



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 12:37 AM UTC

PDB ID : 1ZDV / pdb_00001zdv
Title : Solution Structure of the type 1 pilus assembly platform FimD(25-139)
Authors : Nishiyama, M.; Horst, R.; Herrmann, T.; Vetsch, M.; Bettendorff, P.; Ignatov, O.; Grutter, M.; Wuthrich, K.; Glockshuber, R.; Capitani, G.
Deposited on : 2005-04-15

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

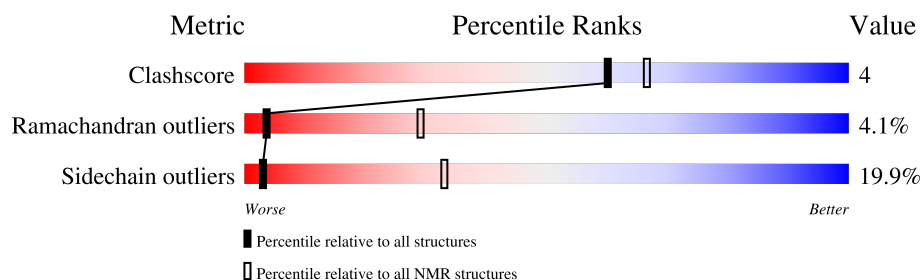
1 Overall quality at a glance

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	115	

2 Ensemble composition and analysis

This entry contains 20 models. Model 3 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *closest to the average*.

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:27-A:137 (111)	0.65	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	1, 4, 6, 8, 9, 10, 12, 13, 15, 17, 18, 20
2	2, 3, 7, 14, 16, 19
3	5, 11

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 1721 atoms, of which 856 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Outer membrane usher protein fimD.

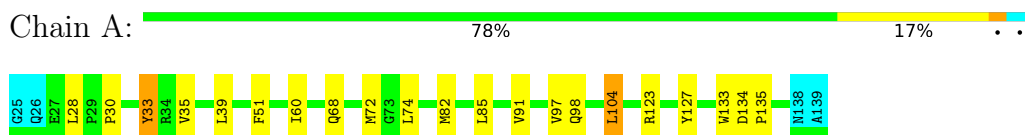
Mol	Chain	Residues	Atoms						Trace
1	A	115	Total	C	H	N	O	S	0
			1721	537	856	151	170	7	

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: Outer membrane usher protein fimD

