



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 04:04 AM UTC

PDB ID : 2KDV / pdb_00002kdv
Title : Solution structure of RNA Pyrophosphohydrolase RppH from Escherichia coli
Authors : Bi, Y.; Li, H.; Jin, C.
Deposited on : 2009-01-19

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

<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 : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
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) : 2.49

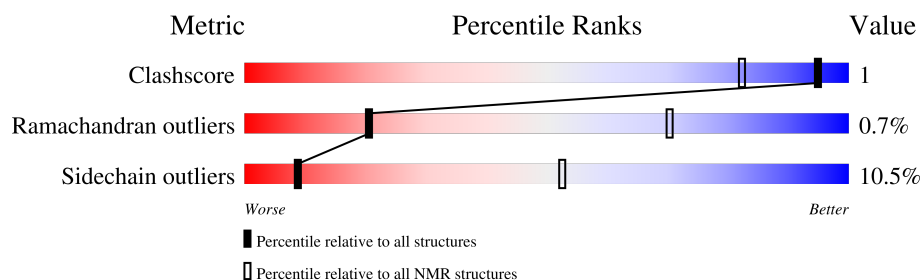
1 Overall quality at a glance

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	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

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	164	 88% 9% .

2 Ensemble composition and analysis

This entry contains 21 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:2-A:160 (159)	0.56	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 2 clusters and 3 single-model clusters were found.

Cluster number	Models
1	1, 2, 3, 4, 5, 6, 7, 9, 11, 13, 14, 15, 17, 18, 19, 20
2	8, 12
Single-model clusters	10; 16; 21

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 2722 atoms, of which 1353 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called RNA pyrophosphohydrolase.

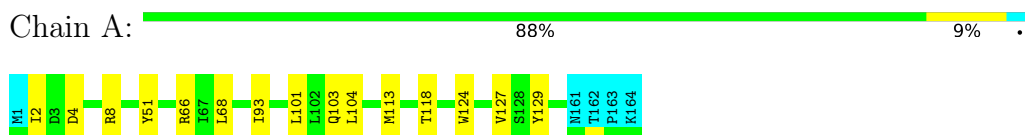
Mol	Chain	Residues	Atoms						Trace
1	A	164	Total	C	H	N	O	S	0
			2722	873	1353	249	239	8	

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: RNA pyrophosphohydrolase

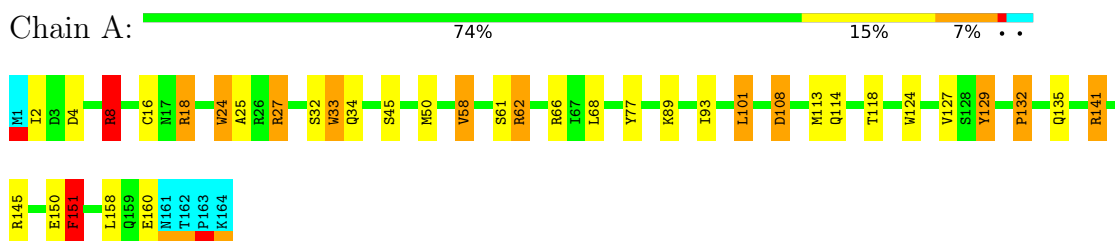


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

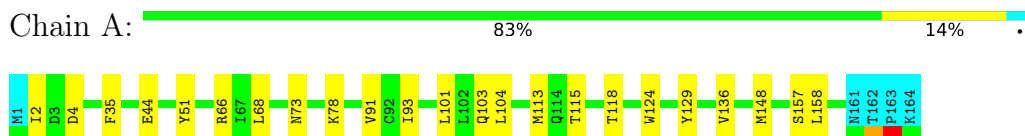
4.2.1 Score per residue for model 1 (medoid)

