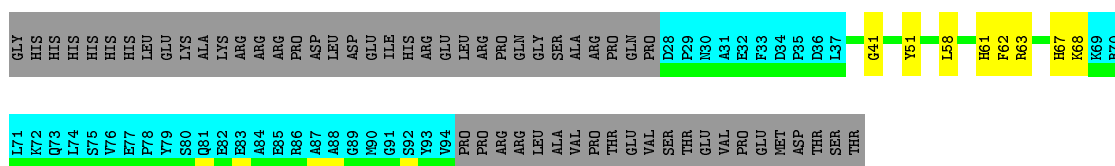
Mol	Chain	Residues	Atoms						Trace
1	A	67	Total	C	H	N	O	S	0
			1038	329	508	97	101	3	

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).

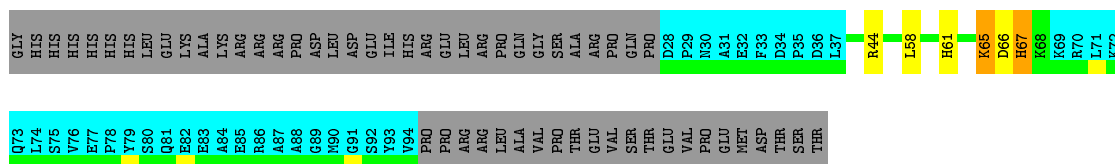
Mol	Chain	Residues	Atoms	
2	A	1	Total	Zn
			1	1



4.2.3 Score per residue for model 3

- Molecule 1: Zinc finger protein 593

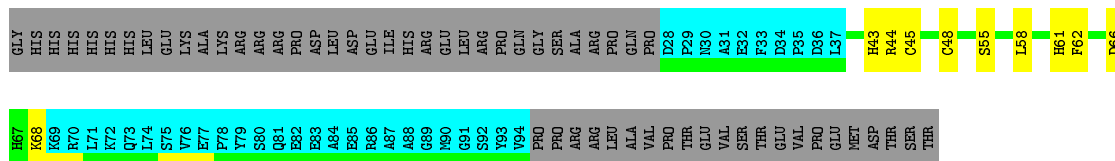
Chain A: 20% 29% 46%



4.2.4 Score per residue for model 4

- Molecule 1: Zinc finger protein 593

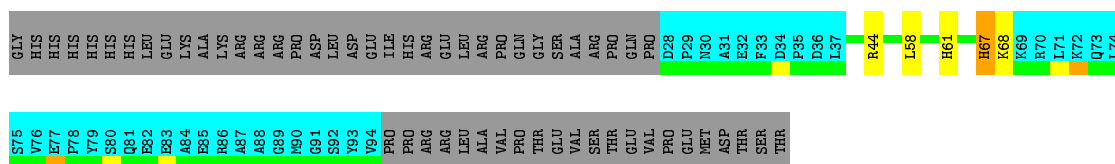
Chain A: 17% 8% 29% 46%



4.2.5 Score per residue for model 5

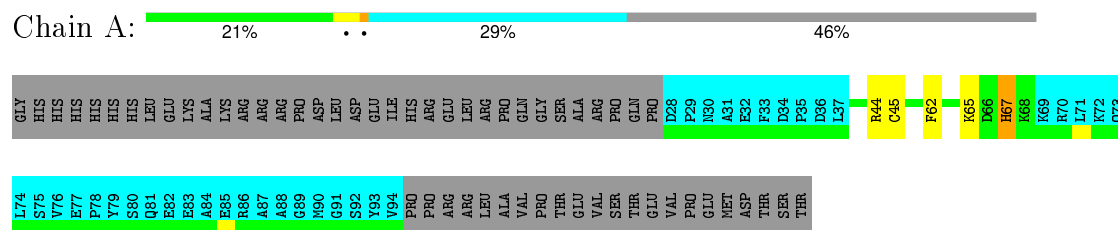
- Molecule 1: Zinc finger protein 593

Chain A: 21% 29% 46%



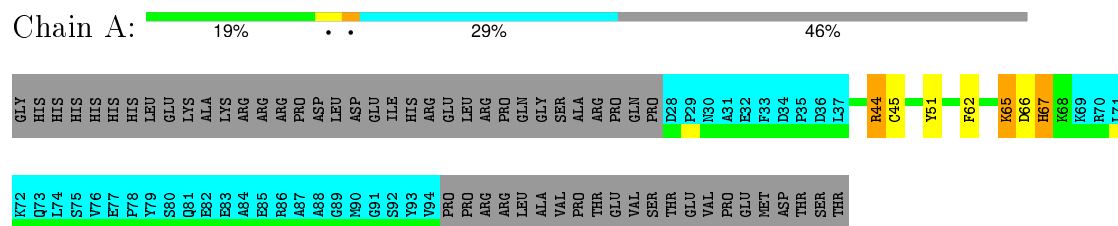
4.2.6 Score per residue for model 6

- Molecule 1: Zinc finger protein 593



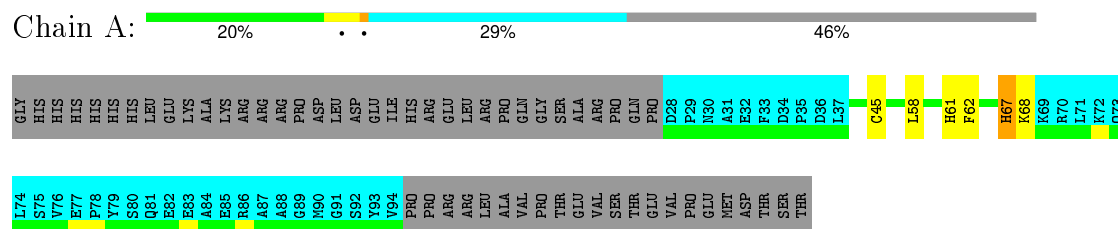
4.2.7 Score per residue for model 7

- Molecule 1: Zinc finger protein 593



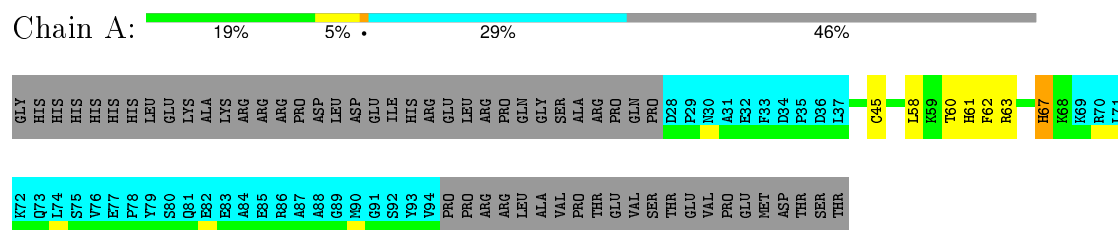
4.2.8 Score per residue for model 8

- Molecule 1: Zinc finger protein 593



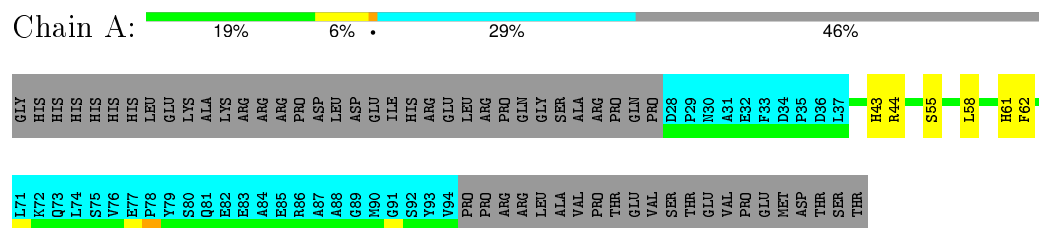