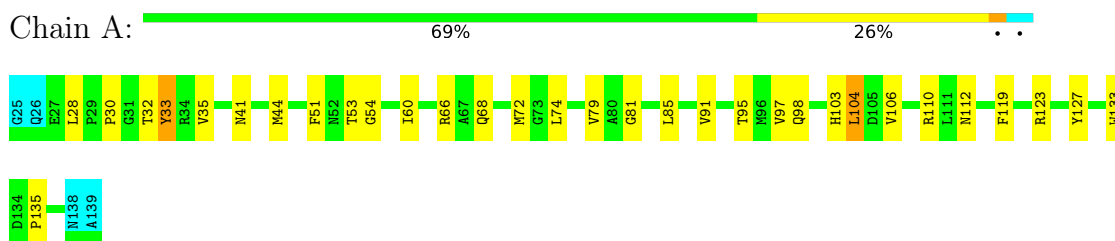


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

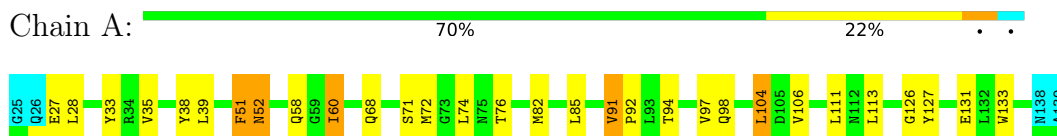
4.2.1 Score per residue for model 1

- Molecule 1: Outer membrane usher protein fimD



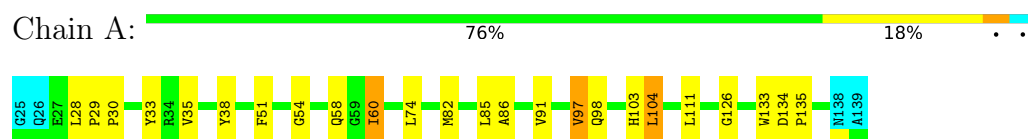
4.2.2 Score per residue for model 2

- Molecule 1: Outer membrane usher protein fimD



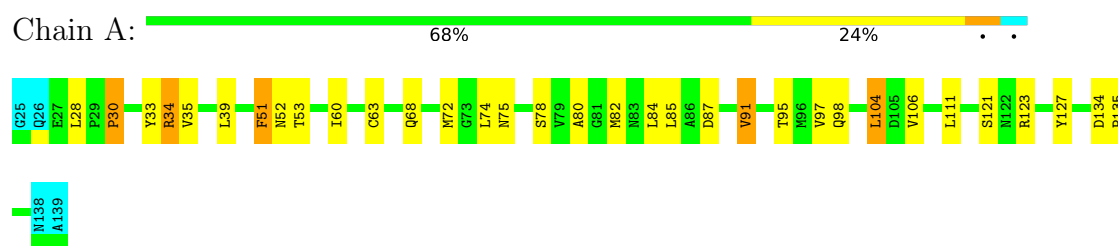
4.2.3 Score per residue for model 3 (medoid)

