



Full wwPDB NMR Structure Validation Report ⓘ

Mar 5, 2026 – 04:38 PM UTC

PDB ID : 1X9A / pdb_00001x9a
Title : Solution NMR Structure of Protein Tm0979 from *Thermotoga maritima*.
Ontario Center for Structural Proteomics Target TM0979_1_87; Northeast
Structural Genomics Consortium Target VT98.
Authors : Gaspar, J.A.; Liu, C.; Vassall, K.A.; Stathopoulos, P.B.; Meglei, G.; Stephen,
R.; Pineda-Lucena, A.; Wu, B.; Yee, A.; Arrowsmith, C.H.; Meiering, E.M.;
Northeast Structural Genomics Consortium (NESG)
Deposited on : 2004-08-20

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

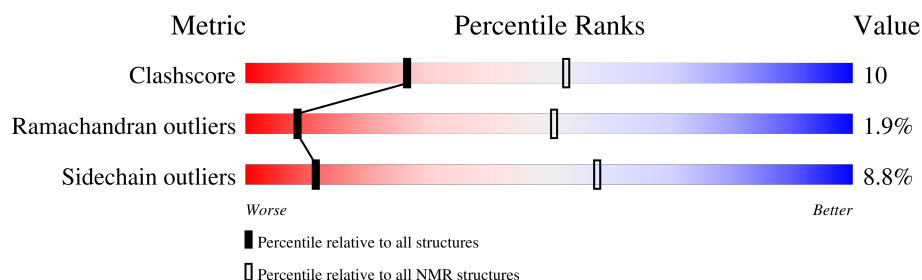
1 Overall quality at a glance

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	107	
1	B	107	

2 Ensemble composition and analysis

This entry contains 10 models. Model 3 is the overall representative, medoid model (most similar to other models). The authors have identified model 9 as representative.

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:87, B:201-B:280 (166)	0.99	3

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 3 clusters. No single-model clusters were found.

Cluster number	Models
1	2, 4, 7, 9
2	3, 6, 8
3	1, 5, 10

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2812 atoms, of which 1416 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called hypothetical protein TM0979.

Mol	Chain	Residues	Atoms						Trace
1	A	87	Total	C	H	N	O	S	0
			1406	455	708	107	135	1	
1	B	87	Total	C	H	N	O	S	0
			1406	455	708	107	135	1	

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q9X074
A	-18	GLY	-	expression tag	UNP Q9X074
A	-17	SER	-	expression tag	UNP Q9X074
A	-16	SER	-	expression tag	UNP Q9X074
A	-15	HIS	-	expression tag	UNP Q9X074
A	-14	HIS	-	expression tag	UNP Q9X074
A	-13	HIS	-	expression tag	UNP Q9X074
A	-12	HIS	-	expression tag	UNP Q9X074
A	-11	HIS	-	expression tag	UNP Q9X074
A	-10	HIS	-	expression tag	UNP Q9X074
A	-9	SER	-	expression tag	UNP Q9X074
A	-8	SER	-	expression tag	UNP Q9X074
A	-7	GLY	-	expression tag	UNP Q9X074
A	-6	LEU	-	expression tag	UNP Q9X074
A	-5	VAL	-	expression tag	UNP Q9X074
A	-4	PRO	-	expression tag	UNP Q9X074
A	-3	ARG	-	expression tag	UNP Q9X074
A	-2	GLY	-	expression tag	UNP Q9X074
A	-1	SER	-	expression tag	UNP Q9X074
A	0	HIS	-	expression tag	UNP Q9X074
B	181	MET	-	expression tag	UNP Q9X074
B	182	GLY	-	expression tag	UNP Q9X074
B	183	SER	-	expression tag	UNP Q9X074
B	184	SER	-	expression tag	UNP Q9X074
B	185	HIS	-	expression tag	UNP Q9X074
B	186	HIS	-	expression tag	UNP Q9X074
B	187	HIS	-	expression tag	UNP Q9X074
B	188	HIS	-	expression tag	UNP Q9X074
B	189	HIS	-	expression tag	UNP Q9X074