- Molecule 1: RNA pyrophosphohydrolase



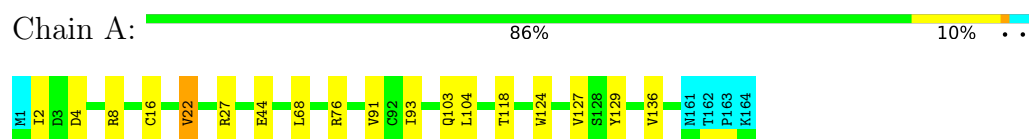
4.2.2 Score per residue for model 2

- Molecule 1: RNA pyrophosphohydrolase



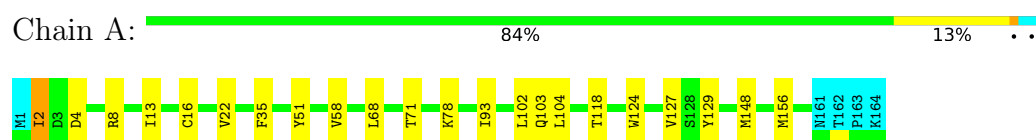
4.2.3 Score per residue for model 3

- Molecule 1: RNA pyrophosphohydrolase



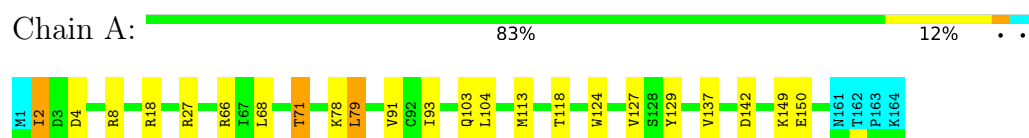
4.2.4 Score per residue for model 4

- Molecule 1: RNA pyrophosphohydrolase



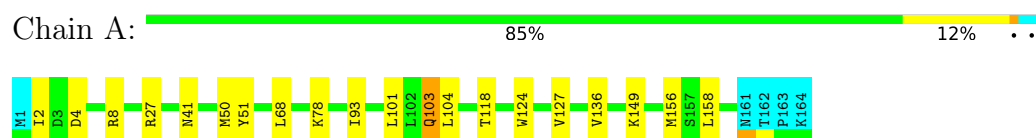
4.2.5 Score per residue for model 5

- Molecule 1: RNA pyrophosphohydrolase



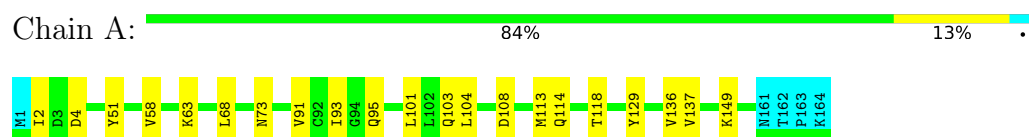
4.2.6 Score per residue for model 6

- Molecule 1: RNA pyrophosphohydrolase



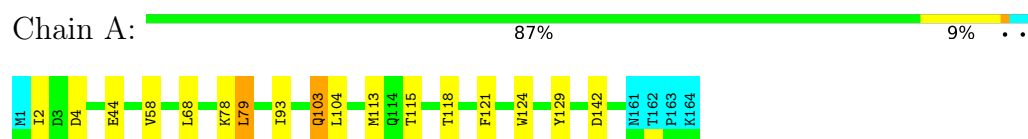
4.2.7 Score per residue for model 7

- Molecule 1: RNA pyrophosphohydrolase



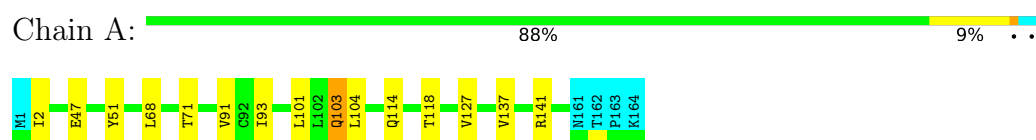
4.2.8 Score per residue for model 8

- Molecule 1: RNA pyrophosphohydrolase



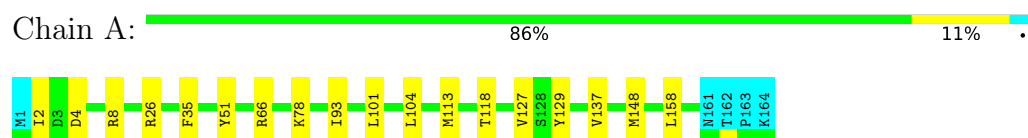
4.2.9 Score per residue for model 9

- Molecule 1: RNA pyrophosphohydrolase



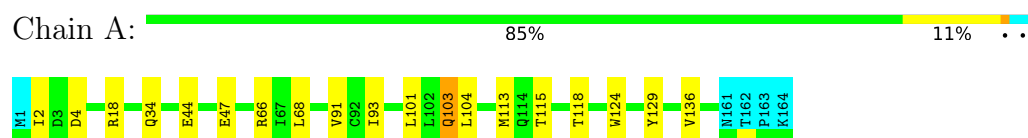