- Molecule 1: Outer membrane usher protein fimD



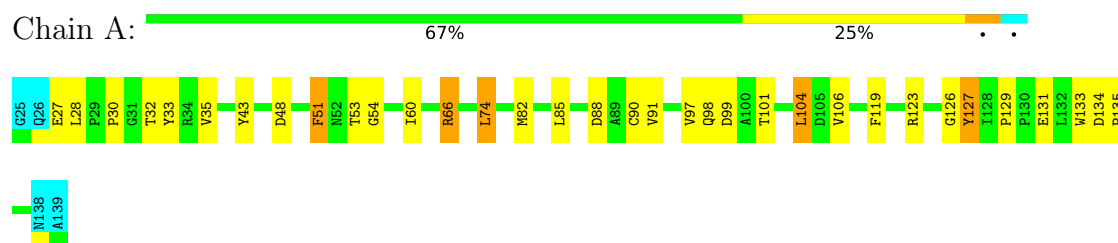
4.2.4 Score per residue for model 4

- Molecule 1: Outer membrane usher protein fimD



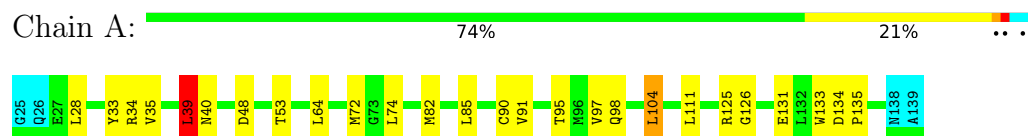
4.2.5 Score per residue for model 5

- Molecule 1: Outer membrane usher protein fimD



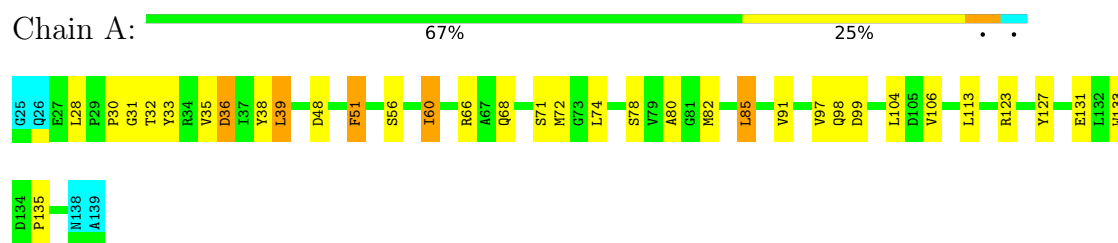
4.2.6 Score per residue for model 6

- Molecule 1: Outer membrane usher protein fimD



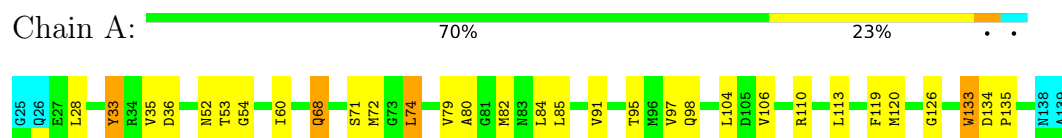
4.2.7 Score per residue for model 7

- Molecule 1: Outer membrane usher protein fimD



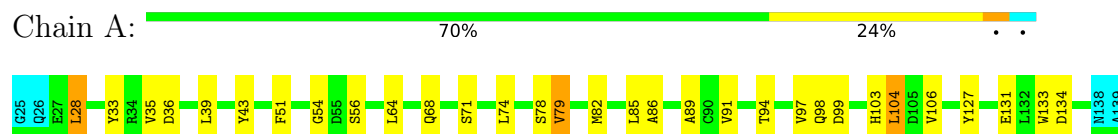
4.2.8 Score per residue for model 8

- Molecule 1: Outer membrane usher protein fimD



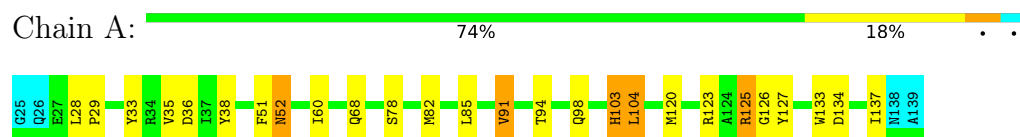
4.2.9 Score per residue for model 9

- Molecule 1: Outer membrane usher protein fimD



4.2.10 Score per residue for model 10

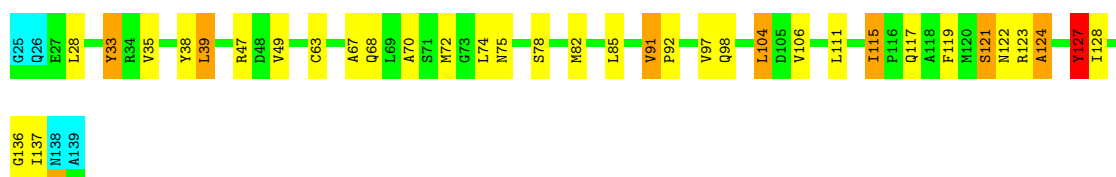
- Molecule 1: Outer membrane usher protein fimD



4.2.11 Score per residue for model 11

- Molecule 1: Outer membrane usher protein fimD

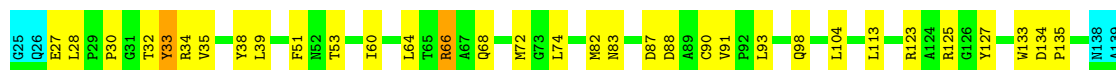




4.2.12 Score per residue for model 12

- Molecule 1: Outer membrane usher protein fimD

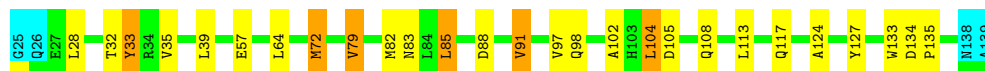
Chain A: 68% 27% . .



4.2.13 Score per residue for model 13

- Molecule 1: Outer membrane usher protein fimD

Chain A: 73% 18% 5% . .



4.2.14 Score per residue for model 14

- Molecule 1: Outer membrane usher protein fimD

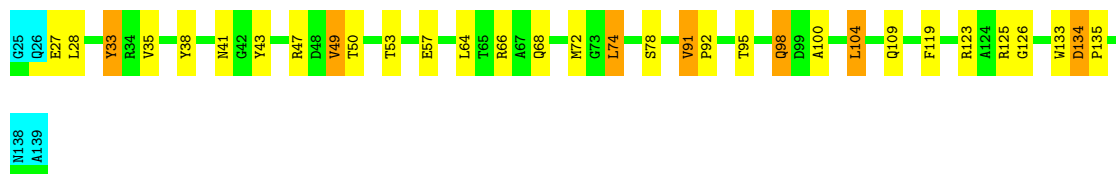
Chain A: 67% 27% . .