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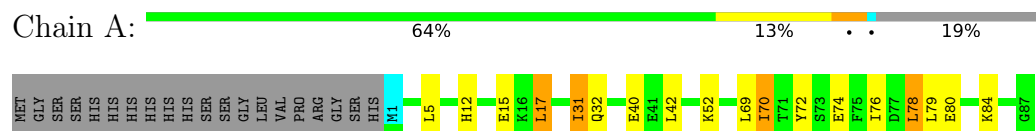
Chain	Residue	Modelled	Actual	Comment	Reference
B	190	HIS	-	expression tag	UNP Q9X074
B	191	SER	-	expression tag	UNP Q9X074
B	192	SER	-	expression tag	UNP Q9X074
B	193	GLY	-	expression tag	UNP Q9X074
B	194	LEU	-	expression tag	UNP Q9X074
B	195	VAL	-	expression tag	UNP Q9X074
B	196	PRO	-	expression tag	UNP Q9X074
B	197	ARG	-	expression tag	UNP Q9X074
B	198	GLY	-	expression tag	UNP Q9X074
B	199	SER	-	expression tag	UNP Q9X074
B	200	HIS	-	expression tag	UNP Q9X074

4 Residue-property plots [i](#)

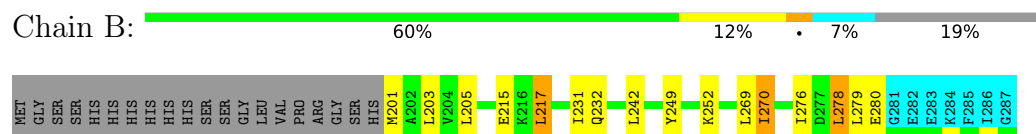
4.1 Average score per residue in the NMR ensemble

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: hypothetical protein TM0979



- Molecule 1: hypothetical protein TM0979

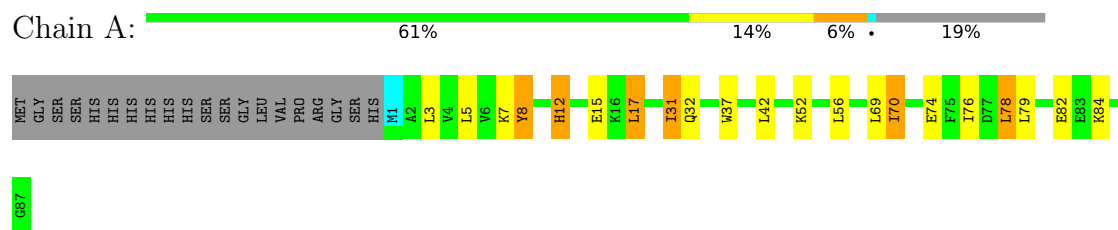


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: hypothetical protein TM0979



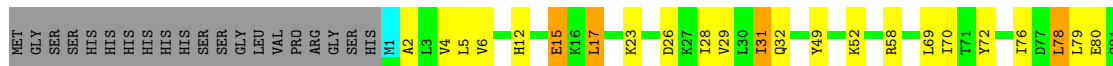
- Molecule 1: hypothetical protein TM0979





4.2.2 Score per residue for model 2

- Molecule 1: hypothetical protein TM0979

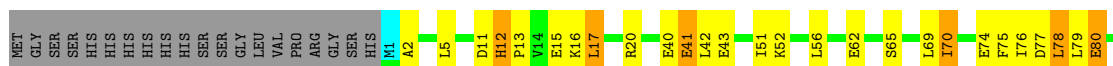


- Molecule 1: hypothetical protein TM0979

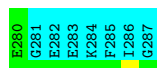


4.2.3 Score per residue for model 3 (medoid)

- Molecule 1: hypothetical protein TM0979

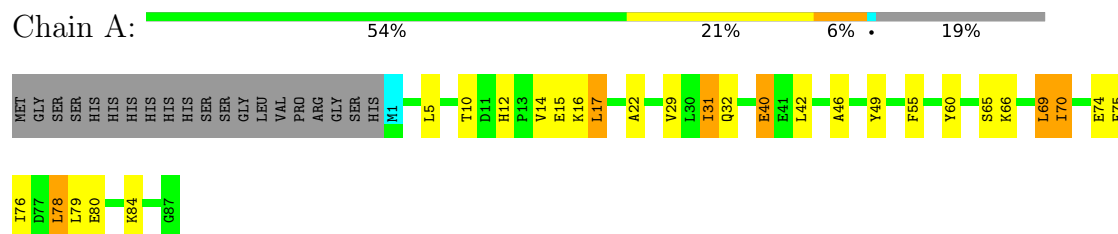


- Molecule 1: hypothetical protein TM0979

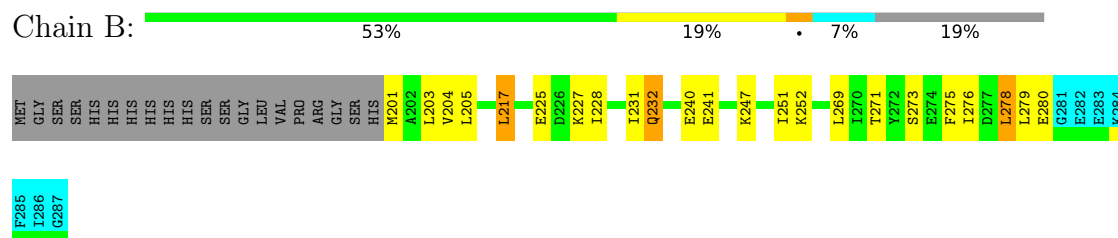


4.2.4 Score per residue for model 4

- Molecule 1: hypothetical protein TM0979

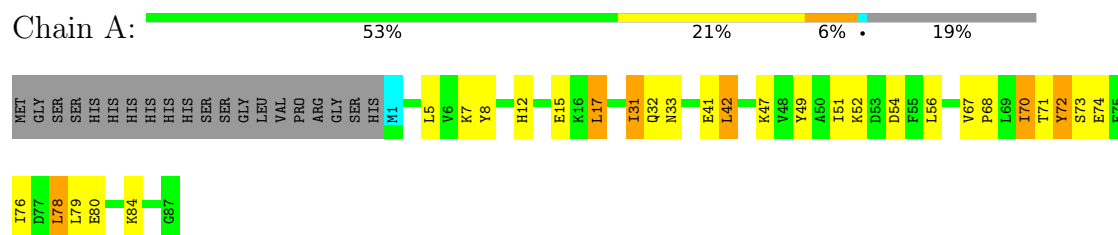


- Molecule 1: hypothetical protein TM0979

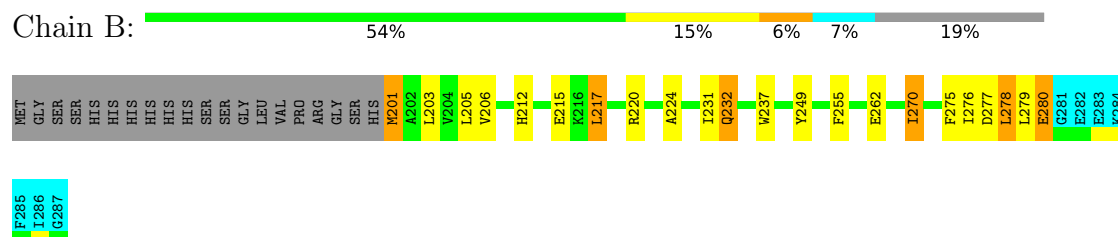


4.2.5 Score per residue for model 5

- Molecule 1: hypothetical protein TM0979

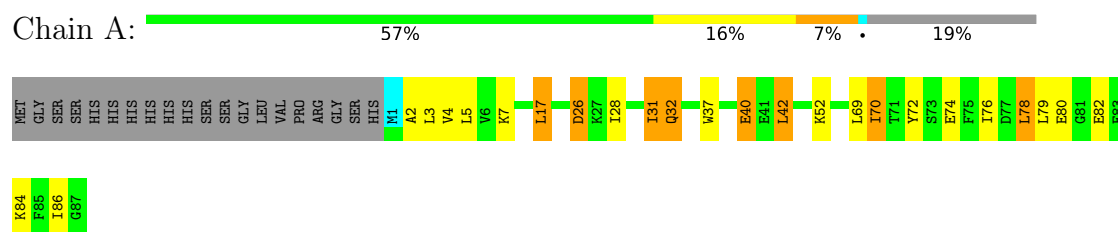


- Molecule 1: hypothetical protein TM0979

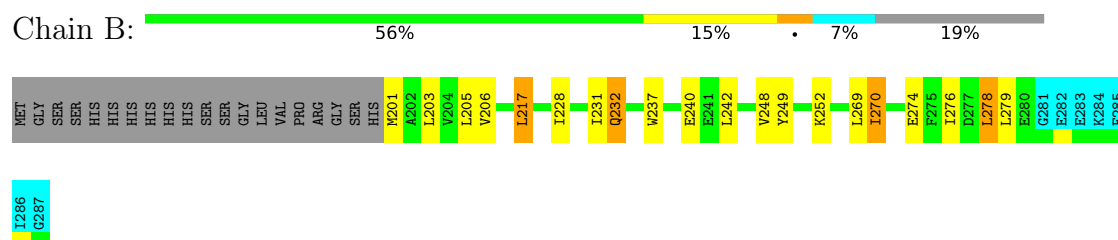


4.2.6 Score per residue for model 6

- Molecule 1: hypothetical protein TM0979

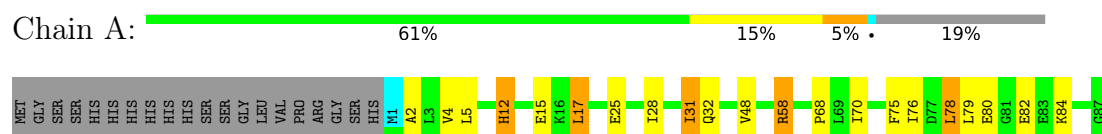