4.2.10 Score per residue for model 10

- Molecule 1: RNA pyrophosphohydrolase



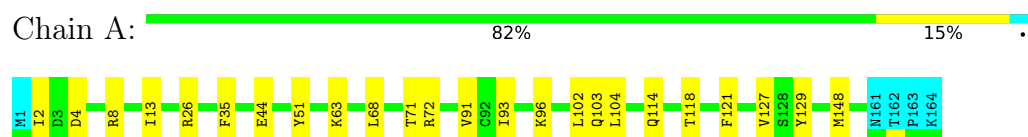
4.2.11 Score per residue for model 11

- Molecule 1: RNA pyrophosphohydrolase



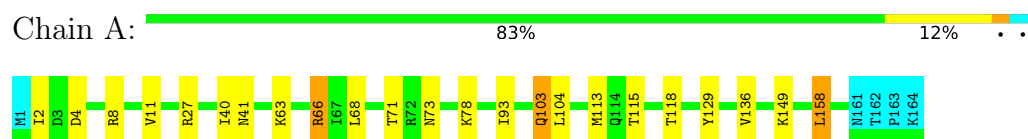
4.2.12 Score per residue for model 12

- Molecule 1: RNA pyrophosphohydrolase



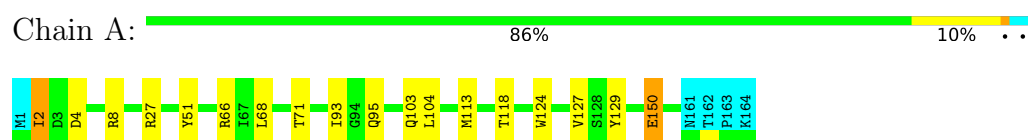
4.2.13 Score per residue for model 13

- Molecule 1: RNA pyrophosphohydrolase



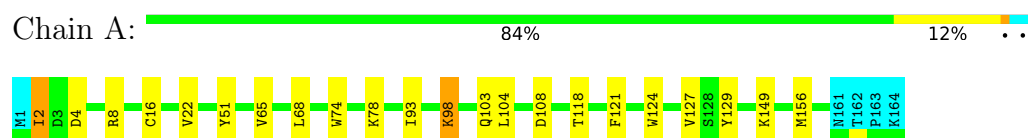
4.2.14 Score per residue for model 14

- Molecule 1: RNA pyrophosphohydrolase



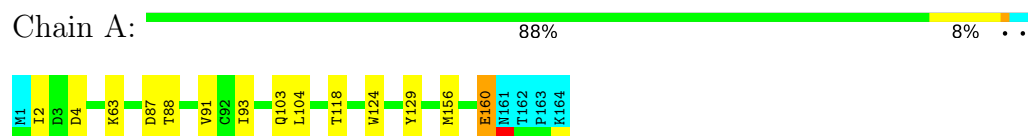
4.2.15 Score per residue for model 15

- Molecule 1: RNA pyrophosphohydrolase



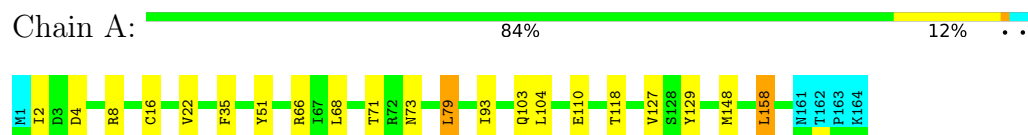
4.2.16 Score per residue for model 16

- Molecule 1: RNA pyrophosphohydrolase



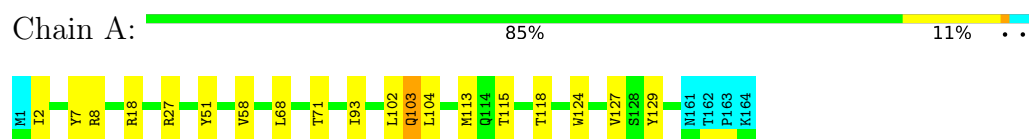
4.2.17 Score per residue for model 17

- Molecule 1: RNA pyrophosphohydrolase



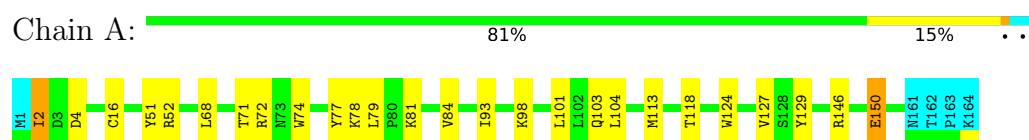
4.2.18 Score per residue for model 18

- Molecule 1: RNA pyrophosphohydrolase



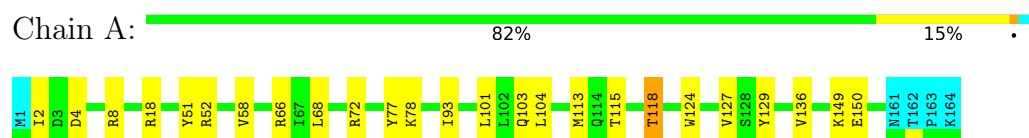
4.2.19 Score per residue for model 19

- Molecule 1: RNA pyrophosphohydrolase



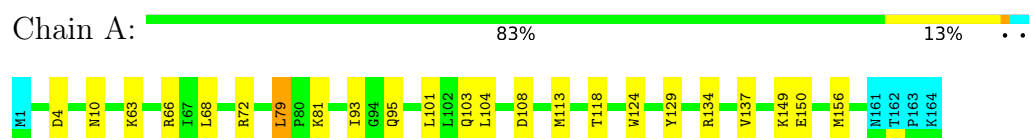
4.2.20 Score per residue for model 20

- Molecule 1: RNA pyrophosphohydrolase



4.2.21 Score per residue for model 21

- Molecule 1: RNA pyrophosphohydrolase



5 Refinement protocol and experimental data overview ⓘ

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

Of the 100 calculated structures, 21 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
Amber	refinement	
CYANA	structure solution	

No chemical shift data was provided.

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	0.69±0.02	0±0/1366 (0.0± 0.0%)	1.29±0.10	2±6/1848 (0.1± 0.3%)
All	All	0.69	2/28686 (0.0%)	1.30	51/38808 (0.1%)

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	1.5±1.5
All	All	0	32

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	24	TRP	NE1-CE2	-6.92	1.29	1.37	1	1
1	A	24	TRP	CD2-CE3	-5.43	1.31	1.40	1	1