4.2.15 Score per residue for model 15

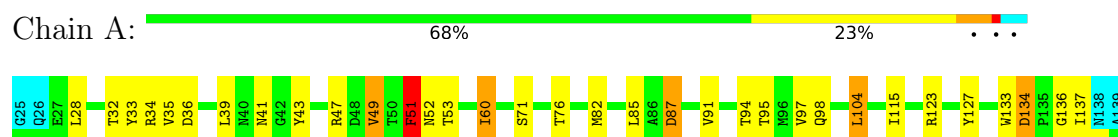
- Molecule 1: Outer membrane usher protein fimD

Chain A: 69% 22% 6% . .



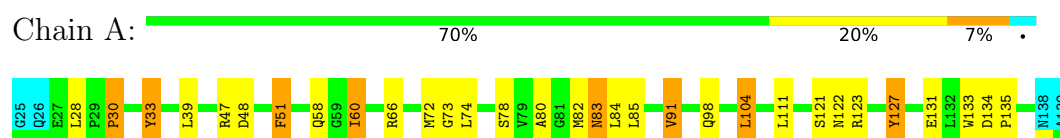
4.2.16 Score per residue for model 16

- Molecule 1: Outer membrane usher protein fimD



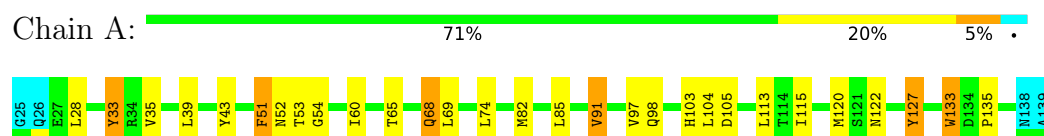
4.2.17 Score per residue for model 17

- Molecule 1: Outer membrane usher protein fimD



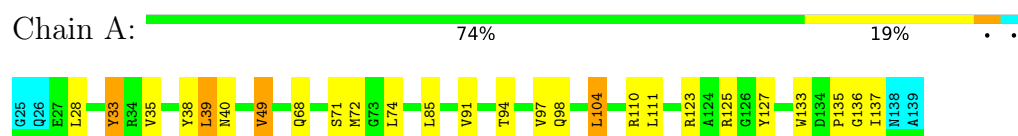
4.2.18 Score per residue for model 18

- Molecule 1: Outer membrane usher protein fimD



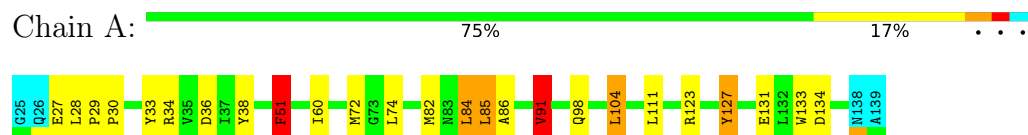
4.2.19 Score per residue for model 19

- Molecule 1: Outer membrane usher protein fimD



4.2.20 Score per residue for model 20

- Molecule 1: Outer membrane usher protein fimD



5 Refinement protocol and experimental data overview ⓘ

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

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *target function*.

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

Software name	Classification	Version
ATHNOS/CANDID	structure solution	
ATHNOS/CANDID	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	0.66±0.02	0±0/856 (0.0± 0.0%)	1.44±0.04	3±2/1172 (0.2± 0.2%)
All	All	0.66	0/17120 (0.0%)	1.44	54/23440 (0.2%)

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	1.4±1.0
All	All	0	27

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	ASP	CA-CB-CG	8.86	121.46	112.60	8	2
1	A	115	ILE	CA-C-N	7.15	128.78	119.84	11	1
1	A	115	ILE	C-N-CA	7.15	128.78	119.84	11	1
1	A	86	ALA	CA-C-N	6.88	129.80	120.38	20	3
1	A	86	ALA	C-N-CA	6.88	129.80	120.38	20	3
1	A	51	PHE	CA-CB-CG	6.83	120.63	113.80	20	1
1	A	52	ASN	CA-CB-CG	6.82	119.42	112.60	2	1
1	A	92	PRO	N-CA-CB	6.60	106.62	102.92	2	2
1	A	29	PRO	N-CA-CB	6.46	106.58	103.22	20	2
1	A	134	ASP	CA-CB-CG	6.44	119.04	112.60	3	5
1	A	47	ARG	CA-C-N	6.18	129.49	120.71	16	1
1	A	47	ARG	C-N-CA	6.18	129.49	120.71	16	1
1	A	87	ASP	CA-CB-CG	6.18	118.78	112.60	16	2
1	A	83	ASN	CA-CB-CG	-6.12	106.48	112.60	12	2
1	A	84	LEU	CB-CA-C	5.93	119.35	109.80	20	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	41	ASN	CA-C-N	5.92	126.27	122.18	14	1
1	A	41	ASN	C-N-CA	5.92	126.27	122.18	14	1
1	A	79	VAL	CB-CA-C	5.90	117.61	111.23	9	1
1	A	103	HIS	CA-CB-CG	5.86	119.66	113.80	10	1
1	A	91	VAL	CB-CA-C	5.83	116.71	110.53	20	1
1	A	121	SER	CA-C-N	5.76	128.26	120.65	17	2
1	A	121	SER	C-N-CA	5.76	128.26	120.65	17	2
1	A	66	ARG	CA-CB-CG	5.72	125.54	114.10	5	1
1	A	103	HIS	CB-CG-CD2	-5.69	123.80	131.20	18	2
1	A	66	ARG	CB-CA-C	-5.64	102.02	110.88	5	1
1	A	39	LEU	CA-C-N	5.58	130.05	122.07	6	1
1	A	39	LEU	C-N-CA	5.58	130.05	122.07	6	1
1	A	126	GLY	N-CA-C	-5.57	105.94	111.85	10	1
1	A	122	ASN	CA-C-N	5.53	132.09	121.54	17	1
1	A	122	ASN	C-N-CA	5.53	132.09	121.54	17	1
1	A	85	LEU	CB-CA-C	5.42	120.00	110.36	16	1
1	A	94	THR	CB-CA-C	5.28	119.41	111.18	9	1
1	A	28	LEU	N-CA-C	5.21	117.06	109.30	9	1
1	A	97	VAL	CB-CA-C	5.12	119.69	111.29	3	1
1	A	75	ASN	CA-CB-CG	5.07	117.67	112.60	4	1
1	A	129	PRO	N-CA-CB	5.04	105.84	103.22	5	1
1	A	29	PRO	CA-C-N	5.02	126.12	119.84	14	1
1	A	29	PRO	C-N-CA	5.02	126.12	119.84	14	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	34	ARG	Sidechain	5
1	A	66	ARG	Sidechain	4
1	A	123	ARG	Sidechain	3
1	A	110	ARG	Sidechain	2
1	A	51	PHE	Sidechain	2
1	A	43	TYR	Sidechain	2
1	A	125	ARG	Peptide,Sidechain	2
1	A	124	ALA	Peptide	2
1	A	127	TYR	Sidechain	2
1	A	47	ARG	Sidechain	2
1	A	33	TYR	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	839	834	822	7±2
All	All	16780	16680	16440	149