- Molecule 1: hypothetical protein TM0979

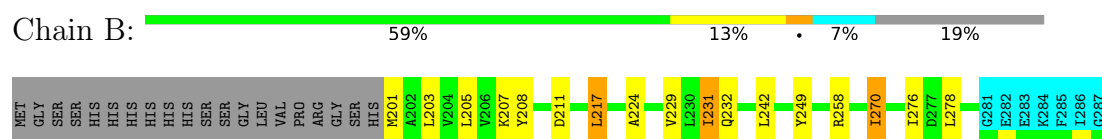


4.2.7 Score per residue for model 7

- Molecule 1: hypothetical protein TM0979

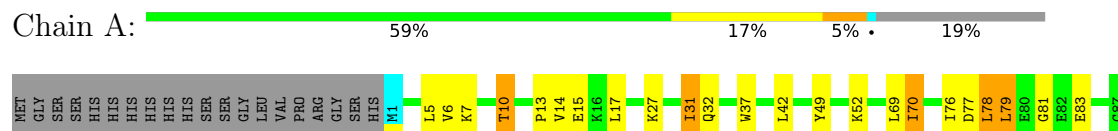


- Molecule 1: hypothetical protein TM0979



4.2.8 Score per residue for model 8

- Molecule 1: hypothetical protein TM0979



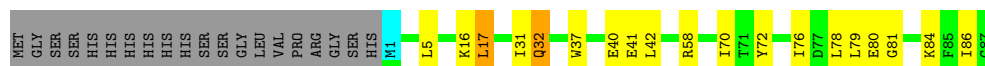
- Molecule 1: hypothetical protein TM0979



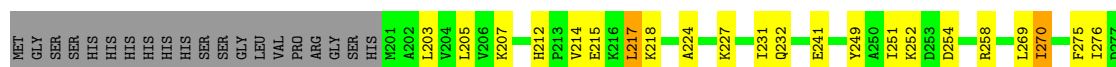


4.2.9 Score per residue for model 9

- Molecule 1: hypothetical protein TM0979

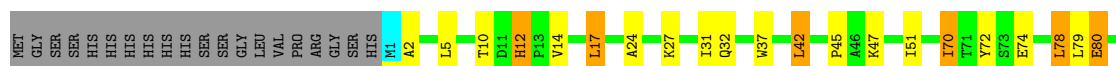


- Molecule 1: hypothetical protein TM0979



4.2.10 Score per residue for model 10

- Molecule 1: hypothetical protein TM0979



- Molecule 1: hypothetical protein TM0979



5 Refinement protocol and experimental data overview

The models were refined using the following method: *distance geometry, simulated annealing, molecular dynamics*.

Of the 20 calculated structures, 10 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
DYANA	structure solution	
CNS	refinement	

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	0.39±0.01	0±0/703 (0.0± 0.0%)	0.73±0.02	0±0/948 (0.0± 0.0%)
1	B	0.40±0.02	0±0/656 (0.0± 0.0%)	0.73±0.02	0±0/889 (0.0± 0.1%)
All	All	0.39	0/13590 (0.0%)	0.73	6/18370 (0.0%)

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	B	232	GLN	CB-CA-C	-5.91	109.77	116.63	4	4
1	A	32	GLN	CB-CA-C	-5.37	110.41	116.63	9	2

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	690	699	698	17±2
1	B	643	657	655	16±3
All	All	13330	13560	13530	268

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 10.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:16:LYS:HE3	1:A:42:LEU:HB3	0.89	1.42	9	2
1:A:76:ILE:HG23	1:B:205:LEU:HD13	0.80	1.48	4	8
1:A:79:LEU:HB3	1:B:203:LEU:HD11	0.80	1.54	9	3
1:A:79:LEU:HD21	1:B:279:LEU:HD13	0.80	1.52	8	3
1:B:212:HIS:HB3	1:B:214:VAL:HG12	0.71	1.61	8	3
1:A:86:ILE:HG22	1:B:218:LYS:HE2	0.69	1.62	9	1
1:A:12:HIS:HB3	1:A:15:GLU:HG2	0.68	1.63	7	4
1:A:78:LEU:HG	1:A:79:LEU:N	0.68	2.03	1	9
1:A:80:GLU:HB3	1:A:84:LYS:HG2	0.65	1.66	3	1
1:B:212:HIS:HD2	1:B:214:VAL:HG12	0.65	1.52	3	1
1:A:70:ILE:HB	1:A:74:GLU:HB2	0.64	1.69	10	6
1:A:8:TYR:HA	1:A:33:ASN:HB2	0.63	1.70	5	1
1:B:278:LEU:HG	1:B:279:LEU:N	0.63	2.09	3	2
1:B:270:ILE:HB	1:B:274:GLU:HB2	0.62	1.70	6	5
1:B:217:LEU:H	1:B:217:LEU:HD13	0.61	1.56	6	9
1:A:52:LYS:HD2	1:A:62:GLU:HG2	0.61	1.73	3	1
1:A:31:ILE:HG21	1:B:276:ILE:HG21	0.60	1.72	7	4
1:B:231:ILE:HD13	1:B:251:ILE:HG13	0.60	1.74	9	2
1:A:12:HIS:HB3	1:A:15:GLU:OE1	0.60	1.95	5	1
1:A:5:LEU:HD13	1:B:276:ILE:HG23	0.59	1.72	3	6
1:B:201:MET:HA	1:B:225:GLU:O	0.59	1.98	2	1
1:A:17:LEU:HD13	1:A:17:LEU:H	0.59	1.58	7	1
1:A:17:LEU:H	1:A:17:LEU:HD13	0.59	1.58	4	7
1:A:10:THR:HA	1:A:37:TRP:CZ2	0.59	2.32	8	2
1:B:278:LEU:C	1:B:278:LEU:HD22	0.58	2.24	4	6
1:B:252:LYS:HG2	1:B:269:LEU:HB3	0.57	1.73	2	1
1:A:12:HIS:HD2	1:A:14:VAL:HG12	0.57	1.59	10	1
1:B:208:TYR:HB3	1:B:211:ASP:HB3	0.57	1.77	7	1
1:B:212:HIS:HB2	1:B:215:GLU:HG2	0.56	1.76	8	1
1:A:79:LEU:HD13	1:B:279:LEU:CD1	0.56	2.30	2	1
1:A:79:LEU:HD11	1:B:279:LEU:HD11	0.56	1.76	6	2
1:A:27:LYS:HD3	1:A:47:LYS:HB2	0.56	1.78	10	1
1:B:255:PHE:HE2	1:B:262:GLU:HA	0.55	1.62	5	1
1:B:254:ASP:O	1:B:258:ARG:HG2	0.55	2.01	3	2
1:B:278:LEU:HD22	1:B:278:LEU:O	0.55	2.02	1	1
1:A:31:ILE:HG21	1:B:276:ILE:HG12	0.55	1.79	9	4
1:A:72:TYR:HD2	1:B:231:ILE:HD11	0.54	1.62	6	2
1:A:29:VAL:HA	1:A:49:TYR:O	0.54	2.02	2	2
1:B:242:LEU:HD12	1:B:267:VAL:HG11	0.54	1.78	2	1
1:A:4:VAL:O	1:A:28:ILE:HA	0.54	2.02	2	3
1:B:220:ARG:HA	1:B:243:GLU:O	0.53	2.03	1	2
1:B:252:LYS:HB2	1:B:269:LEU:HB3	0.53	1.79	9	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:86:ILE:HD13	1:B:205:LEU:HB2	0.53	1.79	10	1
1:A:31:ILE:HD13	1:A:51:ILE:HG13	0.53	1.79	10	1
1:B:227:LYS:HG2	1:B:247:LYS:HB2	0.52	1.82	4	1
1:A:14:VAL:O	1:A:17:LEU:HG	0.52	2.04	8	1
1:B:228:ILE:HB	1:B:248:VAL:HG12	0.52	1.81	6	1
1:B:210:THR:HA	1:B:237:TRP:CZ2	0.51	2.41	8	1
1:B:201:MET:HG3	1:B:225:GLU:O	0.51	2.06	4	1
1:A:5:LEU:HD11	1:B:279:LEU:HD23	0.50	1.84	2	1
1:A:76:ILE:HG23	1:B:205:LEU:HD22	0.50	1.82	1	3
1:B:212:HIS:O	1:B:215:GLU:HG3	0.50	2.06	5	1
1:A:31:ILE:HG13	1:A:32:GLN:HG3	0.49	1.84	2	9
1:A:7:LYS:HA	1:A:31:ILE:O	0.49	2.06	5	2
1:B:249:TYR:HB3	1:B:270:ILE:HD12	0.49	1.84	5	7
1:A:37:TRP:O	1:A:40:GLU:HG3	0.49	2.07	6	1
1:B:207:LYS:HE3	1:B:232:GLN:OE1	0.49	2.07	3	1
1:A:80:GLU:CD	1:A:86:ILE:HG12	0.49	2.33	3	1
1:B:232:GLN:HA	1:B:254:ASP:OD2	0.49	2.07	8	2
1:A:40:GLU:O	1:A:66:LYS:HD3	0.49	2.07	4	1
1:A:80:GLU:CD	1:A:84:LYS:HA	0.49	2.32	7	1
1:A:6:VAL:HG13	1:A:15:GLU:HG3	0.48	1.85	2	1
1:A:40:GLU:O	1:A:41:GLU:HB2	0.48	2.09	3	1
1:A:2:ALA:HA	1:A:26:ASP:HA	0.48	1.85	6	1
1:A:72:TYR:HD1	1:B:231:ILE:HD11	0.48	1.68	2	1
1:B:212:HIS:HB2	1:B:215:GLU:HG3	0.48	1.84	9	1
1:A:80:GLU:OE2	1:A:84:LYS:HB2	0.48	2.09	2	1
1:A:72:TYR:CD2	1:B:231:ILE:HD11	0.48	2.44	9	2
1:B:230:LEU:HG	1:B:237:TRP:HB2	0.48	1.85	2	1
1:A:51:ILE:HB	1:A:54:ASP:HB3	0.48	1.86	5	1
1:A:72:TYR:CD1	1:B:231:ILE:HD11	0.47	2.44	2	1
1:A:48:VAL:O	1:A:68:PRO:HD2	0.47	2.09	7	1
1:A:79:LEU:HD22	1:B:275:PHE:CE2	0.47	2.44	9	1
1:A:5:LEU:CD1	1:B:276:ILE:HG23	0.47	2.38	2	1
1:B:231:ILE:HG13	1:B:232:GLN:HG3	0.47	1.86	3	7
1:A:52:LYS:O	1:A:56:LEU:HD23	0.46	2.10	1	3
1:A:6:VAL:HG13	1:A:15:GLU:HB3	0.46	1.87	8	1
1:A:37:TRP:O	1:A:42:LEU:HD22	0.46	2.10	1	1
1:B:201:MET:HB3	1:B:225:GLU:HB3	0.46	1.88	3	1
1:A:51:ILE:CD1	1:B:276:ILE:HD11	0.46	2.40	3	1
1:A:42:LEU:HD12	1:A:42:LEU:O	0.46	2.10	4	4
1:A:72:TYR:HD1	1:A:73:SER:N	0.46	2.07	5	1
1:A:3:LEU:HD21	1:B:279:LEU:HG	0.46	1.88	6	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:52:LYS:HB2	1:A:69:LEU:HB3	0.46	1.87	2	3
1:A:81:GLY:C	1:A:83:GLU:H	0.46	2.19	8	1
1:A:13:PRO:HA	1:A:37:TRP:CH2	0.46	2.46	8	1
1:A:5:LEU:HD12	1:B:280:GLU:HG2	0.45	1.88	5	1
1:B:252:LYS:CG	1:B:269:LEU:HB3	0.45	2.41	6	1
1:B:206:VAL:HG11	1:B:237:TRP:CD1	0.45	2.46	6	3
1:B:254:ASP:O	1:B:258:ARG:HD3	0.45	2.12	9	1
1:B:207:LYS:HB3	1:B:231:ILE:HG23	0.45	1.87	7	3
1:A:75:PHE:HA	1:A:78:LEU:HD23	0.45	1.89	7	1
1:A:80:GLU:OE2	1:A:86:ILE:HG12	0.45	2.11	3	1
1:A:58:ARG:HD3	1:A:58:ARG:N	0.45	2.28	7	1
1:A:7:LYS:HB3	1:A:31:ILE:HG23	0.44	1.88	1	1
1:A:79:LEU:HD11	1:B:279:LEU:HD22	0.44	1.89	3	1
1:A:86:ILE:HG21	1:B:205:LEU:HB2	0.44	1.89	6	1
1:B:270:ILE:HB	1:B:274:GLU:CB	0.44	2.43	3	1
1:B:275:PHE:HA	1:B:278:LEU:CD1	0.44	2.43	5	1
1:A:76:ILE:HG21	1:B:231:ILE:HG21	0.44	1.90	7	1
1:A:5:LEU:HD12	1:B:280:GLU:HG3	0.43	1.88	2	1
1:A:72:TYR:O	1:A:76:ILE:HD13	0.43	2.13	5	2
1:B:217:LEU:HA	1:B:220:ARG:HD3	0.43	1.90	5	1
1:B:275:PHE:HA	1:B:278:LEU:HD12	0.43	1.89	5	2
1:A:79:LEU:HD22	1:B:279:LEU:HD22	0.43	1.89	5	1
1:B:229:VAL:HA	1:B:249:TYR:O	0.43	2.13	3	2
1:A:65:SER:CB	1:A:69:LEU:HD21	0.43	2.44	4	1
1:B:275:PHE:O	1:B:278:LEU:HD13	0.43	2.14	9	2
1:B:215:GLU:CD	1:B:215:GLU:H	0.43	2.21	1	1
1:A:70:ILE:HB	1:A:74:GLU:CB	0.43	2.44	4	2
1:A:67:VAL:HB	1:A:68:PRO:HD2	0.43	1.91	5	1
1:A:79:LEU:HD22	1:B:279:LEU:HD13	0.42	1.90	1	1
1:A:49:TYR:HB3	1:A:70:ILE:HD12	0.42	1.91	8	1
1:A:16:LYS:HB2	1:A:37:TRP:HZ3	0.42	1.74	9	1
1:A:22:ALA:HB1	1:A:46:ALA:HB2	0.42	1.91	4	1
1:B:239:LEU:HD22	1:B:264:ASP:OD2	0.42	2.13	10	1
1:A:79:LEU:HD13	1:B:279:LEU:HD13	0.42	1.91	2	1
1:A:2:ALA:HB1	1:A:26:ASP:HA	0.42	1.91	2	1
1:B:204:VAL:O	1:B:228:ILE:HA	0.42	2.14	4	1
1:A:52:LYS:CG	1:A:69:LEU:HB3	0.42	2.45	8	1
1:B:278:LEU:HD22	1:B:278:LEU:C	0.42	2.40	1	1
1:A:75:PHE:CE2	1:B:279:LEU:HD21	0.42	2.49	4	1
1:B:215:GLU:HA	1:B:218:LYS:CE	0.42	2.45	10	1
1:A:8:TYR:HB3	1:A:12:HIS:HB2	0.42	1.92	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:256:LEU:HD21	1:B:262:GLU:HG3	0.42	1.91	10	1
1:B:203:LEU:HD21	1:B:205:LEU:HD11	0.42	1.92	5	1
1:B:261:SER:HB2	1:B:264:ASP:OD1	0.42	2.15	8	1
1:A:80:GLU:OE1	1:A:84:LYS:HB3	0.42	2.15	10	1
1:B:201:MET:O	1:B:201:MET:HG2	0.41	2.14	5	1
1:B:226:ASP:O	1:B:246:ALA:HB1	0.41	2.14	10	1
1:A:23:LYS:HA	1:A:23:LYS:HE2	0.41	1.92	2	1
1:A:17:LEU:C	1:A:17:LEU:HD22	0.41	2.40	10	1
1:B:275:PHE:O	1:B:279:LEU:HG	0.41	2.15	5	1
1:B:275:PHE:HA	1:B:278:LEU:HD23	0.41	1.92	10	1
1:A:65:SER:HB3	1:A:69:LEU:HD21	0.41	1.91	3	1
1:A:55:PHE:O	1:A:60:TYR:HB2	0.41	2.15	4	1
1:A:41:GLU:C	1:A:42:LEU:HG	0.41	2.40	5	1
1:A:5:LEU:HD22	1:B:276:ILE:HG23	0.41	1.91	9	2
1:A:79:LEU:C	1:A:81:GLY:H	0.41	2.23	9	1
1:A:14:VAL:HA	1:A:17:LEU:CD1	0.41	2.45	10	1
1:A:17:LEU:HA	1:A:20:ARG:HD3	0.41	1.91	3	1
1:A:24:ALA:HA	1:A:45:PRO:O	0.41	2.14	10	1
1:B:216:LYS:O	1:B:220:ARG:HG3	0.41	2.16	2	1
1:A:11:ASP:O	1:A:13:PRO:HD3	0.41	2.16	3	1
1:A:5:LEU:HD21	1:B:279:LEU:HD23	0.40	1.93	1	1
1:A:75:PHE:O	1:A:79:LEU:HD13	0.40	2.16	3	1
1:B:252:LYS:HD2	1:B:269:LEU:HB3	0.40	1.94	4	1
1:B:214:VAL:HA	1:B:217:LEU:HD11	0.40	1.93	9	1
1:A:14:VAL:HA	1:A:17:LEU:HD11	0.40	1.93	4	1
1:A:47:LYS:HB3	1:A:49:TYR:CE2	0.40	2.52	5	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	85/107 (79%)	76±2 (89±2%)	7±2 (9±2%)	2±1 (2±1%)	7	47
1	B	79/107 (74%)	71±2 (90±2%)	7±2 (8±2%)	1±1 (2±1%)	11	57
All	All	1640/2140 (77%)	1469 (90%)	140 (9%)	31 (2%)	8	51

All 9 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	84	LYS	7
1	B	242	LEU	6
1	A	82	GLU	4
1	B	224	ALA	4
1	A	2	ALA	3
1	A	42	LEU	3
1	A	41	GLU	2
1	B	240	GLU	1
1	B	241	GLU	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	74/92 (80%)	67±1 (90±1%)	7±1 (10±1%)	10	54
1	B	70/92 (76%)	65±1 (92±1%)	5±1 (8±1%)	14	61
All	All	1440/1840 (78%)	1313 (91%)	127 (9%)	11	58

All 41 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	70	ILE	10
1	A	78	LEU	10
1	B	217	LEU	10
1	B	278	LEU	10
1	A	17	LEU	9
1	B	270	ILE	9
1	A	31	ILE	7
1	A	80	GLU	6
1	A	12	HIS	5
1	B	231	ILE	4
1	B	280	GLU	3
1	A	58	ARG	3
1	A	42	LEU	3

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Mol	Chain	Res	Type	Models (Total)
1	A	40	GLU	3
1	B	203	LEU	3
1	B	201	MET	3
1	B	215	GLU	2
1	A	77	ASP	2
1	A	10	THR	2
1	B	271	THR	2
1	A	3	LEU	1
1	A	8	TYR	1
1	A	15	GLU	1
1	A	16	LYS	1
1	A	43	GLU	1
1	B	223	LYS	1
1	A	69	LEU	1
1	B	273	SER	1
1	A	71	THR	1
1	A	72	TYR	1
1	B	277	ASP	1
1	A	26	ASP	1
1	B	240	GLU	1
1	A	25	GLU	1
1	B	258	ARG	1
1	A	7	LYS	1
1	A	27	LYS	1
1	A	79	LEU	1
1	B	210	THR	1
1	B	227	LYS	1
1	B	241	GLU	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