



Full wwPDB NMR Structure Validation Report ⓘ

Mar 6, 2026 – 06:20 PM UTC

PDB ID : 2KW4 / pdb_00002kw4
Title : Solution NMR Structure of the Holo Form of a Ribonuclease H domain from D.hafniense, Northeast Structural Genomics Consortium Target DhR1A
Authors : Mills, J.L.; Eletsky, A.; Hua, J.; Belote, R.L.; Buchwald, W.A.; Ciccocanti, C.; Janjua, H.; Nair, R.; Rost, B.; Acton, T.B.; Xiao, R.; Everett, J.K.; Montelione, G.T.; Szyperski, T.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2010-03-31

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

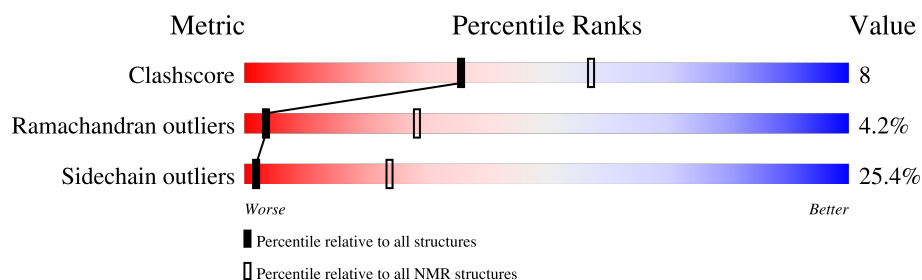
1 Overall quality at a glance

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	147	

2 Ensemble composition and analysis

This entry contains 20 models. Model 8 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:3-A:73, A:78-A:86, A:95-A:137 (123)	0.98	8

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 and 2 single-model clusters were found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2294 atoms, of which 1114 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms						Trace
1	A	147	Total	C	H	N	O	S	0
			2293	750	1114	207	218	4	

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q24WL3
A	140	LEU	-	expression tag	UNP Q24WL3
A	141	GLU	-	expression tag	UNP Q24WL3
A	142	HIS	-	expression tag	UNP Q24WL3
A	143	HIS	-	expression tag	UNP Q24WL3
A	144	HIS	-	expression tag	UNP Q24WL3
A	145	HIS	-	expression tag	UNP Q24WL3
A	146	HIS	-	expression tag	UNP Q24WL3
A	147	HIS	-	expression tag	UNP Q24WL3

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

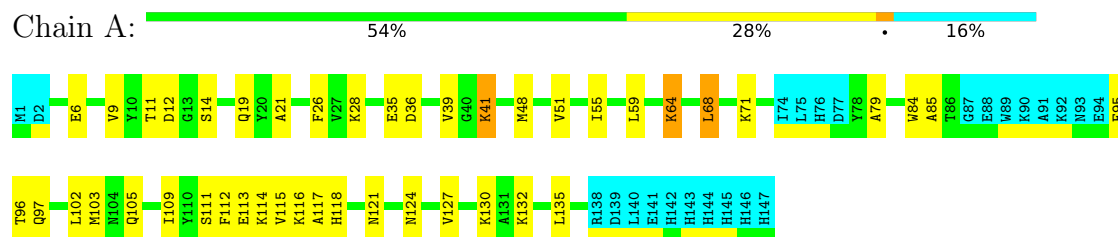
Mol	Chain	Residues	Atoms	
2	A	1	Total	Mg
			1	1

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: Uncharacterized protein



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

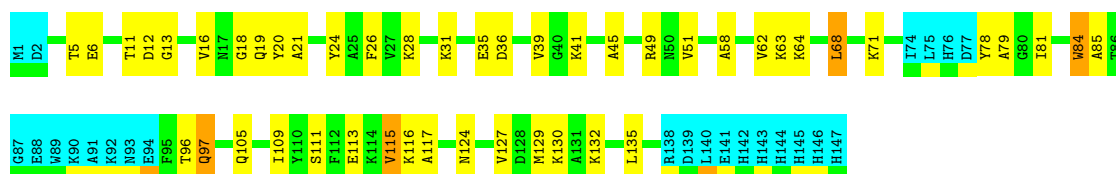
- Molecule 1: Uncharacterized protein



4.2.2 Score per residue for model 2

- Molecule 1: Uncharacterized protein

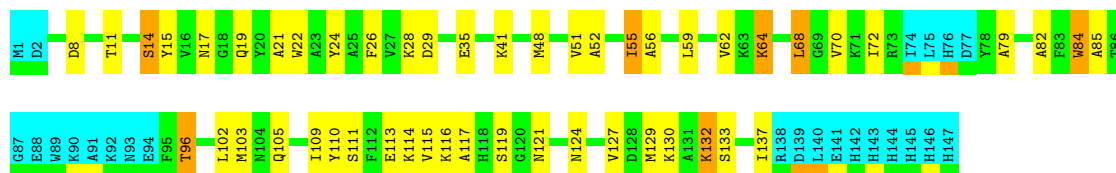




4.2.3 Score per residue for model 3

- Molecule 1: Uncharacterized protein

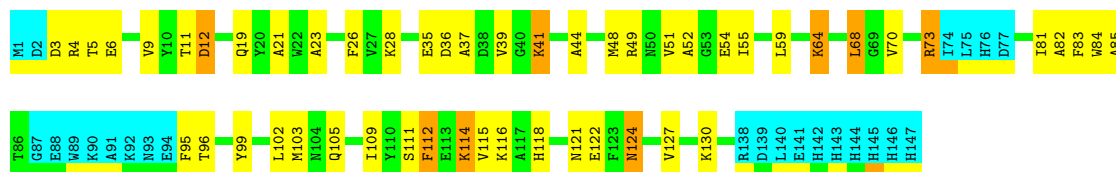
Chain A:



4.2.4 Score per residue for model 4

- Molecule 1: Uncharacterized protein

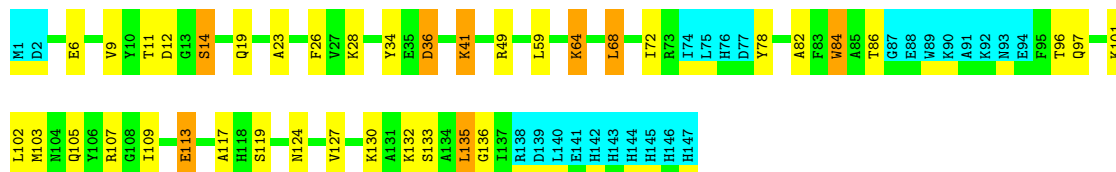
Chain A:



4.2.5 Score per residue for model 5

- Molecule 1: Uncharacterized protein

Chain A:



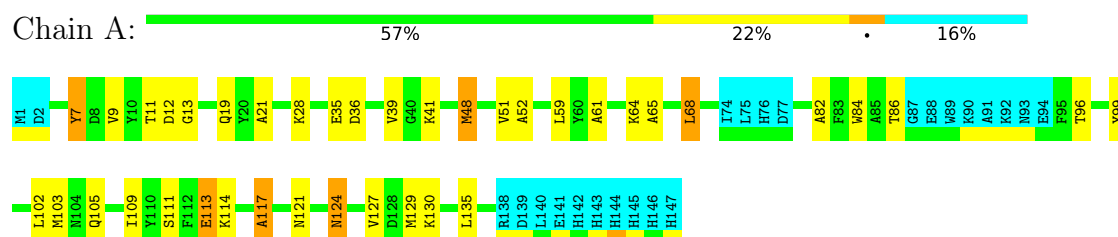
4.2.6 Score per residue for model 6

- Molecule 1: Uncharacterized protein



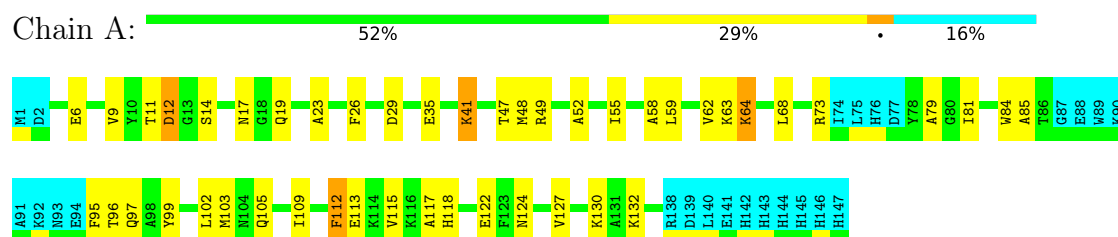
4.2.7 Score per residue for model 7

- Molecule 1: Uncharacterized protein



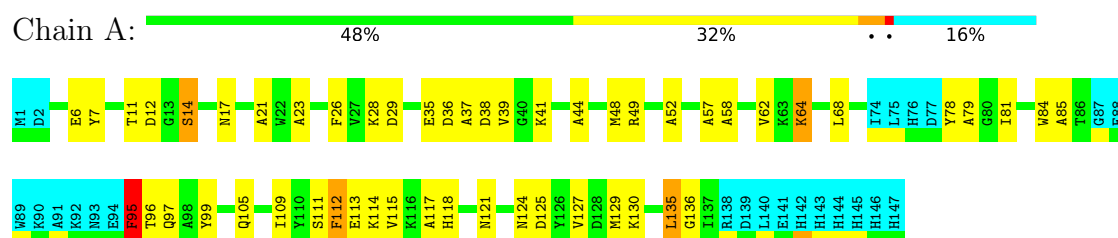
4.2.8 Score per residue for model 8 (medoid)

- Molecule 1: Uncharacterized protein



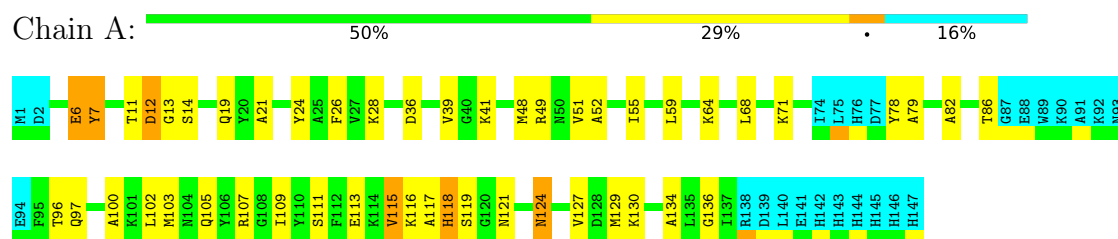
4.2.9 Score per residue for model 9

- Molecule 1: Uncharacterized protein



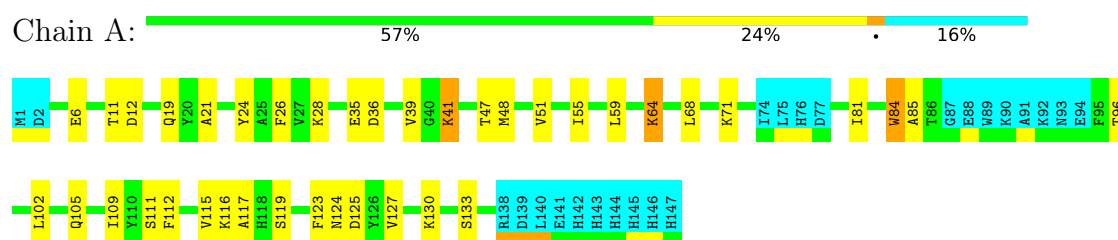
4.2.10 Score per residue for model 10

- Molecule 1: Uncharacterized protein



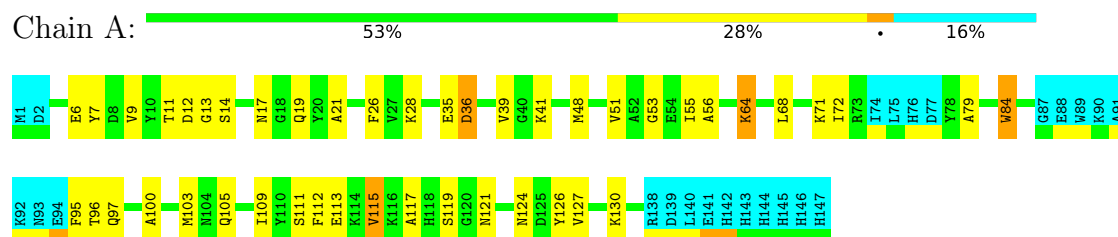
4.2.11 Score per residue for model 11

- Molecule 1: Uncharacterized protein



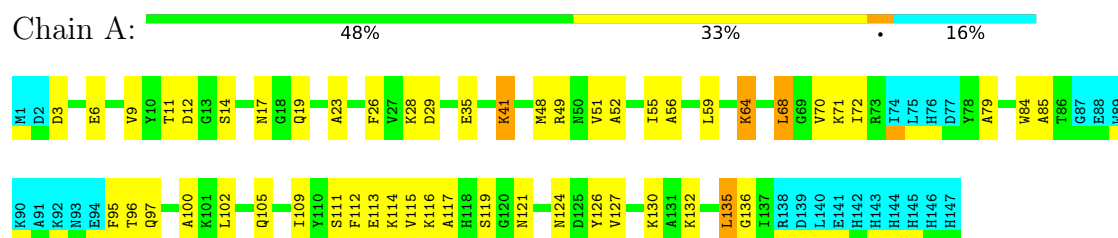
4.2.12 Score per residue for model 12

- Molecule 1: Uncharacterized protein



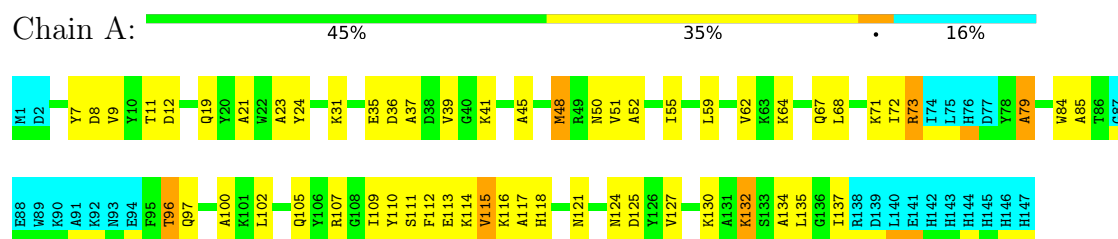
4.2.13 Score per residue for model 13

- Molecule 1: Uncharacterized protein



4.2.14 Score per residue for model 14

- Molecule 1: Uncharacterized protein



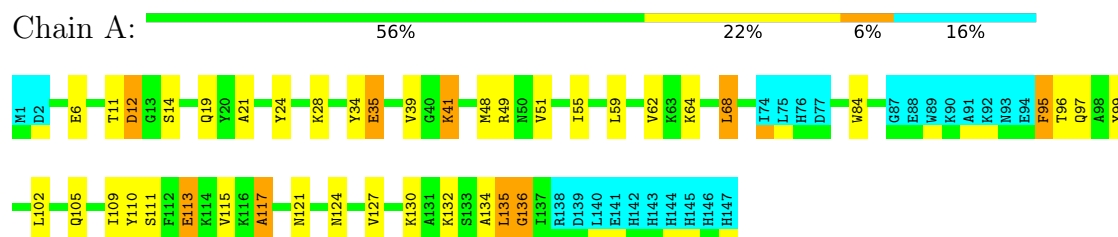
4.2.15 Score per residue for model 15

- Molecule 1: Uncharacterized protein



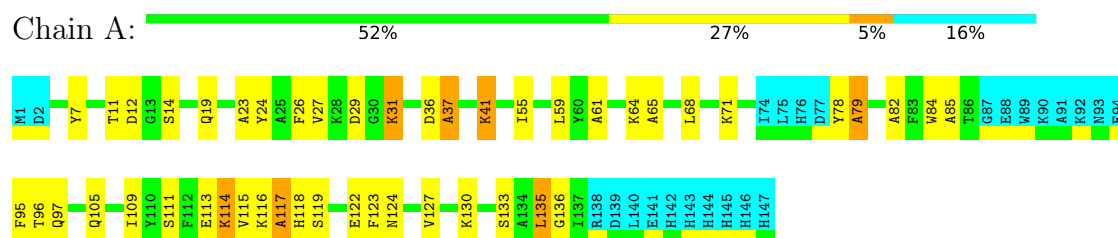
4.2.16 Score per residue for model 16

- Molecule 1: Uncharacterized protein



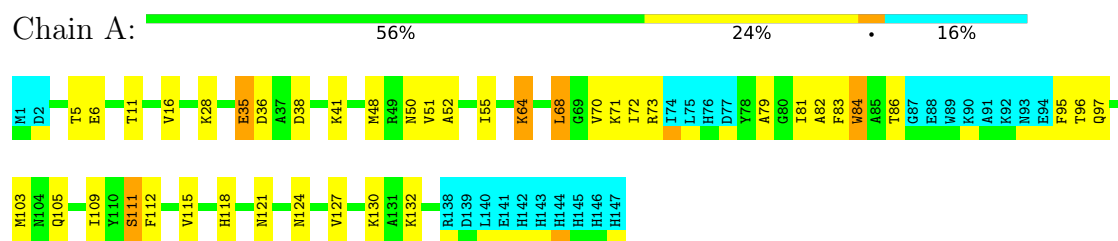
4.2.17 Score per residue for model 17

- Molecule 1: Uncharacterized protein



4.2.18 Score per residue for model 18

- Molecule 1: Uncharacterized protein



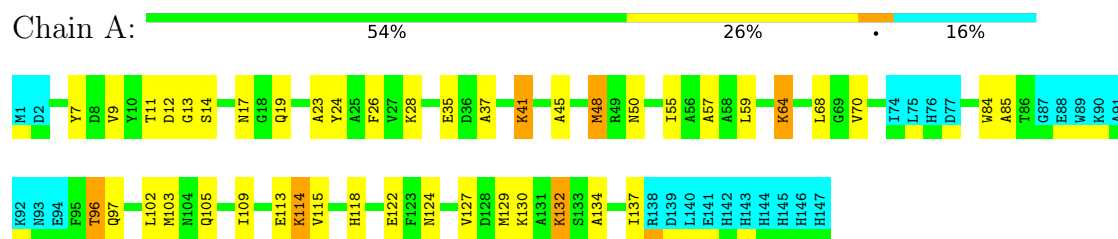
4.2.19 Score per residue for model 19

- Molecule 1: Uncharacterized protein



4.2.20 Score per residue for model 20

- Molecule 1: Uncharacterized protein



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 least restraint violations*.

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

Software name	Classification	Version
MOLMOL	refinement	
TALOS	geometry optimization	

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: MG

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	966	930	930	15±3
All	All	19340	18600	18600	297

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 8.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:21:ALA:HB2	1:A:39:VAL:HG12	0.86	1.48	14	8
1:A:23:ALA:HB2	1:A:37:ALA:HB2	0.83	1.50	20	5
1:A:124:ASN:O	1:A:127:VAL:HG22	0.71	1.84	7	20
1:A:11:THR:O	1:A:127:VAL:HG21	0.69	1.87	4	20
1:A:99:TYR:CD2	1:A:115:VAL:HG21	0.68	2.24	19	1
1:A:64:LYS:O	1:A:68:LEU:HD12	0.66	1.89	10	20
1:A:51:VAL:O	1:A:55:ILE:HD12	0.63	1.94	13	8
1:A:84:TRP:CE3	1:A:115:VAL:HG21	0.63	2.28	20	2
1:A:59:LEU:HD13	1:A:102:LEU:HD13	0.63	1.71	16	9
1:A:7:TYR:CD2	1:A:134:ALA:HB2	0.62	2.29	20	1
1:A:8:ASP:OD2	1:A:79:ALA:HB2	0.62	1.94	14	2
1:A:11:THR:HG22	1:A:57:ALA:HB1	0.61	1.72	9	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:21:ALA:CB	1:A:39:VAL:HG12	0.61	2.24	14	6
1:A:59:LEU:HD13	1:A:102:LEU:CD1	0.60	2.27	11	13
1:A:85:ALA:HB1	1:A:114:LYS:O	0.60	1.96	15	7
1:A:45:ALA:O	1:A:48:MET:HE1	0.60	1.97	20	2
1:A:8:ASP:OD2	1:A:127:VAL:HG12	0.60	1.97	3	1
1:A:85:ALA:HA	1:A:115:VAL:HG23	0.60	1.72	20	2
1:A:100:ALA:HB2	1:A:115:VAL:O	0.59	1.97	19	3
1:A:48:MET:HE2	1:A:52:ALA:HB2	0.59	1.75	10	1
1:A:84:TRP:CZ2	1:A:96:THR:HG22	0.59	2.33	20	1
1:A:82:ALA:O	1:A:86:THR:HG23	0.58	1.98	7	4
1:A:96:THR:HG21	1:A:118:HIS:CD2	0.58	2.33	14	1
1:A:78:TYR:HA	1:A:82:ALA:HB3	0.57	1.74	5	1
1:A:7:TYR:HB2	1:A:134:ALA:HB2	0.57	1.74	14	1
1:A:26:PHE:CE2	1:A:68:LEU:HD11	0.57	2.34	1	12
1:A:68:LEU:O	1:A:70:VAL:HG13	0.57	2.00	3	5
1:A:9:VAL:HG12	1:A:11:THR:HG23	0.57	1.75	20	11
1:A:48:MET:HE2	1:A:51:VAL:HG21	0.57	1.75	7	1
1:A:3:ASP:CG	1:A:5:THR:HG23	0.56	2.26	4	1
1:A:21:ALA:HB2	1:A:39:VAL:HG13	0.56	1.77	9	3
1:A:44:ALA:HB1	1:A:95:PHE:CE1	0.55	2.36	4	1
1:A:15:TYR:HB2	1:A:21:ALA:HB3	0.54	1.79	3	2
1:A:23:ALA:HB2	1:A:37:ALA:CB	0.54	2.32	14	5
1:A:81:ILE:HG21	1:A:112:PHE:CD2	0.54	2.38	9	2
1:A:81:ILE:O	1:A:85:ALA:HB2	0.53	2.04	11	3
1:A:132:LYS:HG2	1:A:137:ILE:HD12	0.53	1.79	20	3
1:A:81:ILE:HG21	1:A:112:PHE:CD1	0.52	2.39	19	2
1:A:48:MET:O	1:A:51:VAL:HG23	0.52	2.04	19	5
1:A:135:LEU:HD23	1:A:136:GLY:N	0.52	2.19	13	7
1:A:51:VAL:HG13	1:A:119:SER:O	0.52	2.04	3	1
1:A:52:ALA:HB1	1:A:99:TYR:OH	0.52	2.03	8	4
1:A:72:ILE:HD12	1:A:73:ARG:NH1	0.51	2.20	18	1
1:A:12:ASP:CG	1:A:81:ILE:HD11	0.51	2.30	1	1
1:A:79:ALA:HB3	1:A:82:ALA:HB3	0.51	1.83	17	1
1:A:44:ALA:HB1	1:A:95:PHE:CE2	0.50	2.41	9	1
1:A:9:VAL:HG21	1:A:72:ILE:HG22	0.50	1.83	12	1
1:A:62:VAL:HG21	1:A:110:TYR:CE1	0.50	2.41	16	3
1:A:132:LYS:HA	1:A:135:LEU:HD12	0.50	1.84	14	3
1:A:23:ALA:HB2	1:A:37:ALA:HB1	0.49	1.83	9	1
1:A:48:MET:HE3	1:A:52:ALA:HB3	0.49	1.83	14	1
1:A:58:ALA:O	1:A:62:VAL:HG23	0.49	2.08	6	4
1:A:52:ALA:O	1:A:56:ALA:HB2	0.49	2.07	19	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:47:THR:HG23	1:A:48:MET:HE2	0.48	1.85	11	1
1:A:7:TYR:CE2	1:A:134:ALA:HB2	0.48	2.43	20	1
1:A:6:GLU:HB3	1:A:134:ALA:HB2	0.48	1.85	16	1
1:A:85:ALA:CA	1:A:115:VAL:HG23	0.48	2.38	4	2
1:A:51:VAL:HG13	1:A:119:SER:HB3	0.48	1.85	6	1
1:A:44:ALA:HB1	1:A:95:PHE:CZ	0.47	2.44	4	1
1:A:9:VAL:HG21	1:A:72:ILE:CG2	0.47	2.40	12	1
1:A:85:ALA:HA	1:A:115:VAL:HG12	0.47	1.85	2	1
1:A:14:SER:O	1:A:21:ALA:HB3	0.47	2.08	9	1
1:A:84:TRP:O	1:A:84:TRP:CD1	0.47	2.68	11	1
1:A:59:LEU:HD11	1:A:99:TYR:CD1	0.47	2.45	4	1
1:A:55:ILE:HD13	1:A:96:THR:HG23	0.47	1.87	3	1
1:A:11:THR:HG22	1:A:23:ALA:O	0.47	2.10	6	4
1:A:48:MET:HE2	1:A:95:PHE:N	0.47	2.25	12	1
1:A:48:MET:CG	1:A:52:ALA:HB2	0.46	2.41	18	1
1:A:100:ALA:HB1	1:A:115:VAL:HG23	0.46	1.87	10	1
1:A:100:ALA:HB2	1:A:115:VAL:HG23	0.45	1.88	13	2
1:A:116:LYS:O	1:A:117:ALA:HB2	0.45	2.10	11	3
1:A:23:ALA:CB	1:A:37:ALA:HB2	0.45	2.35	20	1
1:A:48:MET:HE2	1:A:51:VAL:CG2	0.45	2.42	7	1
1:A:55:ILE:HG22	1:A:59:LEU:HD21	0.45	1.89	17	1
1:A:45:ALA:C	1:A:48:MET:HE1	0.45	2.36	20	1
1:A:84:TRP:HE1	1:A:117:ALA:N	0.45	2.10	5	2
1:A:59:LEU:HD13	1:A:102:LEU:HD12	0.45	1.89	5	1
1:A:24:TYR:N	1:A:24:TYR:CD1	0.45	2.85	10	1
1:A:85:ALA:HB2	1:A:115:VAL:HG12	0.44	1.89	13	1
1:A:55:ILE:HD13	1:A:99:TYR:CD2	0.44	2.46	8	1
1:A:11:THR:HG22	1:A:24:TYR:HB3	0.44	1.90	16	1
1:A:43:PRO:O	1:A:47:THR:HG22	0.44	2.12	19	1
1:A:22:TRP:HB3	1:A:56:ALA:HB3	0.44	1.90	3	2
1:A:99:TYR:CE2	1:A:115:VAL:HG21	0.44	2.48	16	1
1:A:126:TYR:CD1	1:A:126:TYR:C	0.43	2.96	12	3
1:A:85:ALA:HB2	1:A:115:VAL:HG13	0.43	1.91	1	1
1:A:61:ALA:O	1:A:65:ALA:HB2	0.43	2.14	7	2
1:A:20:TYR:CZ	1:A:45:ALA:HB1	0.42	2.49	2	1
1:A:53:GLY:O	1:A:56:ALA:HB3	0.42	2.14	12	1
1:A:47:THR:HG21	1:A:95:PHE:CZ	0.42	2.49	8	1
1:A:82:ALA:HB2	1:A:112:PHE:HE1	0.42	1.75	6	1
1:A:48:MET:CE	1:A:52:ALA:HB3	0.42	2.44	14	1
1:A:16:VAL:O	1:A:16:VAL:HG13	0.42	2.15	2	3
1:A:84:TRP:CD1	1:A:84:TRP:C	0.41	2.93	11	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:135:LEU:HD22	1:A:137:ILE:HD11	0.41	1.93	1	1
1:A:48:MET:HE2	1:A:51:VAL:HB	0.41	1.92	19	1
1:A:48:MET:HG3	1:A:52:ALA:HB2	0.41	1.92	18	1
1:A:84:TRP:CE3	1:A:115:VAL:HG12	0.41	2.51	19	1
1:A:27:VAL:HG13	1:A:31:LYS:O	0.40	2.16	17	1
1:A:81:ILE:HG23	1:A:84:TRP:HZ3	0.40	1.76	18	1
1:A:6:GLU:HB2	1:A:134:ALA:HB2	0.40	1.94	10	1
1:A:3:ASP:OD1	1:A:70:VAL:HG12	0.40	2.15	13	1
1:A:81:ILE:HG23	1:A:84:TRP:CZ3	0.40	2.51	2	1
1:A:11:THR:O	1:A:12:ASP:C	0.40	2.63	17	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	123/147 (84%)	95±2 (78±2%)	22±2 (18±2%)	5±2 (4±2%)	3	28
All	All	2460/2940 (84%)	1908 (78%)	448 (18%)	104 (4%)	3	28

All 25 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	41	LYS	12
1	A	12	ASP	8
1	A	79	ALA	8
1	A	6	GLU	8
1	A	35	GLU	7
1	A	49	ARG	7
1	A	17	ASN	7
1	A	13	GLY	6
1	A	117	ALA	6
1	A	14	SER	5
1	A	113	GLU	4
1	A	7	TYR	4

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Mol	Chain	Res	Type	Models (Total)
1	A	136	GLY	3
1	A	95	PHE	3
1	A	78	TYR	2
1	A	114	LYS	2
1	A	36	ASP	2
1	A	50	ASN	2
1	A	37	ALA	2
1	A	116	LYS	1
1	A	18	GLY	1
1	A	119	SER	1
1	A	73	ARG	1
1	A	118	HIS	1
1	A	5	THR	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	92/114 (81%)	69±2 (75±3%)	23±2 (25±3%)	2	24
All	All	1840/2280 (81%)	1372 (75%)	468 (25%)	2	24

All 60 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	96	THR	20
1	A	105	GLN	20
1	A	109	ILE	20
1	A	130	LYS	20
1	A	19	GLN	18
1	A	41	LYS	18
1	A	84	TRP	16
1	A	28	LYS	16
1	A	113	GLU	15
1	A	12	ASP	14
1	A	97	GLN	14
1	A	36	ASP	13

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Mol	Chain	Res	Type	Models (Total)
1	A	111	SER	13
1	A	115	VAL	12
1	A	132	LYS	12
1	A	35	GLU	12
1	A	68	LEU	11
1	A	103	MET	11
1	A	71	LYS	11
1	A	64	LYS	11
1	A	121	ASN	11
1	A	14	SER	10
1	A	112	PHE	9
1	A	48	MET	9
1	A	135	LEU	8
1	A	118	HIS	8
1	A	116	LYS	7
1	A	24	TYR	7
1	A	129	MET	7
1	A	114	LYS	6
1	A	119	SER	6
1	A	6	GLU	6
1	A	29	ASP	6
1	A	133	SER	5
1	A	73	ARG	5
1	A	122	GLU	5
1	A	107	ARG	5
1	A	31	LYS	4
1	A	55	ILE	4
1	A	95	PHE	4
1	A	78	TYR	3
1	A	83	PHE	3
1	A	124	ASN	3
1	A	34	TYR	3
1	A	49	ARG	3
1	A	7	TYR	3
1	A	125	ASP	3
1	A	63	LYS	2
1	A	3	ASP	2
1	A	26	PHE	2
1	A	38	ASP	2
1	A	123	PHE	2
1	A	5	THR	1
1	A	4	ARG	1

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Mol	Chain	Res	Type	Models (Total)
1	A	54	GLU	1
1	A	101	LYS	1
1	A	47	THR	1
1	A	67	GLN	1
1	A	106	TYR	1
1	A	50	ASN	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