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 4.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:76:THR:HG22	1:A:82:MET:SD	0.73	2.23	16	1
1:A:33:TYR:CE2	1:A:35:VAL:HG11	0.71	2.21	16	18
1:A:82:MET:HE2	1:A:91:VAL:HG11	0.66	1.66	8	12
1:A:33:TYR:CE2	1:A:104:LEU:HD23	0.64	2.27	20	3
1:A:33:TYR:CE2	1:A:104:LEU:HD13	0.62	2.29	4	8
1:A:74:LEU:HD22	1:A:119:PHE:O	0.58	1.97	11	4
1:A:33:TYR:CZ	1:A:104:LEU:HD23	0.54	2.38	19	1
1:A:82:MET:HE2	1:A:91:VAL:HG21	0.54	1.80	7	4
1:A:39:LEU:HD13	1:A:72:MET:HB3	0.53	1.79	13	1
1:A:82:MET:HA	1:A:85:LEU:HD12	0.52	1.81	20	4
1:A:51:PHE:CE2	1:A:60:ILE:HD12	0.52	2.40	20	10
1:A:33:TYR:CZ	1:A:35:VAL:HG21	0.52	2.40	13	1
1:A:127:TYR:C	1:A:128:ILE:HD12	0.51	2.31	11	1
1:A:51:PHE:CD2	1:A:60:ILE:HD12	0.51	2.41	10	4
1:A:33:TYR:CE2	1:A:104:LEU:CD2	0.51	2.94	19	3
1:A:74:LEU:HD23	1:A:115:ILE:HD13	0.51	1.82	11	1
1:A:102:ALA:HA	1:A:113:LEU:HD23	0.50	1.83	13	1
1:A:65:THR:C	1:A:69:LEU:HD13	0.50	2.31	18	1
1:A:39:LEU:HD11	1:A:115:ILE:CD1	0.50	2.37	11	2
1:A:104:LEU:C	1:A:104:LEU:HD22	0.49	2.32	13	4
1:A:30:PRO:HD2	1:A:51:PHE:CZ	0.49	2.43	14	8
1:A:70:ALA:CB	1:A:124:ALA:HB2	0.49	2.38	11	2
1:A:33:TYR:CE2	1:A:35:VAL:CG1	0.48	2.96	14	2
1:A:127:TYR:C	1:A:127:TYR:CD1	0.48	2.91	11	1
1:A:82:MET:HE2	1:A:91:VAL:CG1	0.48	2.38	11	2
1:A:33:TYR:CD1	1:A:49:VAL:HG13	0.48	2.43	16	3
1:A:33:TYR:CD1	1:A:51:PHE:CE1	0.48	3.00	7	7

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:67:ALA:HB2	1:A:128:ILE:HD13	0.48	1.84	11	1
1:A:33:TYR:CZ	1:A:104:LEU:CD2	0.48	2.97	19	1
1:A:82:MET:HE2	1:A:91:VAL:CG2	0.47	2.38	20	2
1:A:32:THR:HG23	1:A:49:VAL:C	0.47	2.34	16	1
1:A:127:TYR:CD1	1:A:127:TYR:C	0.47	2.91	20	2
1:A:128:ILE:HD12	1:A:128:ILE:N	0.46	2.25	14	1
1:A:29:PRO:HB3	1:A:60:ILE:HG21	0.46	1.87	10	1
1:A:104:LEU:HD12	1:A:104:LEU:C	0.45	2.36	15	2
1:A:68:GLN:NE2	1:A:133:TRP:CD2	0.45	2.85	18	2
1:A:33:TYR:CE1	1:A:49:VAL:HG13	0.45	2.46	19	2
1:A:39:LEU:CD2	1:A:120:MET:HE1	0.45	2.42	18	1
1:A:33:TYR:CZ	1:A:104:LEU:HD13	0.44	2.47	4	1
1:A:74:LEU:HD13	1:A:75:ASN:C	0.44	2.37	11	1
1:A:39:LEU:HD22	1:A:40:ASN:H	0.44	1.73	6	2
1:A:33:TYR:CZ	1:A:104:LEU:HG	0.43	2.48	6	1
1:A:74:LEU:HD23	1:A:115:ILE:HG21	0.43	1.89	11	1
1:A:79:VAL:HG21	1:A:82:MET:HE3	0.43	1.90	13	2
1:A:33:TYR:CD2	1:A:35:VAL:HG11	0.43	2.49	2	2
1:A:79:VAL:HG12	1:A:119:PHE:CD2	0.42	2.49	1	1
1:A:39:LEU:HD13	1:A:72:MET:CB	0.42	2.43	13	1
1:A:98:GLN:C	1:A:100:ALA:H	0.42	2.22	15	1
1:A:39:LEU:HD11	1:A:115:ILE:HG13	0.42	1.91	18	1
1:A:86:ALA:HB3	1:A:89:ALA:HB2	0.42	1.92	9	1
1:A:52:ASN:C	1:A:60:ILE:HG22	0.42	2.40	10	1
1:A:82:MET:CE	1:A:91:VAL:HG21	0.42	2.45	3	1
1:A:74:LEU:HD13	1:A:75:ASN:N	0.42	2.29	11	1
1:A:69:LEU:HB3	1:A:74:LEU:HD12	0.41	1.91	14	1
1:A:44:MET:SD	1:A:72:MET:HA	0.41	2.55	1	1
1:A:104:LEU:C	1:A:104:LEU:CD2	0.41	2.93	6	2
1:A:39:LEU:C	1:A:39:LEU:HD13	0.40	2.42	7	1
1:A:91:VAL:HG12	1:A:92:PRO:HD2	0.40	1.93	15	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	111/115 (97%)	88±3 (80±3%)	18±3 (16±3%)	4±2 (4±2%)	4	29
All	All	2220/2300 (97%)	1769 (80%)	361 (16%)	90 (4%)	4	29