4.2.9 Score per residue for model 9

- Molecule 1: Zinc finger protein 593



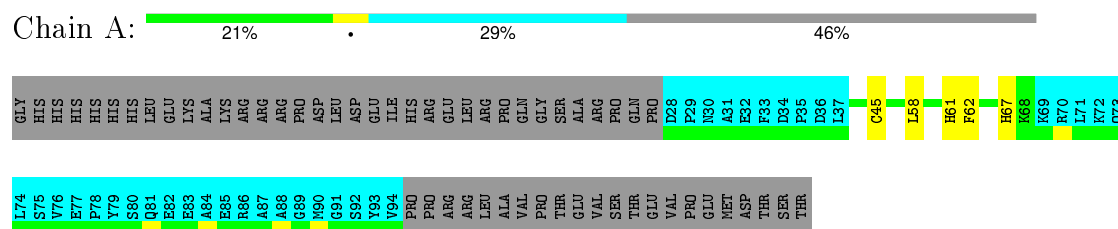
4.2.10 Score per residue for model 10

- Molecule 1: Zinc finger protein 593



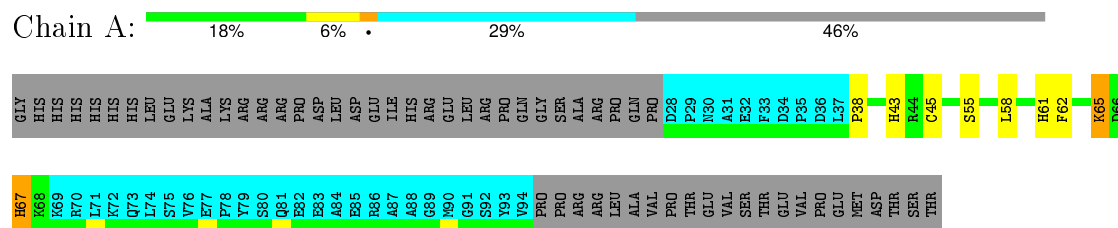
4.2.11 Score per residue for model 11

- Molecule 1: Zinc finger protein 593



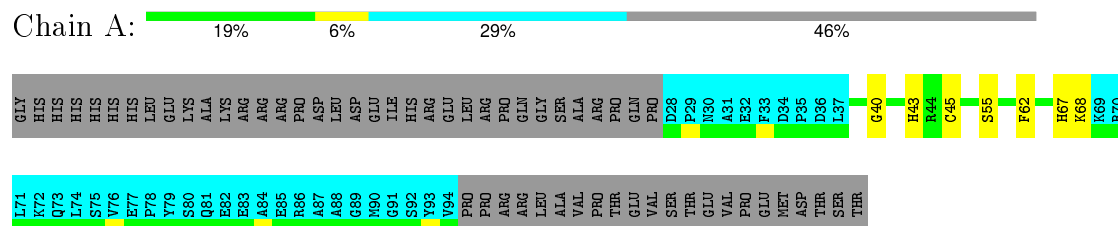
4.2.12 Score per residue for model 12

- Molecule 1: Zinc finger protein 593



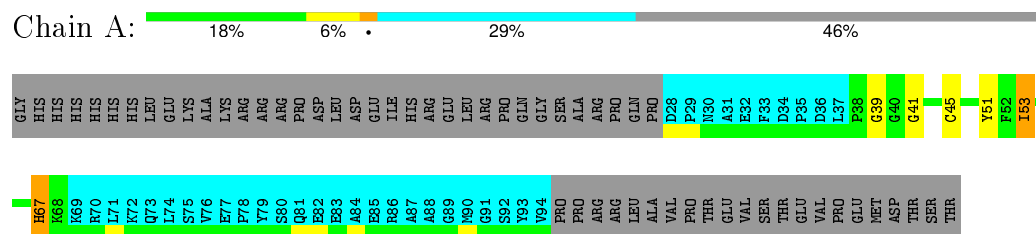
4.2.13 Score per residue for model 13

- Molecule 1: Zinc finger protein 593



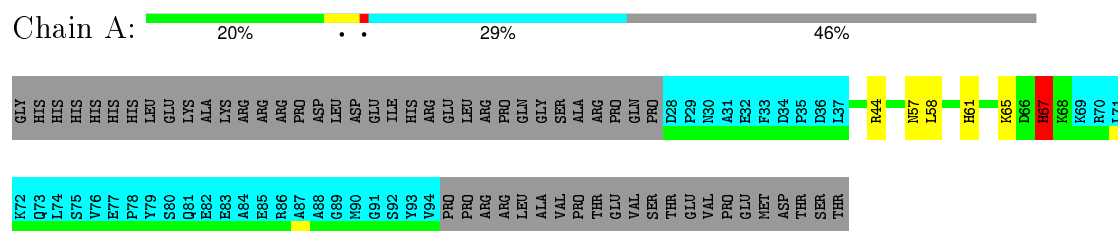
4.2.14 Score per residue for model 14

- Molecule 1: Zinc finger protein 593



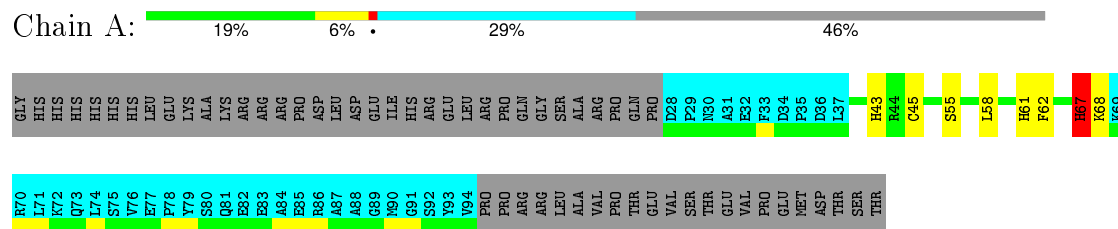
4.2.15 Score per residue for model 15

- Molecule 1: Zinc finger protein 593



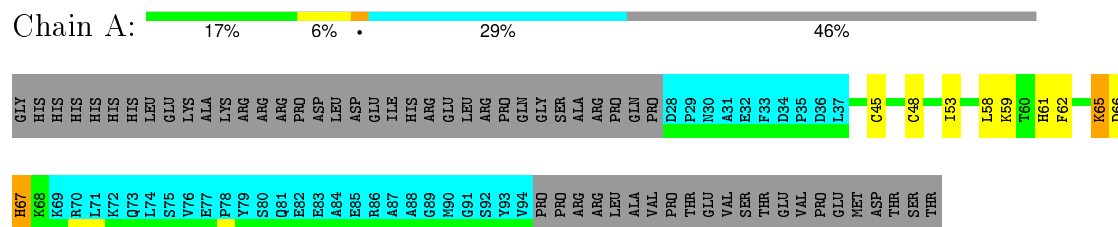
4.2.16 Score per residue for model 16

- Molecule 1: Zinc finger protein 593



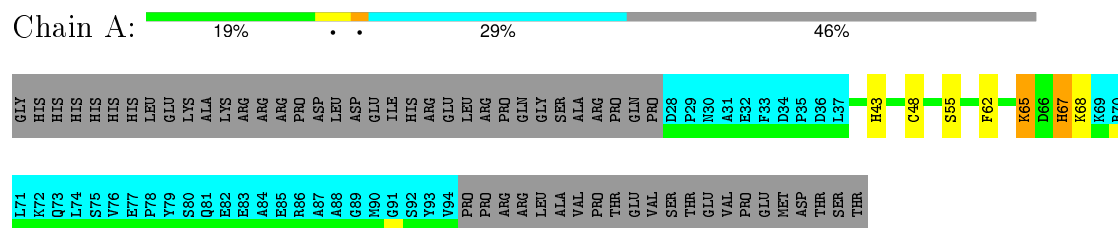
4.2.17 Score per residue for model 17

- Molecule 1: Zinc finger protein 593



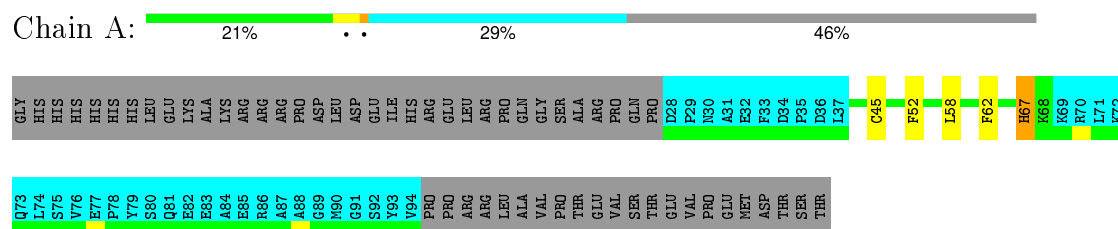
4.2.18 Score per residue for model 18

- Molecule 1: Zinc finger protein 593



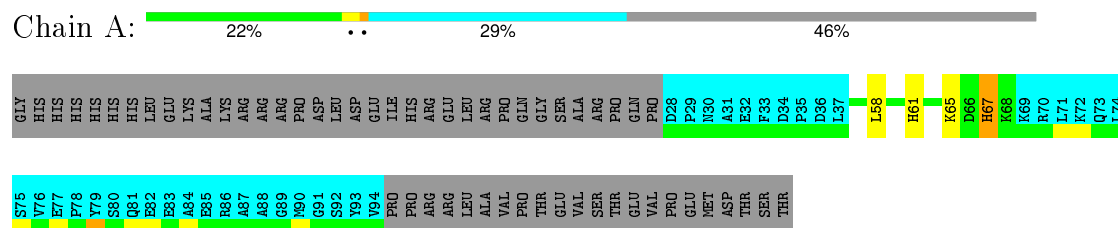
4.2.19 Score per residue for model 19

- Molecule 1: Zinc finger protein 593



4.2.20 Score per residue for model 20

- Molecule 1: Zinc finger protein 593



5 Refinement protocol and experimental data overview

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

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	BMRB entry 6682
Number of chemical shift lists	1
Total number of shifts	750
Number of shifts mapped to atoms	750
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	86%

No validations of the models with respect to experimental NMR restraints is performed at this time.

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 ⓘ

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.

All 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
1:A:65:LYS:HD2	1:A:66:ASP:N	0.50	2.21	3	4
1:A:63:ARG:HA	1:A:63:ARG:NE	0.49	2.22	2	1
1:A:43:HIS:CE1	1:A:55:SER:HB2	0.47	2.44	12	3
1:A:41:GLY:HA2	1:A:51:TYR:CD1	0.47	2.45	2	2
1:A:62:PHE:HB3	1:A:68:LYS:HD2	0.47	1.87	16	3
1:A:39:GLY:N	1:A:53:ILE:HA	0.46	2.26	14	1
1:A:43:HIS:ND1	1:A:55:SER:HB2	0.46	2.25	4	2
1:A:59:LYS:HE2	1:A:59:LYS:HA	0.44	1.88	17	1
1:A:52:PHE:HB2	1:A:58:LEU:CD2	0.44	2.42	19	1
1:A:65:LYS:HD3	1:A:65:LYS:H	0.44	1.72	18	1
1:A:65:LYS:H	1:A:65:LYS:HD3	0.44	1.72	12	1
1:A:60:THR:O	1:A:63:ARG:HG2	0.43	2.13	9	1
1:A:44:ARG:HD2	1:A:51:TYR:CE2	0.41	2.50	7	1
1:A:43:HIS:ND1	1:A:55:SER:HB3	0.41	2.31	18	1

6.3 Torsion angles ⓘ

6.3.1 Protein backbone ⓘ

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%)	1±1 (2±2%)	0±0 (0±1%)	56 85
All	All	620/2480 (25%)	609 (98%)	10 (2%)	1 (0%)	56 85

All 1 unique Ramachandran outliers are listed below.

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

6.3.2 Protein sidechains ⓘ

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%)	45 87
All	All	520/2160 (24%)	501 (96%)	19 (4%)	45 87

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 ⓘ

There are no RNA molecules in this entry.

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

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

6.5 Carbohydrates [i](#)

There are no carbohydrates 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

The completeness of assignment taking into account all chemical shift lists is 86% for the well-defined parts and 87% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 6682

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	750
Number of shifts mapped to atoms	750
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	67	-0.44 ± 0.19	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	62	-0.11 ± 0.18	None needed (< 0.5 ppm)
$^{13}\text{C}'$	65	-0.19 ± 0.23	None needed (< 0.5 ppm)
^{15}N	63	-0.36 ± 0.32	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 86%, i.e. 342 atoms were assigned a chemical shift out of a possible 396. 3 out of 3 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	152/153 (99%)	61/61 (100%)	61/62 (98%)	30/30 (100%)
Sidechain	158/195 (81%)	100/117 (85%)	57/65 (88%)	1/13 (8%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	32/48 (67%)	20/26 (77%)	12/18 (67%)	0/4 (0%)
Overall	342/396 (86%)	181/204 (89%)	130/145 (90%)	31/47 (66%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 87%, i.e. 722 atoms were assigned a chemical shift out of a possible 831. 6 out of 8 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	325/327 (99%)	130/130 (100%)	132/134 (99%)	63/63 (100%)
Sidechain	352/431 (82%)	222/257 (86%)	125/150 (83%)	5/24 (21%)
Aromatic	45/73 (62%)	29/39 (74%)	16/30 (53%)	0/4 (0%)
Overall	722/831 (87%)	381/426 (89%)	273/314 (87%)	68/91 (75%)

7.1.4 Statistically unusual chemical shifts ⓘ

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	86	ARG	NE	111.90	92.63 – 76.73	17.1

7.1.5 Random Coil Index (RCI) plots ⓘ

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition.

Random coil index (RCI) for chain A:

