



# Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 09:36 PM BST

PDB ID : 2JYB  
Title : binary hvDHFR1:folate complex  
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Deposited on : 2007-12-10

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  
<http://wwpdb.org/validation/2016/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:

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

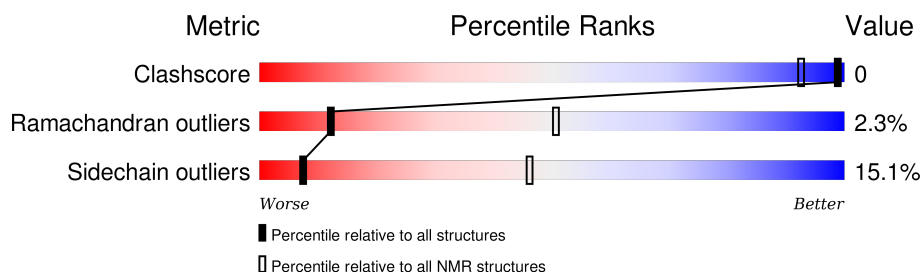
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment 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	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  $\geq 3$ , 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	162	

## 2 Ensemble composition and analysis

This entry contains 20 models. Model 18 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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:10, A:27-A:65, A:71-A:122, A:129-A:158 (131)	0.26	18

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 2 clusters and 12 single-model clusters were found.

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

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2467 atoms, of which 1202 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Dihydrofolate reductase.

Mol	Chain	Residues	Atoms						Trace
1	A	162	Total	C	H	N	O	S	0
			2467	788	1202	218	255	4	

There is a discrepancy between the modelled and reference sequences:

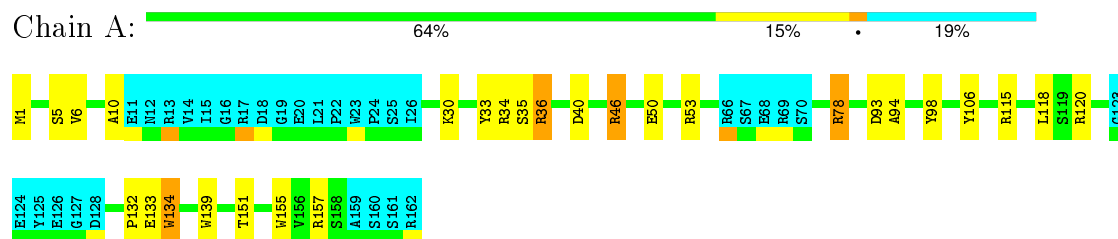
Chain	Residue	Modelled	Actual	Comment	Reference
A	37	VAL	ILE	CONFLICT	UNP P15093

## 4 Residue-property plots [i](#)

### 4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA and DNA 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: Dihydrofolate reductase

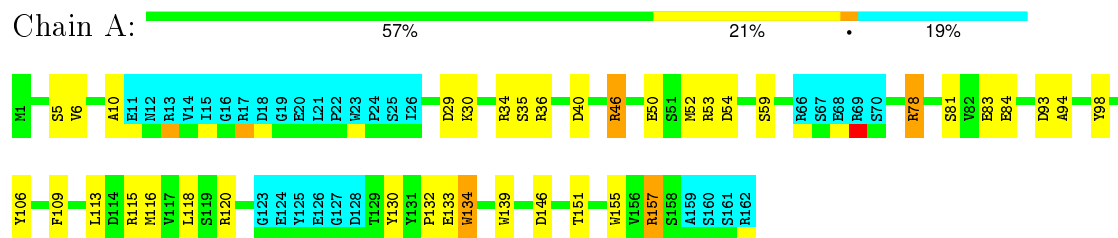


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

#### 4.2.1 Score per residue for model 1

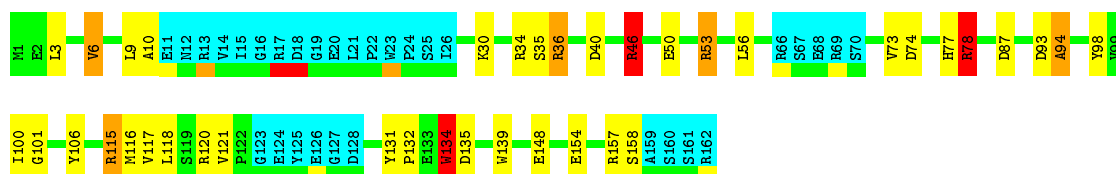
- Molecule 1: Dihydrofolate reductase



#### 4.2.2 Score per residue for model 2

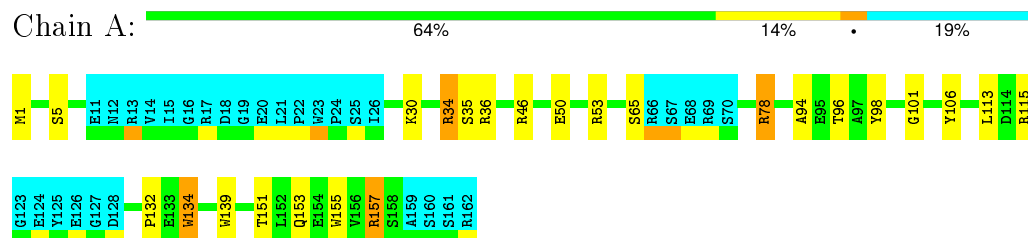
- Molecule 1: Dihydrofolate reductase





#### 4.2.3 Score per residue for model 3

- Molecule 1: Dihydrofolate reductase



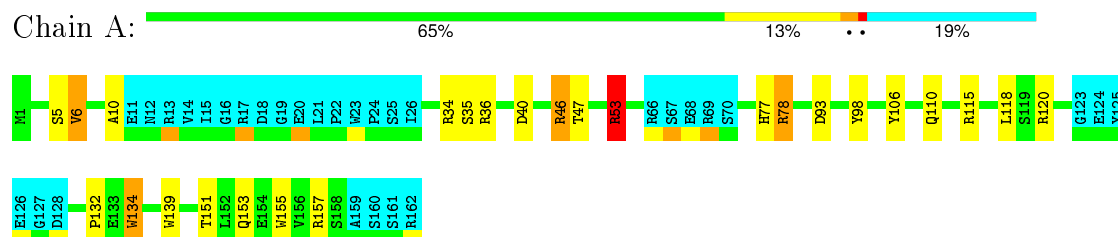
#### 4.2.4 Score per residue for model 4

- Molecule 1: Dihydrofolate reductase



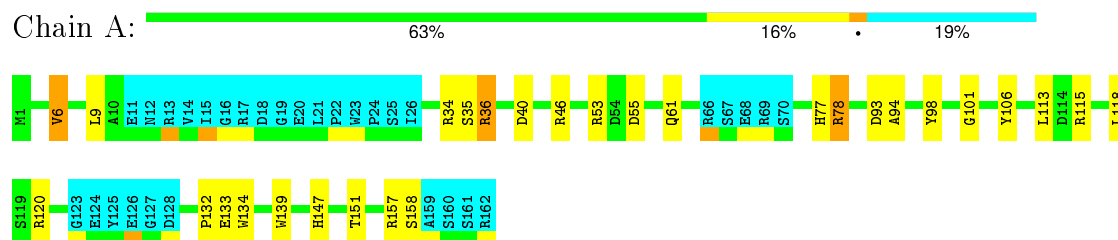
#### 4.2.5 Score per residue for model 5

- Molecule 1: Dihydrofolate reductase



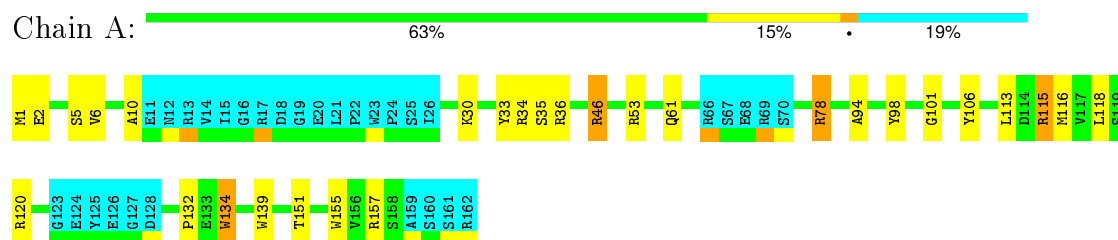
#### 4.2.6 Score per residue for model 6

- Molecule 1: Dihydrofolate reductase



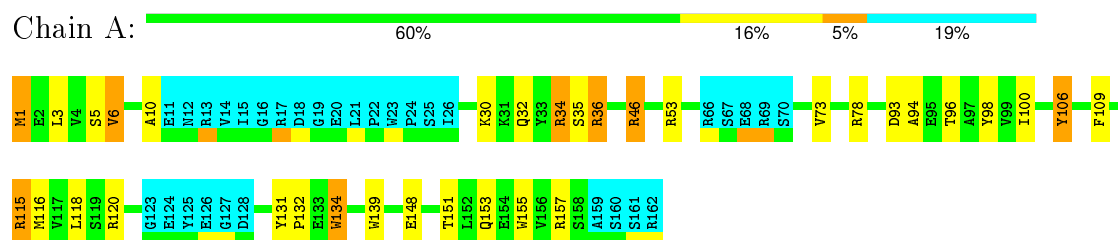
#### 4.2.7 Score per residue for model 7

- Molecule 1: Dihydrofolate reductase



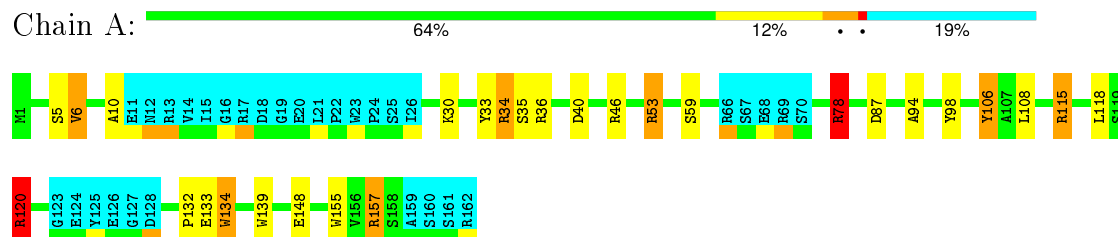
#### 4.2.8 Score per residue for model 8

- Molecule 1: Dihydrofolate reductase



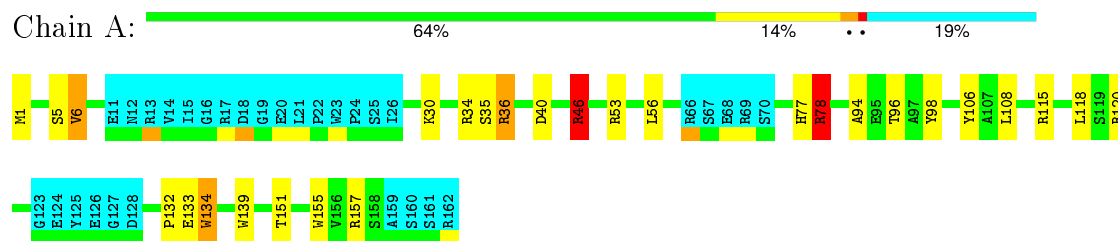
#### 4.2.9 Score per residue for model 9

- Molecule 1: Dihydrofolate reductase



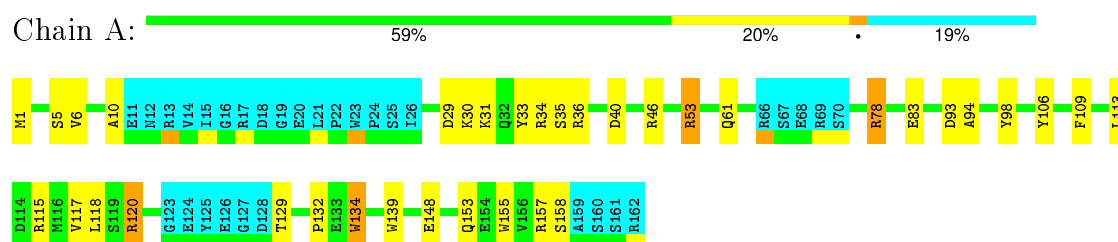
### 4.2.10 Score per residue for model 10

- Molecule 1: Dihydrofolate reductase



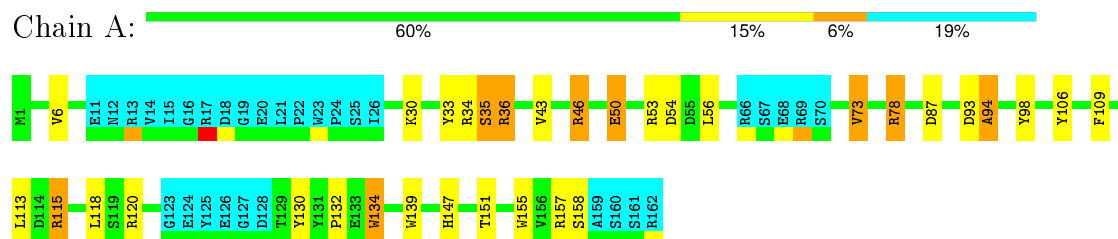
### 4.2.11 Score per residue for model 11

- Molecule 1: Dihydrofolate reductase



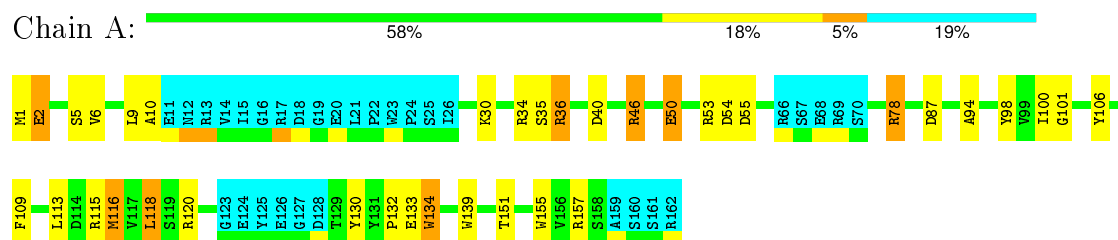
### 4.2.12 Score per residue for model 12

- Molecule 1: Dihydrofolate reductase



### 4.2.13 Score per residue for model 13

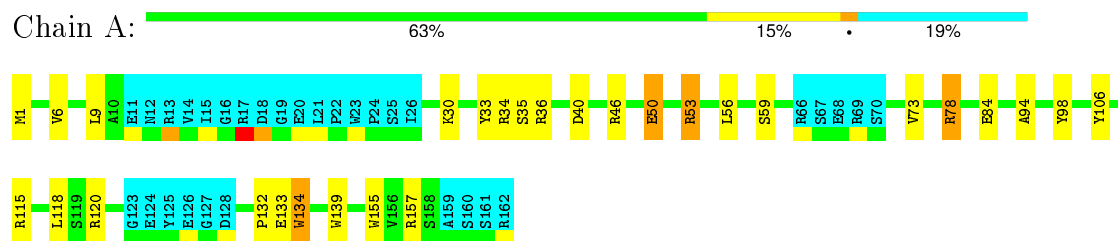
- Molecule 1: Dihydrofolate reductase





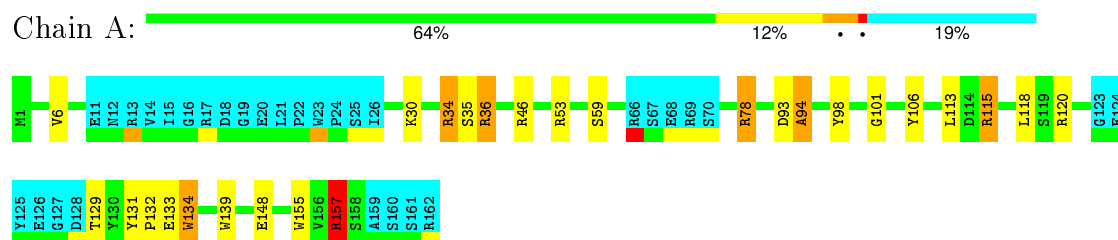
## 4.2.14 Score per residue for model 14

- Molecule 1: Dihydrofolate reductase



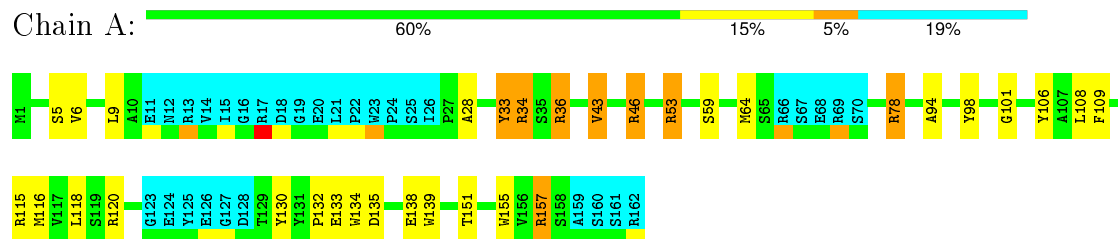
## 4.2.15 Score per residue for model 15

- Molecule 1: Dihydrofolate reductase



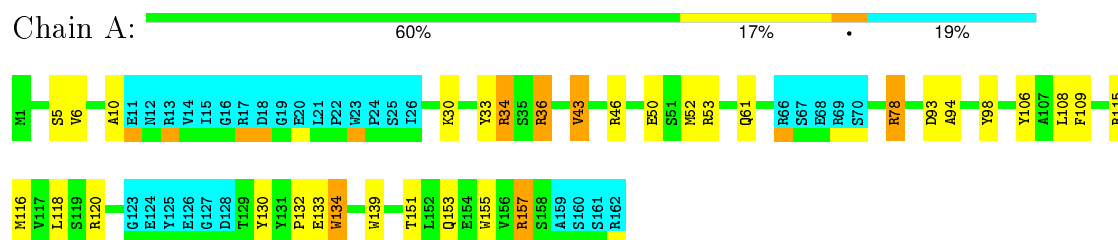
## 4.2.16 Score per residue for model 16

- Molecule 1: Dihydrofolate reductase



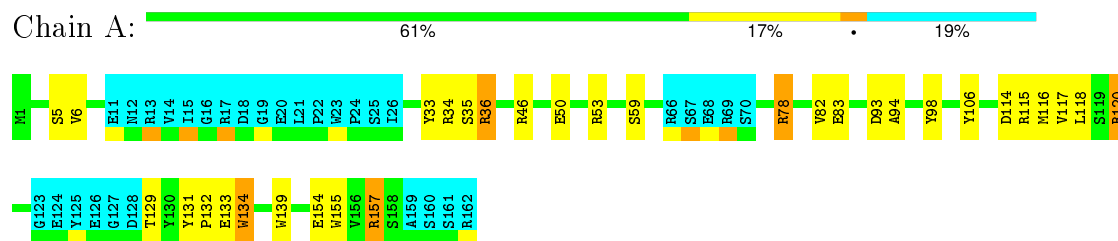
## 4.2.17 Score per residue for model 17

- Molecule 1: Dihydrofolate reductase



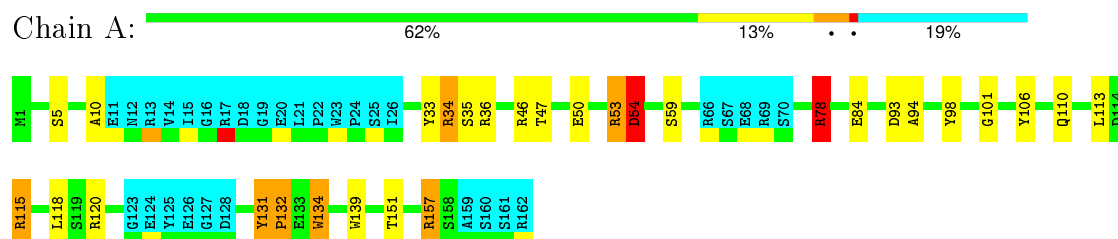
#### 4.2.18 Score per residue for model 18 (medoid)

- Molecule 1: Dihydrofolate reductase



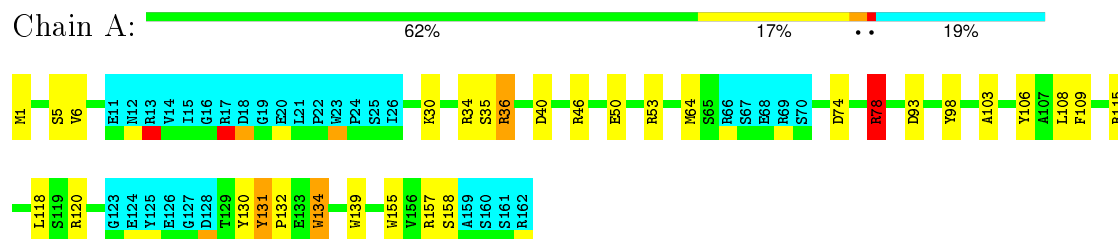
#### 4.2.19 Score per residue for model 19

- Molecule 1: Dihydrofolate reductase



#### 4.2.20 Score per residue for model 20

- Molecule 1: Dihydrofolate reductase



## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing*.

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
CNS	refinement	

No chemical shift data was provided. 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

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	1.32±0.01	0±0/1043 (0.0±0.0%)	1.65±0.01	16±1/1424 (1.1±0.1%)
All	All	1.32	0/20860 (0.0%)	1.65	321/28480 (1.1%)

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	5.5±2.2
All	All	0	111

There are no bond-length outliers.

All unique angle 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(°)	Ideal(°)	Models	
								Worst	Total
1	A	36	ARG	NE-CZ-NH1	11.93	126.26	120.30	20	20
1	A	115	ARG	NE-CZ-NH1	11.92	126.26	120.30	2	20
1	A	78	ARG	NE-CZ-NH1	11.90	126.25	120.30	6	18
1	A	53	ARG	NE-CZ-NH1	10.15	125.38	120.30	17	20
1	A	120	ARG	NE-CZ-NH1	9.89	125.25	120.30	20	20
1	A	157	ARG	NE-CZ-NH1	9.27	124.93	120.30	12	18
1	A	34	ARG	NE-CZ-NH1	9.23	124.92	120.30	15	20
1	A	50	GLU	CA-CB-CG	9.01	133.22	113.40	13	1
1	A	157	ARG	NE-CZ-NH2	-8.84	115.88	120.30	12	6
1	A	115	ARG	NE-CZ-NH2	-8.70	115.95	120.30	2	7
1	A	46	ARG	NE-CZ-NH1	8.61	124.60	120.30	7	19
1	A	120	ARG	NE-CZ-NH2	-8.39	116.10	120.30	14	8
1	A	53	ARG	NE-CZ-NH2	-7.46	116.57	120.30	5	4
1	A	139	TRP	CD1-NE1-CE2	-7.39	102.35	109.00	9	20
1	A	36	ARG	NE-CZ-NH2	-7.38	116.61	120.30	20	10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	46	ARG	NE-CZ-NH2	-6.99	116.81	120.30	16	7
1	A	10	ALA	N-CA-CB	-6.85	100.51	110.10	5	7
1	A	94	ALA	N-CA-CB	-6.69	100.73	110.10	12	3
1	A	78	ARG	NE-CZ-NH2	-6.69	116.96	120.30	5	4
1	A	10	ALA	CB-CA-C	6.45	119.77	110.10	4	5
1	A	33	TYR	CB-CG-CD2	-6.42	117.15	121.00	9	1
1	A	50	GLU	CB-CA-C	6.27	122.93	110.40	12	2
1	A	34	ARG	NE-CZ-NH2	-6.06	117.27	120.30	18	7
1	A	155	TRP	CD1-NE1-CE2	-6.01	103.59	109.00	12	17
1	A	6	VAL	CA-CB-CG2	5.98	119.87	110.90	12	12
1	A	134	TRP	CD1-NE1-CE2	-5.98	103.62	109.00	20	20
1	A	43	VAL	CA-CB-CG2	5.69	119.43	110.90	17	3
1	A	37	VAL	CA-CB-CG2	5.66	119.40	110.90	4	1
1	A	2	GLU	CB-CA-C	5.64	121.67	110.40	13	1
1	A	96	THR	CA-CB-CG2	5.46	120.05	112.40	10	3
1	A	101	GLY	O-C-N	-5.44	113.95	123.20	2	5
1	A	54	ASP	CB-CG-OD2	5.44	123.20	118.30	19	1
1	A	52	MET	CA-CB-CG	5.43	122.54	113.30	1	1
1	A	133	GLU	CB-CA-C	5.39	121.18	110.40	4	2
1	A	121	VAL	CA-CB-CG2	5.30	118.86	110.90	2	1
1	A	151	THR	CA-CB-CG2	5.29	119.81	112.40	8	1
1	A	73	VAL	CA-CB-CG2	5.21	118.72	110.90	14	2
1	A	117	VAL	CA-CB-CG2	5.17	118.66	110.90	11	1
1	A	131	TYR	CB-CG-CD2	-5.11	117.93	121.00	15	1
1	A	6	VAL	CA-CB-CG1	5.10	118.55	110.90	4	1
1	A	157	ARG	N-CA-CB	5.06	119.70	110.60	15	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	35	SER	Peptide	17
1	A	30	LYS	Peptide	11
1	A	115	ARG	Sidechain	7
1	A	130	TYR	Sidechain	7
1	A	78	ARG	Sidechain	7
1	A	50	GLU	Peptide	7
1	A	34	ARG	Sidechain,Peptide	6
1	A	53	ARG	Sidechain	5
1	A	9	LEU	Peptide	4

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Mol	Chain	Res	Type	Group	Models (Total)
1	A	77	HIS	Sidechain	4
1	A	33	TYR	Sidechain	3
1	A	131	TYR	Sidechain	3
1	A	157	ARG	Sidechain	3
1	A	46	ARG	Sidechain	3
1	A	101	GLY	Peptide	3
1	A	147	HIS	Sidechain	2
1	A	120	ARG	Sidechain	2
1	A	100	ILE	Peptide	2
1	A	1	MET	Peptide	2
1	A	106	TYR	Sidechain	2
1	A	36	ARG	Sidechain	2
1	A	103	ALA	Peptide	1
1	A	138	GLU	Peptide	1
1	A	82	VAL	Peptide	1
1	A	109	PHE	Sidechain	1
1	A	28	ALA	Peptide	1
1	A	134	TRP	Peptide	1
1	A	54	ASP	Peptide	1
1	A	148	GLU	Peptide	1

## 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	1019	976	973	0±1
All	All	20380	19520	19460	5

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 0.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:116:MET:SD	1:A:118:LEU:HD12	0.54	2.41	13	1
1:A:117:VAL:HG12	1:A:154:GLU:CG	0.54	2.33	4	2
1:A:117:VAL:HG12	1:A:154:GLU:HG2	0.42	1.91	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:73:VAL:HG12	1:A:74:ASP:H	0.42	1.75	2	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	130/162 (80%)	105±3 (81±3%)	22±3 (17±3%)	3±1 (2±1%)	12	51
All	All	2600/3240 (80%)	2107 (81%)	433 (17%)	60 (2%)	12	51

All 13 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	132	PRO	20
1	A	94	ALA	18
1	A	93	ASP	6
1	A	158	SER	5
1	A	148	GLU	2
1	A	73	VAL	2
1	A	46	ARG	1
1	A	83	GLU	1
1	A	102	GLY	1
1	A	157	ARG	1
1	A	30	LYS	1
1	A	52	MET	1
1	A	35	SER	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	106/132 (80%)	90±3 (85±3%)	16±3 (15±3%)	7	46
All	All	2120/2640 (80%)	1799 (85%)	321 (15%)	7	46

All 56 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	98	TYR	20
1	A	106	TYR	20
1	A	118	LEU	20
1	A	78	ARG	19
1	A	134	TRP	18
1	A	5	SER	15
1	A	6	VAL	12
1	A	151	THR	12
1	A	40	ASP	11
1	A	1	MET	9
1	A	36	ARG	9
1	A	133	GLU	9
1	A	113	LEU	9
1	A	46	ARG	8
1	A	116	MET	8
1	A	109	PHE	7
1	A	59	SER	7
1	A	33	TYR	7
1	A	157	ARG	7
1	A	93	ASP	6
1	A	108	LEU	6
1	A	54	ASP	5
1	A	153	GLN	5
1	A	61	GLN	5
1	A	50	GLU	4
1	A	131	TYR	4
1	A	56	LEU	4
1	A	53	ARG	4
1	A	87	ASP	4
1	A	129	THR	3
1	A	84	GLU	3
1	A	30	LYS	3
1	A	120	ARG	3
1	A	3	LEU	2
1	A	43	VAL	2
1	A	34	ARG	2

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Mol	Chain	Res	Type	Models (Total)
1	A	55	ASP	2
1	A	64	MET	2
1	A	2	GLU	2
1	A	29	ASP	2
1	A	110	GLN	2
1	A	47	THR	2
1	A	148	GLU	2
1	A	135	ASP	2
1	A	83	GLU	2
1	A	100	ILE	1
1	A	114	ASP	1
1	A	32	GLN	1
1	A	65	SER	1
1	A	74	ASP	1
1	A	132	PRO	1
1	A	31	LYS	1
1	A	146	ASP	1
1	A	81	SER	1
1	A	9	LEU	1
1	A	115	ARG	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 ⓘ

There are no carbohydrates in this entry.

### 6.6 Ligand geometry ⓘ

There are no ligands in this entry.

## 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