All 24 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	135	PRO	14
1	A	134	ASP	11
1	A	127	TYR	7
1	A	123	ARG	7
1	A	54	GLY	6
1	A	126	GLY	6
1	A	32	THR	5
1	A	80	ALA	4
1	A	137	ILE	4
1	A	90	CYS	3
1	A	99	ASP	3
1	A	43	TYR	3
1	A	136	GLY	3
1	A	63	CYS	2
1	A	27	GLU	2
1	A	122	ASN	2
1	A	81	GLY	1
1	A	31	GLY	1
1	A	85	LEU	1
1	A	121	SER	1
1	A	109	GLN	1
1	A	73	GLY	1
1	A	83	ASN	1
1	A	30	PRO	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	91/93 (98%)	73±3 (80±3%)	18±3 (20±3%)	3	33
All	All	1820/1860 (98%)	1458 (80%)	362 (20%)	3	33

All 59 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	28	LEU	20
1	A	98	GLN	20
1	A	104	LEU	18
1	A	133	TRP	18
1	A	74	LEU	16
1	A	85	LEU	15
1	A	97	VAL	15
1	A	68	GLN	13
1	A	72	MET	13
1	A	91	VAL	12
1	A	38	TYR	10
1	A	39	LEU	10
1	A	33	TYR	9
1	A	53	THR	9
1	A	127	TYR	9
1	A	60	ILE	8
1	A	106	VAL	8
1	A	111	LEU	8
1	A	131	GLU	7
1	A	51	PHE	7
1	A	78	SER	7
1	A	95	THR	6
1	A	52	ASN	6
1	A	71	SER	6
1	A	113	LEU	6
1	A	36	ASP	6
1	A	94	THR	5
1	A	64	LEU	5
1	A	125	ARG	5
1	A	49	VAL	5
1	A	84	LEU	4
1	A	48	ASP	4
1	A	41	ASN	3
1	A	66	ARG	3
1	A	103	HIS	3
1	A	27	GLU	3
1	A	58	GLN	3
1	A	30	PRO	3
1	A	88	ASP	3
1	A	56	SER	3
1	A	105	ASP	3

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Mol	Chain	Res	Type	Models (Total)
1	A	123	ARG	3
1	A	34	ARG	2
1	A	120	MET	2
1	A	79	VAL	2
1	A	117	GLN	2
1	A	87	ASP	2
1	A	112	ASN	1
1	A	76	THR	1
1	A	121	SER	1
1	A	101	THR	1
1	A	47	ARG	1
1	A	93	LEU	1
1	A	108	GLN	1
1	A	46	THR	1
1	A	50	THR	1
1	A	57	GLU	1
1	A	83	ASN	1
1	A	110	ARG	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](#)

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 ⓘ

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided