



# Full wwPDB NMR Structure Validation Report ⓘ

Mar 6, 2022 – 08:55 PM EST

PDB ID : 2OD1  
Title : Solution structure of the MYND domain from human AML1-ETO  
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Deposited on : 2006-12-21

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) 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 : 2.27  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.27

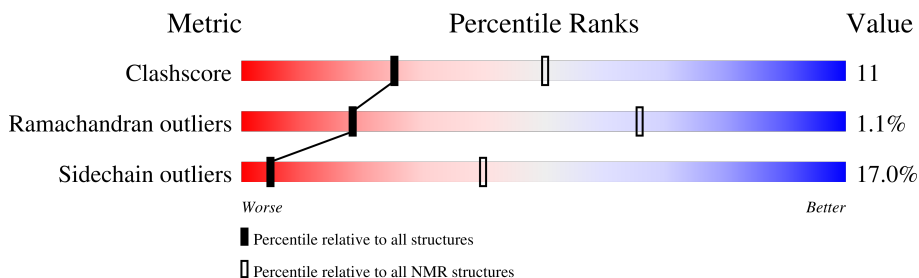
# 1 Overall quality at a glance

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	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  $\geq 3$ , 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	60	

## 2 Ensemble composition and analysis i

This entry contains 31 models. Model 26 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *fewest violations*.

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:662-A:699 (38)	0.07	26

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 5 clusters. No single-model clusters were found.

Cluster number	Models
1	3, 5, 7, 13, 23, 25, 26, 29, 30, 31
2	1, 2, 17, 19, 20, 21, 22, 24
3	4, 6, 8, 9, 12, 16
4	10, 14, 15, 27
5	11, 18, 28

### 3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 732 atoms, of which 340 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Protein CBFA2T1.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
1	A	50	730	230	340	76	77	7	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	648	GLY	-	cloning artifact	UNP Q06455
A	649	SER	-	cloning artifact	UNP Q06455
A	650	PRO	-	cloning artifact	UNP Q06455
A	651	ASN	-	cloning artifact	UNP Q06455
A	652	SER	-	cloning artifact	UNP Q06455
A	653	GLY	-	cloning artifact	UNP Q06455
A	654	SER	-	cloning artifact	UNP Q06455
A	655	PRO	-	cloning artifact	UNP Q06455
A	656	ASN	-	cloning artifact	UNP Q06455
A	657	SER	-	cloning artifact	UNP Q06455

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

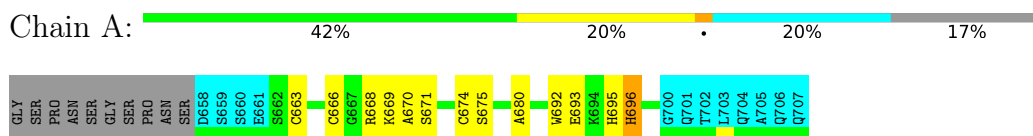
Mol	Chain	Residues	Atoms	
			Total	Zn
2	A	2	2	2

## 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: Protein CBFA2T1

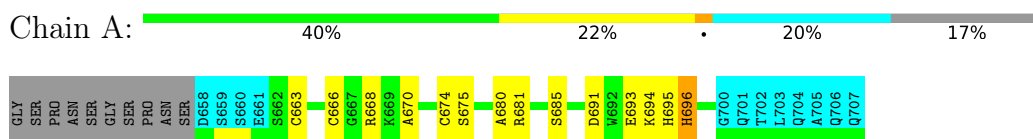


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

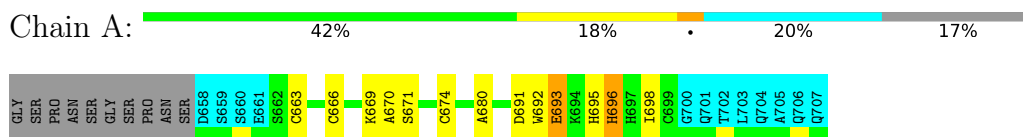
#### 4.2.1 Score per residue for model 1

- Molecule 1: Protein CBFA2T1



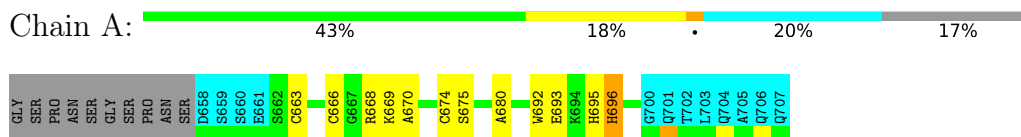
#### 4.2.2 Score per residue for model 2

- Molecule 1: Protein CBFA2T1



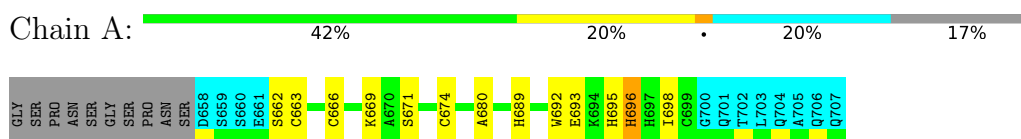
### 4.2.3 Score per residue for model 3

- Molecule 1: Protein CBFA2T1



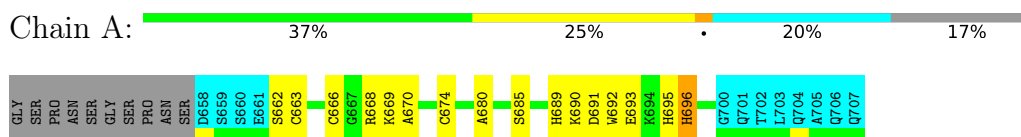
### 4.2.4 Score per residue for model 4

- Molecule 1: Protein CBFA2T1



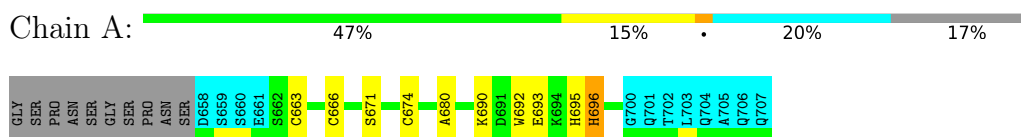
### 4.2.5 Score per residue for model 5

- Molecule 1: Protein CBFA2T1



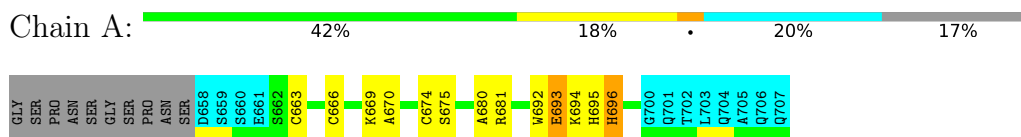
### 4.2.6 Score per residue for model 6

- Molecule 1: Protein CBFA2T1



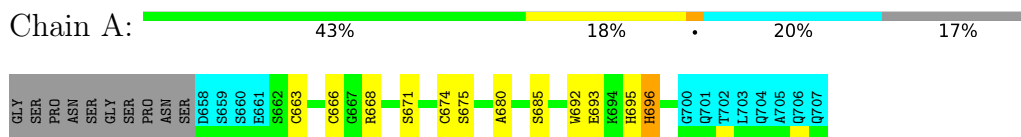
### 4.2.7 Score per residue for model 7

- Molecule 1: Protein CBFA2T1



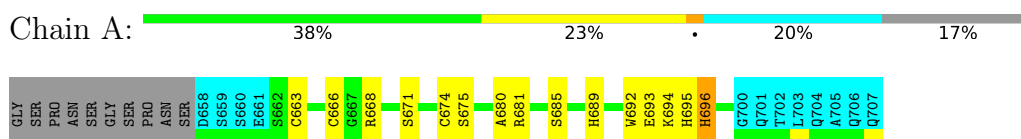
### 4.2.8 Score per residue for model 8

- Molecule 1: Protein CBFA2T1



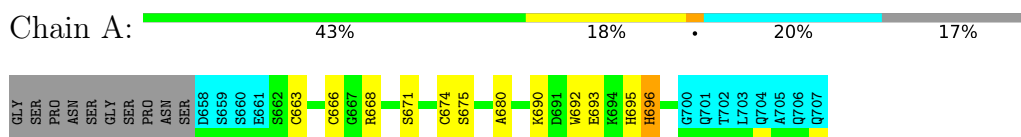
### 4.2.9 Score per residue for model 9

- Molecule 1: Protein CBFA2T1



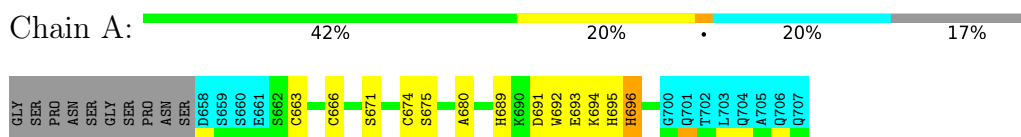
### 4.2.10 Score per residue for model 10

- Molecule 1: Protein CBFA2T1



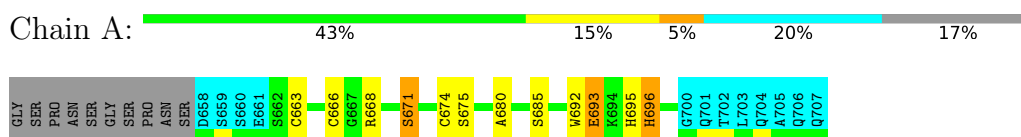
### 4.2.11 Score per residue for model 11

- Molecule 1: Protein CBFA2T1



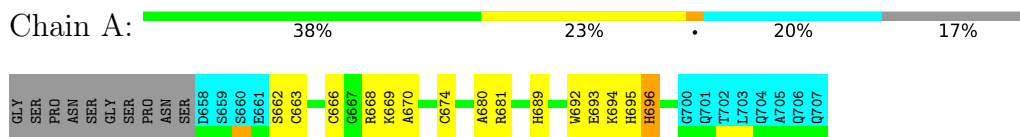
### 4.2.12 Score per residue for model 12

- Molecule 1: Protein CBFA2T1



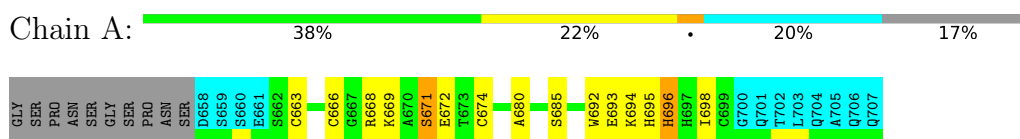
### 4.2.13 Score per residue for model 13

- Molecule 1: Protein CBFA2T1



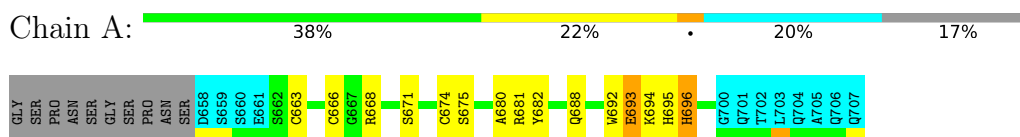
### 4.2.14 Score per residue for model 14

- Molecule 1: Protein CBFA2T1



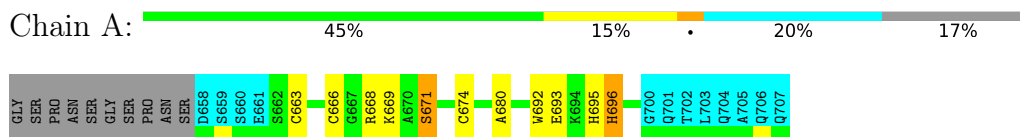
### 4.2.15 Score per residue for model 15

- Molecule 1: Protein CBFA2T1



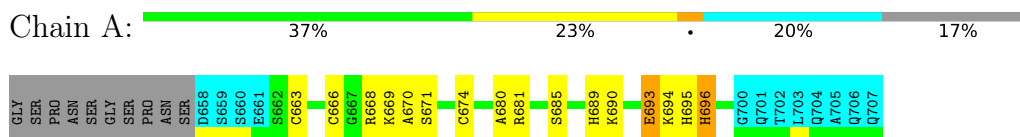
### 4.2.16 Score per residue for model 16

- Molecule 1: Protein CBFA2T1



### 4.2.17 Score per residue for model 17

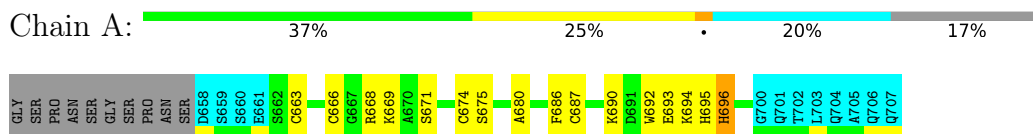
- Molecule 1: Protein CBFA2T1





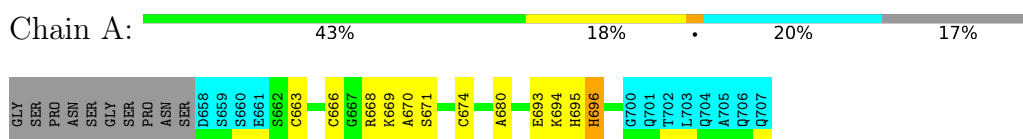
#### 4.2.18 Score per residue for model 18

- Molecule 1: Protein CBFA2T1



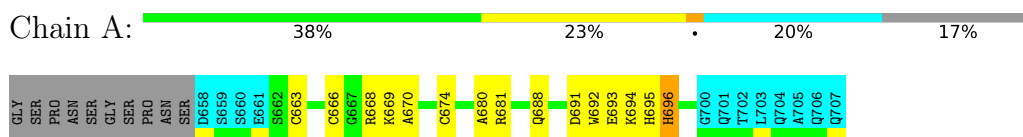
#### 4.2.19 Score per residue for model 19

- Molecule 1: Protein CBFA2T1



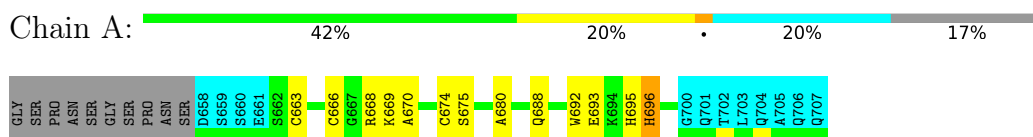
#### 4.2.20 Score per residue for model 20

- Molecule 1: Protein CBFA2T1



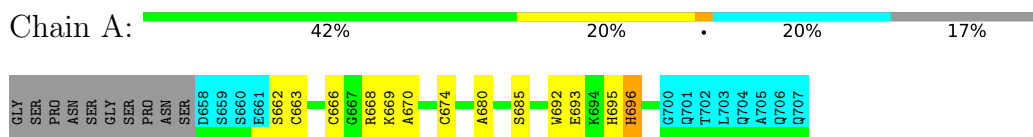
#### 4.2.21 Score per residue for model 21

- Molecule 1: Protein CBFA2T1



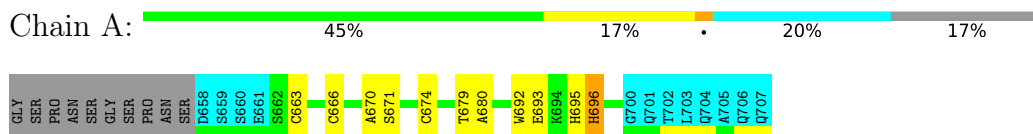
#### 4.2.22 Score per residue for model 22

- Molecule 1: Protein CBFA2T1



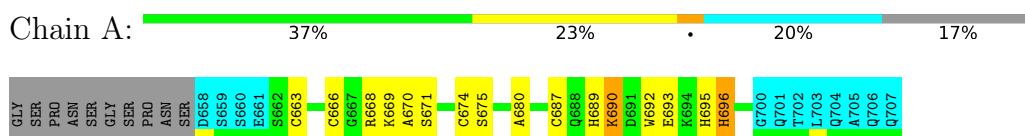
### 4.2.23 Score per residue for model 23

- Molecule 1: Protein CBFA2T1



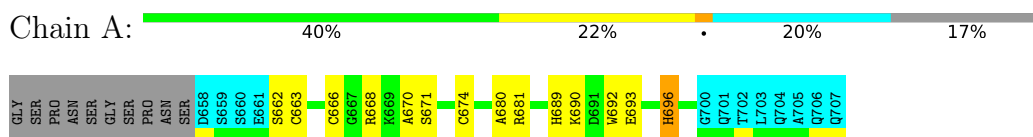
### 4.2.24 Score per residue for model 24

- Molecule 1: Protein CBFA2T1



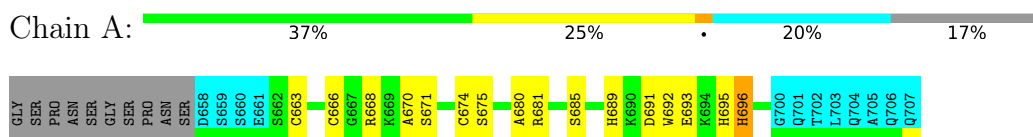
### 4.2.25 Score per residue for model 25

- Molecule 1: Protein CBFA2T1



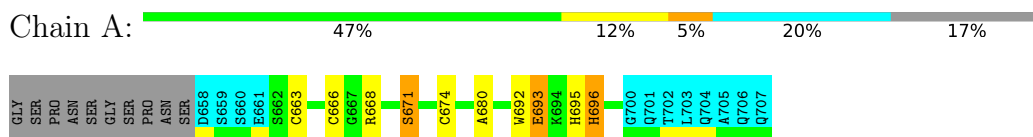
### 4.2.26 Score per residue for model 26 (medoid)

- Molecule 1: Protein CBFA2T1



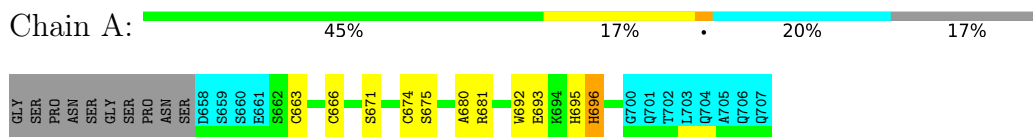
### 4.2.27 Score per residue for model 27

- Molecule 1: Protein CBFA2T1



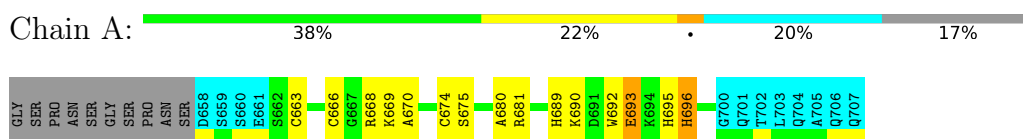
#### 4.2.28 Score per residue for model 28

- Molecule 1: Protein CBFA2T1



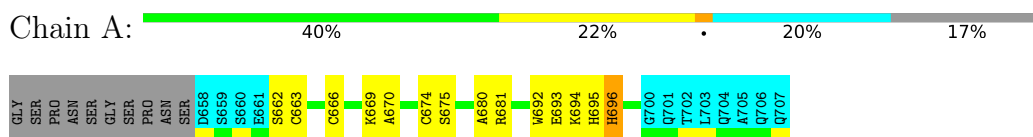
#### 4.2.29 Score per residue for model 29

- Molecule 1: Protein CBFA2T1



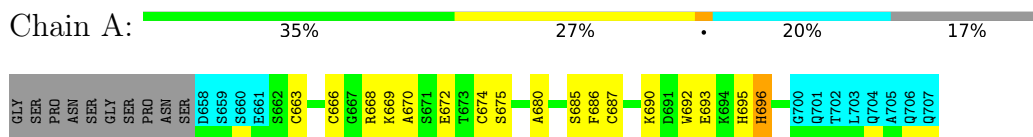
#### 4.2.30 Score per residue for model 30

- Molecule 1: Protein CBFA2T1



#### 4.2.31 Score per residue for model 31

- Molecule 1: Protein CBFA2T1



## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing, torsion angle dynamics*.

Of the 100 calculated structures, 31 were deposited, based on the following criterion: *structures with the least restraint violations*.

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

Software name	Classification	Version
CNS	structure solution	
CNS	refinement	

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

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	301	262	259	6±1
All	All	9393	8122	8029	200

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 11.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:674:CYS:HB2	1:A:680:ALA:HB3	0.71	1.61	27	31
1:A:695:HIS:CE1	1:A:698:ILE:HD11	0.58	2.33	4	3
1:A:663:CYS:HA	1:A:670:ALA:HB2	0.58	1.76	26	18
1:A:680:ALA:HB2	1:A:695:HIS:CE1	0.57	2.34	12	30
1:A:663:CYS:CA	1:A:670:ALA:HB2	0.53	2.34	17	18
1:A:693:GLU:O	1:A:696:HIS:CD2	0.50	2.65	17	3
1:A:663:CYS:HB3	1:A:666:CYS:HB2	0.50	1.84	11	31
1:A:693:GLU:O	1:A:696:HIS:CE1	0.49	2.66	22	13
1:A:698:ILE:HD12	1:A:698:ILE:C	0.47	2.30	2	3
1:A:682:TYR:CD2	1:A:688:GLN:CG	0.46	2.98	15	1
1:A:693:GLU:O	1:A:696:HIS:ND1	0.43	2.51	28	15
1:A:692:TRP:CE3	1:A:692:TRP:O	0.42	2.73	7	24

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:682:TYR:CD2	1:A:688:GLN:HG2	0.41	2.50	15	1
1:A:692:TRP:O	1:A:692:TRP:CE3	0.41	2.74	28	4
1:A:687:CYS:O	1:A:690:LYS:HG2	0.41	2.16	24	1
1:A:671:SER:OG	1:A:672:GLU:N	0.41	2.54	14	1
1:A:688:GLN:O	1:A:692:TRP:CB	0.40	2.69	20	1
1:A:686:PHE:CE2	1:A:687:CYS:SG	0.40	3.15	18	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	38/60 (63%)	34±1 (90±3%)	3±0 (9±1%)	0±0 (1±1%)	18	66
All	All	1178/1860 (63%)	1059 (90%)	106 (9%)	13 (1%)	18	66

All 1 unique Ramachandran outliers are listed below.

Mol	Chain	Res	Type	Models (Total)
1	A	671	SER	13

### 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	33/51 (65%)	27±2 (83±5%)	6±2 (17±5%)	5	40
All	All	1023/1581 (65%)	849 (83%)	174 (17%)	5	40

All 16 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	696	HIS	31
1	A	668	ARG	23
1	A	669	LYS	18
1	A	675	SER	17
1	A	681	ARG	12
1	A	694	LYS	12
1	A	685	SER	10
1	A	671	SER	10
1	A	689	HIS	10
1	A	690	LYS	9
1	A	693	GLU	7
1	A	691	ASP	6
1	A	662	SER	6
1	A	688	GLN	1
1	A	679	THR	1
1	A	672	GLU	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 2 ligands modelled in this entry, 2 are 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