



Full wwPDB NMR Structure Validation Report ⓘ

Nov 9, 2024 – 01:06 PM EST

PDB ID : 1ZNF
Title : THREE-DIMENSIONAL SOLUTION STRUCTURE OF A SINGLE ZINC FINGER DNA-BINDING DOMAIN
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Deposited on : 1989-09-25

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

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

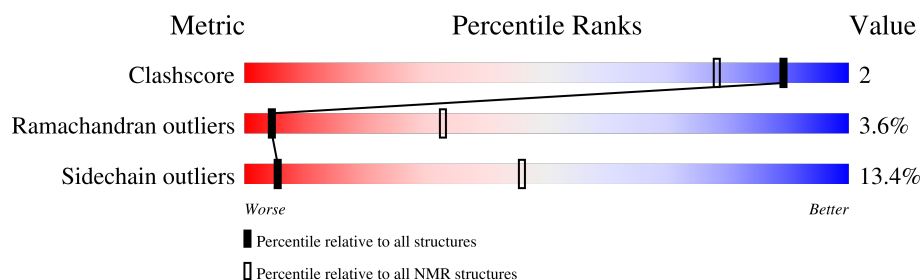
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment 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	210492	14027
Ramachandran outliers	207382	12486
Sidechain outliers	206894	12463

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	27	

2 Ensemble composition and analysis

This entry contains 37 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:1-A:24 (24)	0.60	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 4 single-model clusters were found.

Cluster number	Models
1	1, 2, 3, 5, 6, 7, 9, 10, 11, 12, 13, 15, 18, 19, 20, 23, 25, 28, 29, 30, 31, 32, 33, 34, 36
2	8, 24, 26
3	21, 22, 27
4	14, 17
Single-model clusters	4; 16; 35; 37

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 424 atoms, of which 212 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called 31ST ZINC FINGER FROM XFIN.

Mol	Chain	Residues	Atoms						Trace
1	A	27	Total	C	H	N	O	S	1
			423	129	212	44	36	2	

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

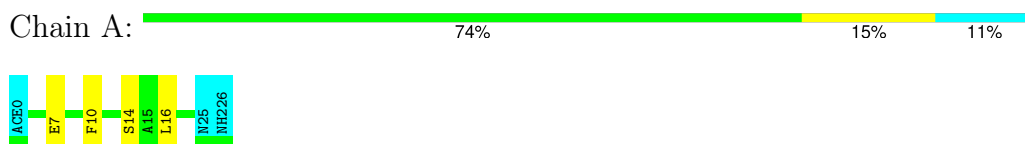
Mol	Chain	Residues	Atoms	
2	A	1	Total	Zn
			1	1

4 Residue-property plots

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: 31ST ZINC FINGER FROM XFIN

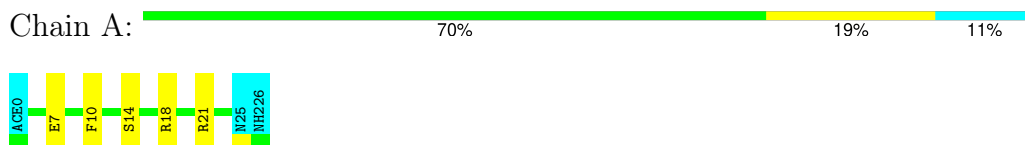


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

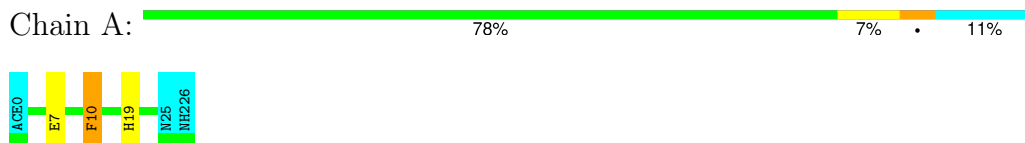
4.2.1 Score per residue for model 1 (medoid)

- Molecule 1: 31ST ZINC FINGER FROM XFIN



4.2.2 Score per residue for model 2

- Molecule 1: 31ST ZINC FINGER FROM XFIN



4.2.3 Score per residue for model 3

- Molecule 1: 31ST ZINC FINGER FROM XFIN

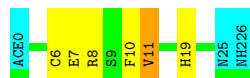
Chain A:  70% 15% 11%



4.2.4 Score per residue for model 4

- Molecule 1: 31ST ZINC FINGER FROM XFIN

Chain A:  67% 19% 11%



4.2.5 Score per residue for model 5

- Molecule 1: 31ST ZINC FINGER FROM XFIN

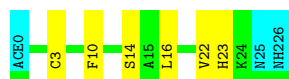
Chain A:  70% 19% 11%



4.2.6 Score per residue for model 6


- Molecule 1: 31ST ZINC FINGER FROM XFIN

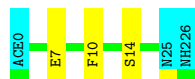
Chain A:  67% 22% 11%



4.2.7 Score per residue for model 7

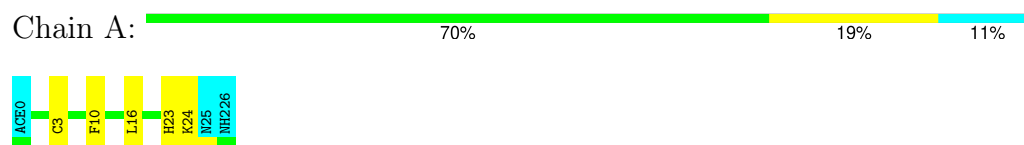
- Molecule 1: 31ST ZINC FINGER FROM XFIN

Chain A:  78% 11% 11%



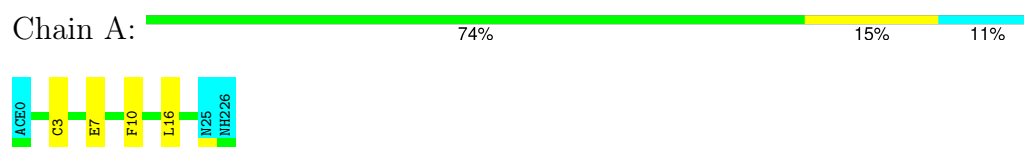
4.2.8 Score per residue for model 8

- Molecule 1: 31ST ZINC FINGER FROM XFIN



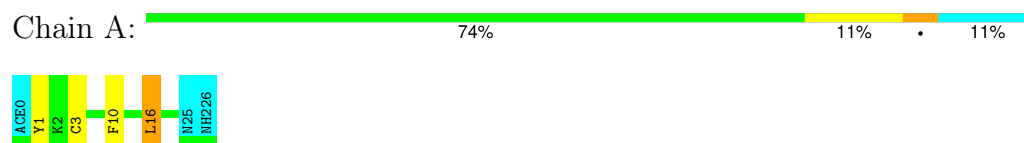
4.2.9 Score per residue for model 9

- Molecule 1: 31ST ZINC FINGER FROM XFIN



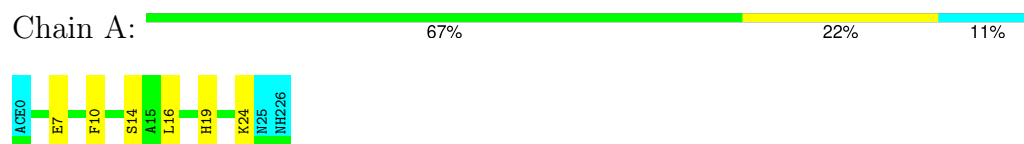
4.2.10 Score per residue for model 10

- Molecule 1: 31ST ZINC FINGER FROM XFIN



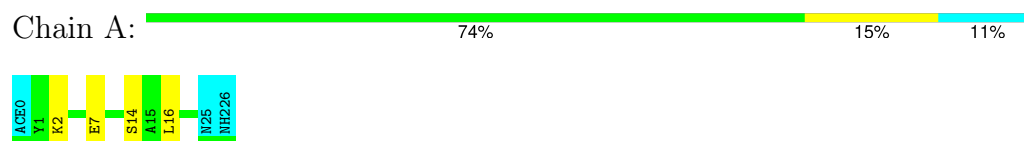
4.2.11 Score per residue for model 11

- Molecule 1: 31ST ZINC FINGER FROM XFIN



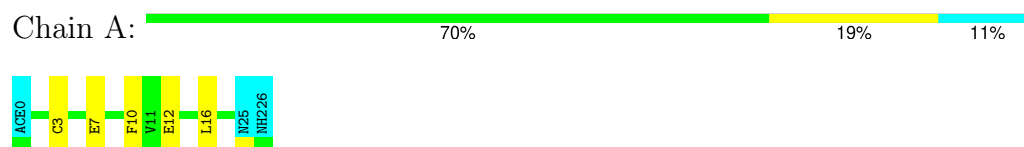
4.2.12 Score per residue for model 12

- Molecule 1: 31ST ZINC FINGER FROM XFIN



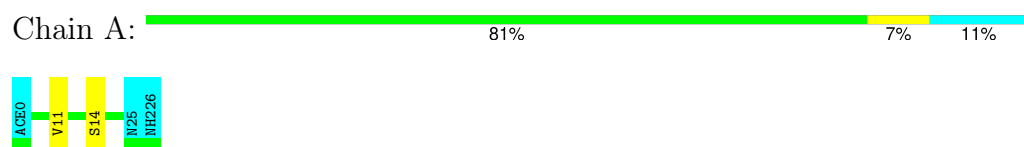
4.2.13 Score per residue for model 13

- Molecule 1: 31ST ZINC FINGER FROM XFIN



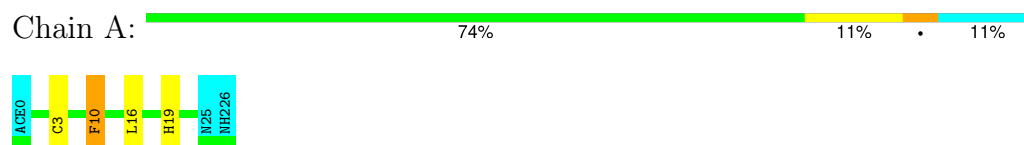
4.2.14 Score per residue for model 14

- Molecule 1: 31ST ZINC FINGER FROM XFIN



4.2.15 Score per residue for model 15

- Molecule 1: 31ST ZINC FINGER FROM XFIN



4.2.16 Score per residue for model 16

- Molecule 1: 31ST ZINC FINGER FROM XFIN



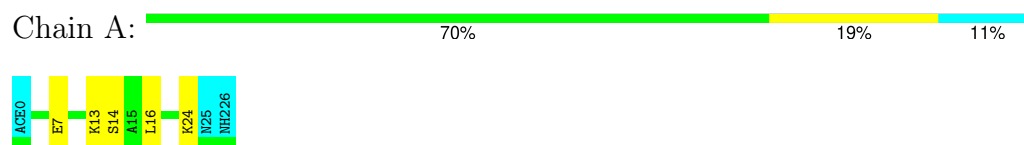
4.2.17 Score per residue for model 17

- Molecule 1: 31ST ZINC FINGER FROM XFIN



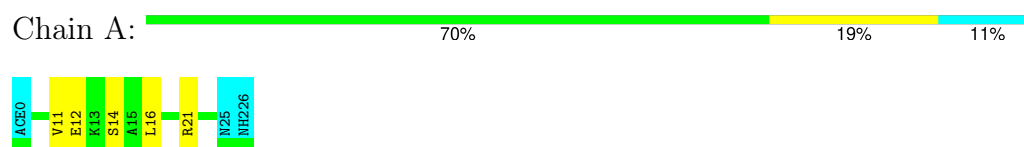
4.2.18 Score per residue for model 18

- Molecule 1: 31ST ZINC FINGER FROM XFIN



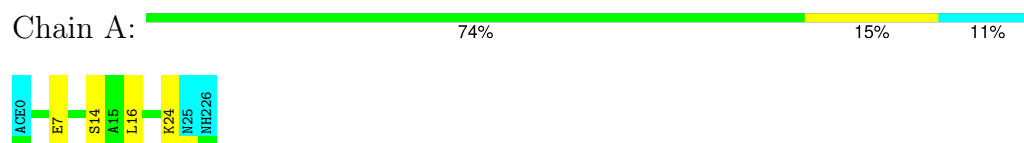
4.2.19 Score per residue for model 19

- Molecule 1: 31ST ZINC FINGER FROM XFIN



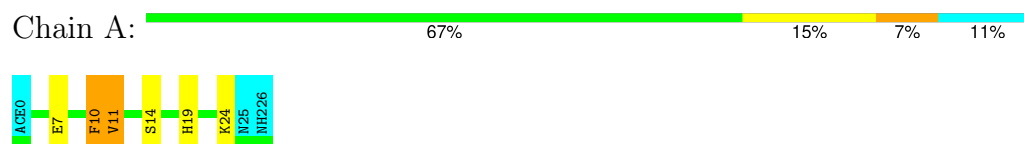
4.2.20 Score per residue for model 20

- Molecule 1: 31ST ZINC FINGER FROM XFIN



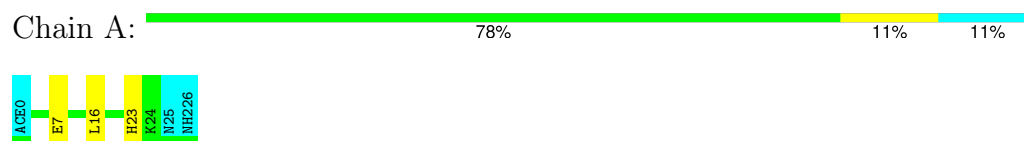
4.2.21 Score per residue for model 21

- Molecule 1: 31ST ZINC FINGER FROM XFIN



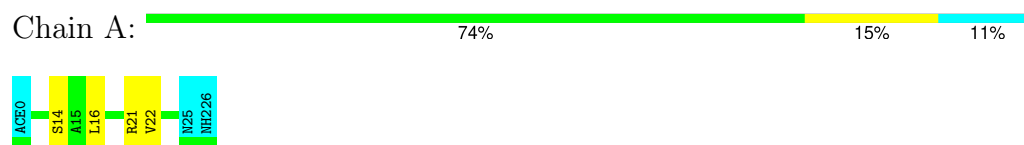
4.2.22 Score per residue for model 22

- Molecule 1: 31ST ZINC FINGER FROM XFIN



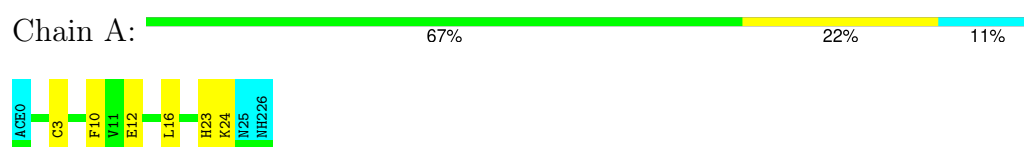
4.2.23 Score per residue for model 23

- Molecule 1: 31ST ZINC FINGER FROM XFIN



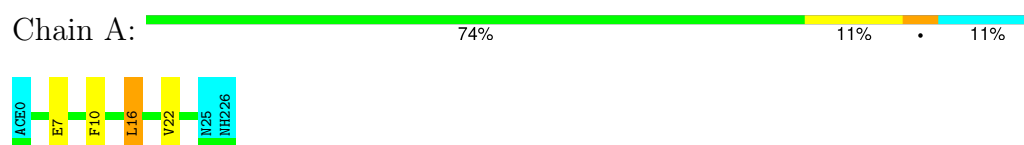
4.2.24 Score per residue for model 24

- Molecule 1: 31ST ZINC FINGER FROM XFIN



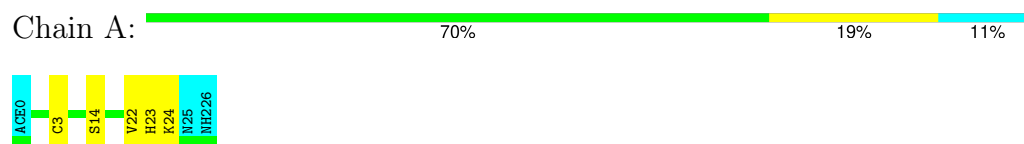
4.2.25 Score per residue for model 25

- Molecule 1: 31ST ZINC FINGER FROM XFIN



4.2.26 Score per residue for model 26

- Molecule 1: 31ST ZINC FINGER FROM XFIN



4.2.27 Score per residue for model 27

- Molecule 1: 31ST ZINC FINGER FROM XFIN



4.2.28 Score per residue for model 28


- Molecule 1: 31ST ZINC FINGER FROM XFIN

Chain A:  70% 19% 11%



4.2.29 Score per residue for model 29

- Molecule 1: 31ST ZINC FINGER FROM XFIN

Chain A:  78% 11% 11%



4.2.30 Score per residue for model 30

- Molecule 1: 31ST ZINC FINGER FROM XFIN

Chain A:  67% 22% 11%



4.2.31 Score per residue for model 31

- Molecule 1: 31ST ZINC FINGER FROM XFIN

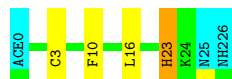
Chain A:  70% 19% 11%



4.2.32 Score per residue for model 32

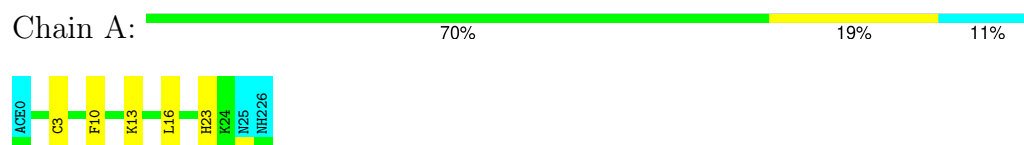
- Molecule 1: 31ST ZINC FINGER FROM XFIN

Chain A:  74% 11% 11%



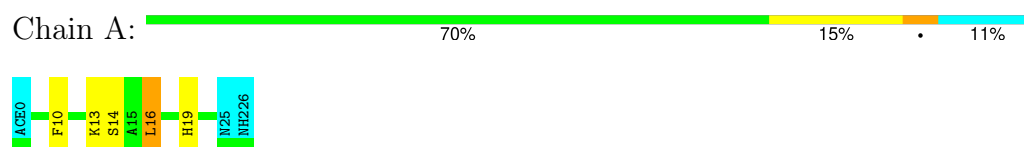
4.2.33 Score per residue for model 33

- Molecule 1: 31ST ZINC FINGER FROM XFIN



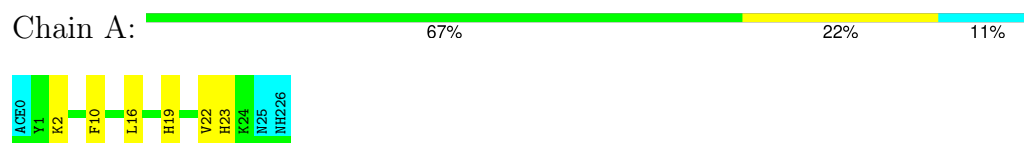
4.2.34 Score per residue for model 34

- Molecule 1: 31ST ZINC FINGER FROM XFIN



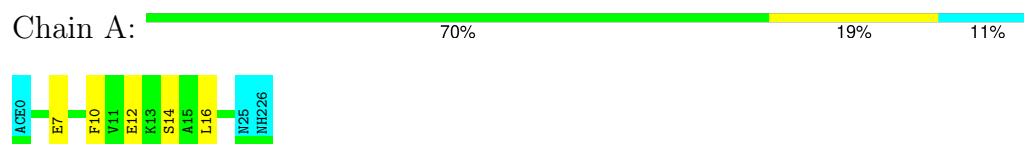
4.2.35 Score per residue for model 35

- Molecule 1: 31ST ZINC FINGER FROM XFIN



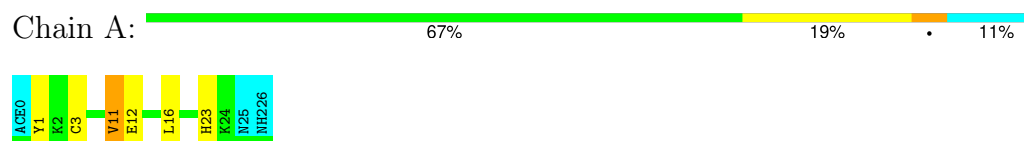
4.2.36 Score per residue for model 36

- Molecule 1: 31ST ZINC FINGER FROM XFIN



4.2.37 Score per residue for model 37

- Molecule 1: 31ST ZINC FINGER FROM XFIN



5 Refinement protocol and experimental data overview ⓘ

Of the ? calculated structures, 37 were deposited, based on the following criterion: ?.

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

Software name	Classification	Version
DISGEO	refinement	
Amber	refinement	

No chemical shift data was provided.

6 Model quality

6.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, ACE, NH2

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	0.56±0.01	0±0/203 (0.0± 0.0%)	1.01±0.06	0±0/269 (0.0± 0.1%)
All	All	0.56	0/7511 (0.0%)	1.02	1/9953 (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	0.4±0.5
All	All	0	15

There are no bond-length outliers.

All unique angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	16	LEU	CA-CB-CG	-5.05	103.67	115.30	35	1

There are no chirality outliers.

All unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	10	PHE	Sidechain	10
1	A	1	TYR	Sidechain	3
1	A	19	HIS	Sidechain	2

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	199	201	201	1±1
All	All	7400	7437	7437	25

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 2.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:10:PHE:CD1	1:A:16:LEU:HD12	0.67	2.23	34	1
1:A:10:PHE:CD2	1:A:16:LEU:HD23	0.63	2.29	24	6
1:A:10:PHE:CE2	1:A:16:LEU:HD23	0.60	2.32	24	4
1:A:10:PHE:CD1	1:A:16:LEU:CD1	0.49	2.94	34	2
1:A:10:PHE:CZ	1:A:19:HIS:CG	0.48	3.02	2	8
1:A:10:PHE:CD2	1:A:16:LEU:HD12	0.45	2.46	25	1
1:A:10:PHE:CE1	1:A:16:LEU:HD12	0.44	2.47	34	1
1:A:10:PHE:CE2	1:A:19:HIS:CG	0.43	3.05	35	1
1:A:10:PHE:CD1	1:A:16:LEU:CD2	0.41	3.04	10	1

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	24/27 (89%)	20±2 (83±7%)	3±2 (13±6%)	1±1 (4±4%)	4	33
All	All	888/999 (89%)	737 (83%)	119 (13%)	32 (4%)	4	33

All 8 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	7	GLU	15
1	A	11	VAL	7
1	A	24	LYS	3
1	A	6	CYS	2
1	A	8	ARG	2
1	A	20	GLN	1
1	A	23	HIS	1
1	A	12	GLU	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	22/23 (96%)	19±1 (87±6%)	3±1 (13±6%)	5	46
All	All	814/851 (96%)	705 (87%)	109 (13%)	5	46

All 15 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	14	SER	20
1	A	16	LEU	17
1	A	3	CYS	14
1	A	23	HIS	11
1	A	24	LYS	8
1	A	22	VAL	8
1	A	11	VAL	5
1	A	2	LYS	5
1	A	13	LYS	5
1	A	21	ARG	4
1	A	7	GLU	4
1	A	12	GLU	4
1	A	18	ARG	2
1	A	17	SER	1
1	A	20	GLN	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 oligosaccharides 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

No chemical shift data were provided