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	24	TRP	CZ3-CH2-CZ2	-13.71	103.67	121.50	1	1
1	A	24	TRP	CE2-CD2-CE3	-12.26	106.54	118.80	1	1
1	A	33	TRP	CA-CB-CG	11.66	135.75	113.60	1	1
1	A	151	PHE	CD1-CG-CD2	-9.39	104.52	118.60	1	1
1	A	24	TRP	CH2-CZ2-CE2	-8.96	105.86	117.50	1	1
1	A	8	ARG	NE-CZ-NH2	-8.85	111.24	119.20	1	1
1	A	8	ARG	CD-NE-CZ	8.42	136.19	124.40	1	1
1	A	24	TRP	NE1-CE2-CD2	-8.29	96.62	107.40	1	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	24	TRP	CB-CG-CD2	8.05	138.07	126.80	1	1
1	A	108	ASP	CA-CB-CG	7.82	120.42	112.60	1	1
1	A	160	GLU	CA-C-N	7.71	136.26	121.54	16	1
1	A	160	GLU	C-N-CA	7.71	136.26	121.54	16	1
1	A	79	LEU	CB-CA-C	7.56	119.74	108.86	21	5
1	A	141	ARG	CD-NE-CZ	-6.93	114.70	124.40	1	1
1	A	141	ARG	CB-CG-CD	6.85	127.05	111.30	1	1
1	A	27	ARG	NE-CZ-NH2	-6.50	113.35	119.20	1	1
1	A	89	LYS	CG-CD-CE	-6.43	96.52	111.30	1	1
1	A	157	SER	CA-C-N	6.29	128.71	120.28	2	1
1	A	157	SER	C-N-CA	6.29	128.71	120.28	2	1
1	A	132	PRO	CA-CB-CG	6.28	116.42	104.50	1	1
1	A	18	ARG	NE-CZ-NH2	-6.08	113.73	119.20	1	1
1	A	33	TRP	N-CA-CB	-6.06	100.71	111.08	1	1
1	A	145	ARG	NE-CZ-NH2	-6.03	113.77	119.20	1	1
1	A	62	ARG	CB-CG-CD	-5.97	97.58	111.30	1	1
1	A	121	PHE	CA-CB-CG	5.94	119.74	113.80	12	2
1	A	18	ARG	NE-CZ-NH1	5.87	127.37	121.50	1	1
1	A	18	ARG	CD-NE-CZ	5.85	132.59	124.40	1	1
1	A	121	PHE	N-CA-CB	-5.72	102.11	111.66	8	1
1	A	58	VAL	CA-CB-CG2	5.68	120.06	110.40	1	1
1	A	114	GLN	CA-C-N	5.60	128.24	120.63	9	3
1	A	114	GLN	C-N-CA	5.60	128.24	120.63	9	3
1	A	88	THR	CA-CB-OG1	5.49	117.83	109.60	16	1
1	A	62	ARG	CD-NE-CZ	-5.35	116.91	124.40	1	1
1	A	142	ASP	CA-CB-CG	-5.34	107.26	112.60	5	2
1	A	24	TRP	CD2-CE2-CZ2	5.33	127.73	122.40	1	1
1	A	118	THR	CB-CA-C	5.33	116.87	108.88	20	1
1	A	8	ARG	NE-CZ-NH1	5.28	126.78	121.50	1	1
1	A	151	PHE	CG-CD1-CE1	-5.24	111.79	120.70	1	1
1	A	104	LEU	CB-CA-C	5.03	117.94	109.89	10	1
1	A	145	ARG	CD-NE-CZ	5.03	131.44	124.40	1	1
1	A	24	TRP	CE2-CD2-CG	5.00	113.20	107.20	1	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	51	TYR	Sidechain	12
1	A	27	ARG	Sidechain	7

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Mol	Chain	Res	Type	Group	Models (Total)
1	A	77	TYR	Sidechain	3
1	A	8	ARG	Sidechain	1
1	A	62	ARG	Sidechain	1
1	A	129	TYR	Sidechain	1
1	A	141	ARG	Sidechain	1
1	A	151	PHE	Sidechain	1
1	A	18	ARG	Sidechain	1
1	A	26	ARG	Sidechain	1
1	A	66	ARG	Sidechain	1
1	A	52	ARG	Sidechain	1
1	A	146	ARG	Sidechain	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	1329	1309	1309	3±3
All	All	27909	27489	27489	59

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 1.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:24:TRP:HE1	1:A:33:TRP:HB3	1.04	1.11	1	1
1:A:24:TRP:NE1	1:A:33:TRP:HB3	0.94	1.77	1	1
1:A:24:TRP:CE3	1:A:132:PRO:HB3	0.89	2.02	1	1
1:A:16:CYS:SG	1:A:101:LEU:HD11	0.80	2.17	1	2
1:A:24:TRP:HE3	1:A:132:PRO:HB3	0.80	1.32	1	1
1:A:24:TRP:CE3	1:A:132:PRO:CB	0.64	2.81	1	1
1:A:24:TRP:CZ3	1:A:132:PRO:CB	0.61	2.83	1	1
1:A:101:LEU:HB2	1:A:151:PHE:HD2	0.61	1.54	1	1
1:A:68:LEU:HD11	1:A:103:GLN:HB2	0.61	1.72	13	18
1:A:24:TRP:CE2	1:A:33:TRP:HB3	0.58	2.32	1	1
1:A:68:LEU:HB2	1:A:101:LEU:HD23	0.51	1.82	1	1
1:A:150:GLU:HB3	1:A:151:PHE:CE1	0.51	2.41	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:24:TRP:C	1:A:24:TRP:CD1	0.50	2.89	1	1
1:A:68:LEU:HD11	1:A:103:GLN:CB	0.49	2.38	17	4
1:A:16:CYS:SG	1:A:22:VAL:HG12	0.49	2.48	17	4
1:A:74:TRP:CE2	1:A:98:LYS:HE3	0.46	2.45	15	2
1:A:24:TRP:HD1	1:A:25:ALA:N	0.45	2.10	1	1
1:A:74:TRP:CD2	1:A:98:LYS:HE3	0.44	2.48	15	1
1:A:71:THR:HG22	1:A:150:GLU:HG3	0.44	1.89	5	2
1:A:24:TRP:CD1	1:A:25:ALA:N	0.44	2.85	1	1
1:A:11:VAL:HG13	1:A:40:ILE:HD11	0.44	1.90	13	1
1:A:35:PHE:CZ	1:A:148:MET:HE3	0.44	2.47	17	5
1:A:71:THR:HG22	1:A:150:GLU:HG2	0.43	1.90	14	1
1:A:68:LEU:HD13	1:A:158:LEU:HG	0.42	1.90	17	2
1:A:50:MET:C	1:A:50:MET:SD	0.41	3.03	1	1
1:A:81:LYS:HA	1:A:84:VAL:HG23	0.41	1.92	19	1
1:A:24:TRP:HE1	1:A:33:TRP:CB	0.40	2.04	1	1
1:A:50:MET:HE1	1:A:51:TYR:CE1	0.40	2.51	6	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	159/164 (97%)	153±1 (96±1%)	5±1 (3±1%)	1±0 (1±0%)	20	70
All	All	3339/3444 (97%)	3214 (96%)	100 (3%)	25 (1%)	20	70

All 3 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	4	ASP	19
1	A	2	ILE	5
1	A	7	TYR	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	142/147 (97%)	127±3 (89±2%)	15±3 (11±2%)	9	52
All	All	2982/3087 (97%)	2668 (89%)	314 (11%)	9	52

All 55 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	93	ILE	21
1	A	118	THR	21
1	A	2	ILE	20
1	A	129	TYR	19
1	A	104	LEU	19
1	A	124	TRP	15
1	A	127	VAL	14
1	A	8	ARG	13
1	A	113	MET	13
1	A	66	ARG	10
1	A	78	LYS	10
1	A	101	LEU	9
1	A	91	VAL	8
1	A	136	VAL	7
1	A	71	THR	7
1	A	149	LYS	7
1	A	103	GLN	7
1	A	58	VAL	6
1	A	115	THR	6
1	A	44	GLU	5
1	A	158	LEU	5
1	A	156	MET	5
1	A	137	VAL	5
1	A	63	LYS	5
1	A	18	ARG	4
1	A	108	ASP	4
1	A	73	ASN	4
1	A	79	LEU	4
1	A	72	ARG	4

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Mol	Chain	Res	Type	Models (Total)
1	A	150	GLU	4
1	A	102	LEU	3
1	A	95	GLN	3
1	A	34	GLN	2
1	A	13	ILE	2
1	A	41	ASN	2
1	A	47	GLU	2
1	A	32	SER	1
1	A	45	SER	1
1	A	61	SER	1
1	A	114	GLN	1
1	A	135	GLN	1
1	A	151	PHE	1
1	A	22	VAL	1
1	A	76	ARG	1
1	A	141	ARG	1
1	A	26	ARG	1
1	A	96	LYS	1
1	A	65	VAL	1
1	A	98	LYS	1
1	A	87	ASP	1
1	A	110	GLU	1
1	A	52	ARG	1
1	A	10	ASN	1
1	A	81	LYS	1
1	A	134	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 oligosaccharides in this entry.

6.6 Ligand geometry

There are no ligands in this entry.

6.7 Other polymers

There are no such molecules in this entry.

6.8 Polymer linkage issues

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided