



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:25 PM UTC

PDB ID : 7SGN / pdb_00007sgn
Title : Crystal structure of periplasmic domain of Helicobacter pylori FliL (residues 81 to 183) (crystal form A)
Authors : Chan, K.L.; Peterson, B.; Khan, M.F.; Roujeinikova, A.
Deposited on : 2021-10-06
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray 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/XrayValidationReportHelp>

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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

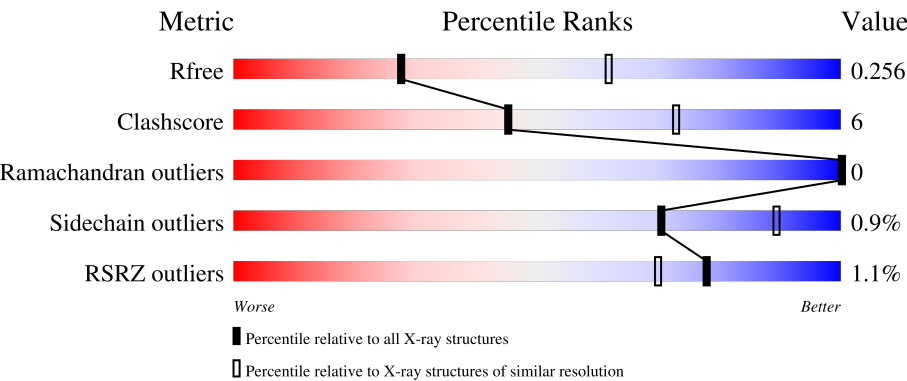
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

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)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	113	% 76%13%11%
1	B	113	2% 84%5%11%
1	C	113	% 75%19%.
1	D	113	% 78%12%11%
1	E	113	% 77%12%12%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	113	85%11%.
1	G	113	2% 80%11%.9%
1	H	113	% 79%11%11%
1	I	113	% 74%16%10%
1	J	113	2% 81%15%.
1	K	113	% 65%20%.13%
1	L	113	63%26%.11%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10096 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Flagellar protein FliL.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	101	Total	C	N	O	0	2	0
			806	519	129	158			
1	B	101	Total	C	N	O	0	2	0
			802	518	129	155			
1	C	108	Total	C	N	O	0	2	0
			854	553	135	166			
1	D	101	Total	C	N	O	0	0	0
			795	514	128	153			
1	E	100	Total	C	N	O	0	3	0
			810	520	131	159			
1	F	108	Total	C	N	O	0	1	0
			855	553	136	166			
1	G	103	Total	C	N	O	0	2	0
			823	532	133	158			
1	H	101	Total	C	N	O	0	2	0
			817	525	133	159			
1	I	102	Total	C	N	O	0	2	0
			812	525	130	157			
1	J	108	Total	C	N	O	0	1	0
			851	552	138	161			
1	K	98	Total	C	N	O	0	1	0
			777	504	124	149			
1	L	101	Total	C	N	O	0	1	0
			798	514	129	155			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	71	SER	-	expression tag	UNP A0A293T1P9
A	72	GLN	-	expression tag	UNP A0A293T1P9
A	73	ASN	-	expression tag	UNP A0A293T1P9
A	74	GLY	-	expression tag	UNP A0A293T1P9
A	75	ARG	-	expression tag	UNP A0A293T1P9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	76	ILE	-	expression tag	UNP A0A293T1P9
A	77	ASP	-	expression tag	UNP A0A293T1P9
A	78	PRO	-	expression tag	UNP A0A293T1P9
A	79	PHE	-	expression tag	UNP A0A293T1P9
A	80	THR	-	expression tag	UNP A0A293T1P9
B	71	SER	-	expression tag	UNP A0A293T1P9
B	72	GLN	-	expression tag	UNP A0A293T1P9
B	73	ASN	-	expression tag	UNP A0A293T1P9
B	74	GLY	-	expression tag	UNP A0A293T1P9
B	75	ARG	-	expression tag	UNP A0A293T1P9
B	76	ILE	-	expression tag	UNP A0A293T1P9
B	77	ASP	-	expression tag	UNP A0A293T1P9
B	78	PRO	-	expression tag	UNP A0A293T1P9
B	79	PHE	-	expression tag	UNP A0A293T1P9
B	80	THR	-	expression tag	UNP A0A293T1P9
C	71	SER	-	expression tag	UNP A0A293T1P9
C	72	GLN	-	expression tag	UNP A0A293T1P9
C	73	ASN	-	expression tag	UNP A0A293T1P9
C	74	GLY	-	expression tag	UNP A0A293T1P9
C	75	ARG	-	expression tag	UNP A0A293T1P9
C	76	ILE	-	expression tag	UNP A0A293T1P9
C	77	ASP	-	expression tag	UNP A0A293T1P9
C	78	PRO	-	expression tag	UNP A0A293T1P9
C	79	PHE	-	expression tag	UNP A0A293T1P9
C	80	THR	-	expression tag	UNP A0A293T1P9
D	71	SER	-	expression tag	UNP A0A293T1P9
D	72	GLN	-	expression tag	UNP A0A293T1P9
D	73	ASN	-	expression tag	UNP A0A293T1P9
D	74	GLY	-	expression tag	UNP A0A293T1P9
D	75	ARG	-	expression tag	UNP A0A293T1P9
D	76	ILE	-	expression tag	UNP A0A293T1P9
D	77	ASP	-	expression tag	UNP A0A293T1P9
D	78	PRO	-	expression tag	UNP A0A293T1P9
D	79	PHE	-	expression tag	UNP A0A293T1P9
D	80	THR	-	expression tag	UNP A0A293T1P9
E	71	SER	-	expression tag	UNP A0A293T1P9
E	72	GLN	-	expression tag	UNP A0A293T1P9
E	73	ASN	-	expression tag	UNP A0A293T1P9
E	74	GLY	-	expression tag	UNP A0A293T1P9
E	75	ARG	-	expression tag	UNP A0A293T1P9
E	76	ILE	-	expression tag	UNP A0A293T1P9
E	77	ASP	-	expression tag	UNP A0A293T1P9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	78	PRO	-	expression tag	UNP A0A293T1P9
E	79	PHE	-	expression tag	UNP A0A293T1P9
E	80	THR	-	expression tag	UNP A0A293T1P9
F	71	SER	-	expression tag	UNP A0A293T1P9
F	72	GLN	-	expression tag	UNP A0A293T1P9
F	73	ASN	-	expression tag	UNP A0A293T1P9
F	74	GLY	-	expression tag	UNP A0A293T1P9
F	75	ARG	-	expression tag	UNP A0A293T1P9
F	76	ILE	-	expression tag	UNP A0A293T1P9
F	77	ASP	-	expression tag	UNP A0A293T1P9
F	78	PRO	-	expression tag	UNP A0A293T1P9
F	79	PHE	-	expression tag	UNP A0A293T1P9
F	80	THR	-	expression tag	UNP A0A293T1P9
G	71	SER	-	expression tag	UNP A0A293T1P9
G	72	GLN	-	expression tag	UNP A0A293T1P9
G	73	ASN	-	expression tag	UNP A0A293T1P9
G	74	GLY	-	expression tag	UNP A0A293T1P9
G	75	ARG	-	expression tag	UNP A0A293T1P9
G	76	ILE	-	expression tag	UNP A0A293T1P9
G	77	ASP	-	expression tag	UNP A0A293T1P9
G	78	PRO	-	expression tag	UNP A0A293T1P9
G	79	PHE	-	expression tag	UNP A0A293T1P9
G	80	THR	-	expression tag	UNP A0A293T1P9
H	71	SER	-	expression tag	UNP A0A293T1P9
H	72	GLN	-	expression tag	UNP A0A293T1P9
H	73	ASN	-	expression tag	UNP A0A293T1P9
H	74	GLY	-	expression tag	UNP A0A293T1P9
H	75	ARG	-	expression tag	UNP A0A293T1P9
H	76	ILE	-	expression tag	UNP A0A293T1P9
H	77	ASP	-	expression tag	UNP A0A293T1P9
H	78	PRO	-	expression tag	UNP A0A293T1P9
H	79	PHE	-	expression tag	UNP A0A293T1P9
H	80	THR	-	expression tag	UNP A0A293T1P9
I	71	SER	-	expression tag	UNP A0A293T1P9
I	72	GLN	-	expression tag	UNP A0A293T1P9
I	73	ASN	-	expression tag	UNP A0A293T1P9
I	74	GLY	-	expression tag	UNP A0A293T1P9
I	75	ARG	-	expression tag	UNP A0A293T1P9
I	76	ILE	-	expression tag	UNP A0A293T1P9
I	77	ASP	-	expression tag	UNP A0A293T1P9
I	78	PRO	-	expression tag	UNP A0A293T1P9
I	79	PHE	-	expression tag	UNP A0A293T1P9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
I	80	THR	-	expression tag	UNP A0A293T1P9
J	71	SER	-	expression tag	UNP A0A293T1P9
J	72	GLN	-	expression tag	UNP A0A293T1P9
J	73	ASN	-	expression tag	UNP A0A293T1P9
J	74	GLY	-	expression tag	UNP A0A293T1P9
J	75	ARG	-	expression tag	UNP A0A293T1P9
J	76	ILE	-	expression tag	UNP A0A293T1P9
J	77	ASP	-	expression tag	UNP A0A293T1P9
J	78	PRO	-	expression tag	UNP A0A293T1P9
J	79	PHE	-	expression tag	UNP A0A293T1P9
J	80	THR	-	expression tag	UNP A0A293T1P9
K	71	SER	-	expression tag	UNP A0A293T1P9
K	72	GLN	-	expression tag	UNP A0A293T1P9
K	73	ASN	-	expression tag	UNP A0A293T1P9
K	74	GLY	-	expression tag	UNP A0A293T1P9
K	75	ARG	-	expression tag	UNP A0A293T1P9
K	76	ILE	-	expression tag	UNP A0A293T1P9
K	77	ASP	-	expression tag	UNP A0A293T1P9
K	78	PRO	-	expression tag	UNP A0A293T1P9
K	79	PHE	-	expression tag	UNP A0A293T1P9
K	80	THR	-	expression tag	UNP A0A293T1P9
L	71	SER	-	expression tag	UNP A0A293T1P9
L	72	GLN	-	expression tag	UNP A0A293T1P9
L	73	ASN	-	expression tag	UNP A0A293T1P9
L	74	GLY	-	expression tag	UNP A0A293T1P9
L	75	ARG	-	expression tag	UNP A0A293T1P9
L	76	ILE	-	expression tag	UNP A0A293T1P9
L	77	ASP	-	expression tag	UNP A0A293T1P9
L	78	PRO	-	expression tag	UNP A0A293T1P9
L	79	PHE	-	expression tag	UNP A0A293T1P9
L	80	THR	-	expression tag	UNP A0A293T1P9

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	38	Total O 38 38	0	0
2	B	29	Total O 29 29	0	0
2	C	39	Total O 39 39	0	0
2	D	24	Total O 24 24	0	0

Continued on next page...

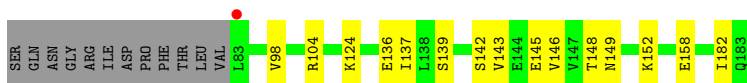
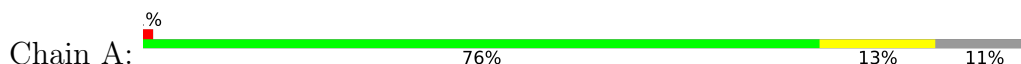
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	22	Total 22	O 22	0	0
2	F	35	Total 35	O 35	0	0
2	G	23	Total 23	O 23	0	0
2	H	17	Total 17	O 17	0	0
2	I	17	Total 17	O 17	0	0
2	J	21	Total 21	O 21	0	0
2	K	16	Total 16	O 16	0	0
2	L	15	Total 15	O 15	0	0

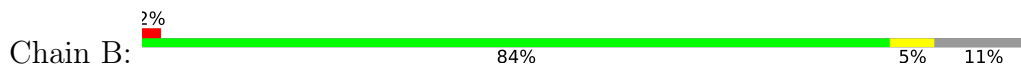
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

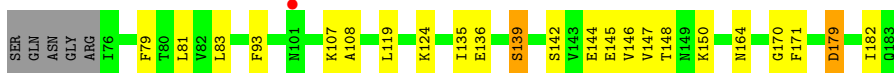
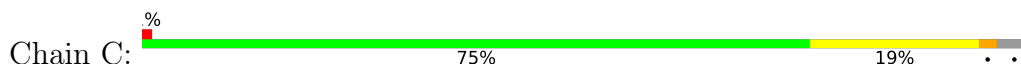
- Molecule 1: Flagellar protein FliL



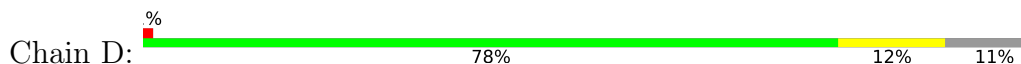
- Molecule 1: Flagellar protein FliL



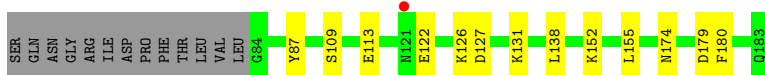
- Molecule 1: Flagellar protein FliL



- Molecule 1: Flagellar protein FliL



- Molecule 1: Flagellar protein FliL



- Molecule 1: Flagellar protein FliL

Chain F:  85% 11% .



- Molecule 1: Flagellar protein FliL

Chain G:  80% 11% 9% .



- Molecule 1: Flagellar protein FliL

Chain H:  79% 11% 11% .



- Molecule 1: Flagellar protein FliL

Chain I:  74% 16% 10% .



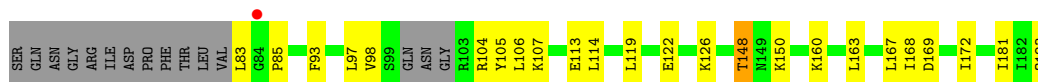
- Molecule 1: Flagellar protein FliL

Chain J:  81% 15% .



- Molecule 1: Flagellar protein FliL

Chain K:  65% 20% 13% .



- Molecule 1: Flagellar protein FliL

Chain L:  63% 26% 11% .





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.50Å 82.55Å 97.84Å 67.72° 83.39° 72.82°	Depositor
Resolution (Å)	46.90 – 2.80 46.90 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.1 (46.90-2.80) 98.1 (46.90-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.168 , 0.258 0.169 , 0.256	Depositor DCC
R_{free} test set	1905 reflections (4.47%)	wwPDB-VP
Wilson B-factor (Å ²)	35.6	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 57.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.015 for -h,-k,-k+l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10096	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.58% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 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 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	0.38	0/816	0.57	0/1099
1	B	0.36	0/812	0.54	0/1093
1	C	0.38	0/866	0.55	0/1168
1	D	0.38	0/805	0.55	0/1083
1	E	0.34	0/820	0.54	0/1103
1	F	0.35	0/867	0.56	0/1169
1	G	0.36	0/833	0.54	0/1121
1	H	0.31	0/827	0.51	0/1112
1	I	0.34	0/822	0.51	0/1107
1	J	0.35	0/863	0.58	0/1163
1	K	0.34	0/786	0.56	0/1056
1	L	0.32	0/808	0.53	0/1087
All	All	0.35	0/9925	0.55	0/13361

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 the 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 within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	806	0	826	8	0
1	B	802	0	827	3	0
1	C	854	0	873	14	0
1	D	795	0	828	7	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	810	0	831	7	0
1	F	855	0	879	8	0
1	G	823	0	856	8	0
1	H	817	0	845	7	0
1	I	812	0	837	12	0
1	J	851	0	880	10	0
1	K	777	0	804	17	0
1	L	798	0	823	22	0
2	A	38	0	0	2	0
2	B	29	0	0	0	0
2	C	39	0	0	0	0
2	D	24	0	0	0	0
2	E	22	0	0	0	0
2	F	35	0	0	0	0
2	G	23	0	0	0	0
2	H	17	0	0	0	0
2	I	17	0	0	1	0
2	J	21	0	0	0	0
2	K	16	0	0	1	0
2	L	15	0	0	1	0
All	All	10096	0	10109	118	0

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 6.

All (118) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:106:LEU:HD12	1:J:182:ILE:HG12	1.70	0.72
1:K:119:LEU:HD11	1:K:167:LEU:HD13	1.75	0.69
1:L:115:SER:HA	1:L:173:LYS:HD2	1.75	0.68
1:G:115:SER:HA	1:G:173:LYS:HD2	1.77	0.66
1:L:149:ASN:O	1:L:153:ASN:ND2	2.25	0.65
1:I:109:SER:H	1:I:179[B]:ASP:HB3	1.62	0.64
1:K:104:ARG:NH2	2:K:201:HOH:O	2.31	0.62
1:G:104:ARG:HD3	1:G:143:VAL:HG21	1.82	0.62
1:D:104:ARG:HD3	1:D:143:VAL:HG21	1.81	0.61
1:K:104:ARG:HD2	1:K:183:GLN:OXT	2.01	0.60
1:L:86:LEU:HD11	1:L:113:GLU:HG3	1.84	0.59
1:K:160:LYS:HE3	1:K:172:ILE:O	2.03	0.58
1:C:135:ILE:O	1:C:139[A]:SER:OG	2.22	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:138:LEU:HD23	1:E:155:LEU:HD11	1.88	0.56
1:L:93:PHE:O	1:L:107:LYS:HA	2.06	0.56
1:G:109:SER:HB2	1:G:179[B]:ASP:HB2	1.88	0.54
1:K:85:PRO:HB2	1:K:114:LEU:HB2	1.90	0.54
1:L:150:LYS:N	1:L:150:LYS:HD2	2.23	0.54
1:J:77:ASP:O	1:J:81:LEU:HD23	2.07	0.54
1:A:124:LYS:NZ	2:A:204:HOH:O	2.40	0.54
1:L:150:LYS:HD2	1:L:150:LYS:H	1.73	0.53
1:L:143:VAL:HG23	1:L:144:GLU:H	1.73	0.53
1:I:97:LEU:HD11	1:I:106:LEU:HB2	1.91	0.53
1:B:108:ALA:HA	1:B:179[A]:ASP:O	2.09	0.53
1:G:117:GLU:HG2	1:K:98:VAL:HG11	1.90	0.53
1:E:87:TYR:OH	1:E:127:ASP:OD1	2.21	0.52
1:H:90[A]:ASP:OD1	1:H:90[A]:ASP:N	2.42	0.52
1:F:108:ALA:HA	1:F:179[A]:ASP:O	2.10	0.51
1:J:156:LYS:O	1:J:160:LYS:HG2	2.09	0.51
1:I:181:ILE:HD11	1:K:105:TYR:CD2	2.45	0.51
1:L:85:PRO:HD2	1:L:114:LEU:O	2.10	0.51
1:J:108:ALA:HA	1:J:179[A]:ASP:O	2.11	0.50
1:K:104:ARG:HH11	1:K:183:GLN:C	2.19	0.50
1:D:122:GLU:OE1	1:D:169:ASP:HB2	2.11	0.50
1:L:126:LYS:O	1:L:130:ILE:HG13	2.12	0.50
1:G:121:ASN:CG	1:I:144:GLU:HG2	2.36	0.50
1:K:148:THR:OG1	1:K:150:LYS:HB3	2.11	0.50
1:F:142:SER:OG	1:F:145:GLU:HG3	2.12	0.50
1:K:97:LEU:HD11	1:K:106:LEU:HB2	1.92	0.49
1:I:181:ILE:HD11	1:K:105:TYR:HD2	1.76	0.49
1:L:104:ARG:NH1	1:L:183:GLN:OXT	2.45	0.49
1:D:146:VAL:HG23	1:D:182:ILE:HD11	1.95	0.49
1:K:168:ILE:HG13	1:K:169:ASP:N	2.27	0.49
1:F:142:SER:HG	1:F:145:GLU:HG3	1.77	0.49
1:D:127:ASP:O	1:D:131:LYS:HG3	2.14	0.48
1:E:122:GLU:HG2	1:E:126:LYS:HE3	1.96	0.48
1:E:152:LYS:HE3	1:E:180:PHE:CE1	2.49	0.48
1:F:119:LEU:O	1:F:123:VAL:HG13	2.14	0.48
1:H:168:ILE:HG13	1:H:169:ASP:OD1	2.14	0.48
1:G:119:LEU:O	1:G:123:VAL:HG13	2.15	0.47
1:L:141:LYS:HG3	1:L:155:LEU:HD13	1.96	0.47
1:D:122:GLU:HG2	1:D:126:LYS:HE3	1.96	0.47
1:A:146:VAL:HG23	1:A:182:ILE:HD11	1.97	0.46
1:C:108:ALA:HA	1:C:179[A]:ASP:O	2.15	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:89:ILE:HD13	1:G:93:PHE:HZ	1.80	0.46
1:B:127:ASP:OD2	1:B:131:LYS:HE3	2.15	0.46
1:A:104:ARG:HD3	1:A:143:VAL:HG21	1.96	0.46
1:F:93:PHE:O	1:F:107:LYS:HA	2.16	0.46
1:D:97:LEU:HD11	1:D:106:LEU:HB2	1.97	0.46
1:K:93:PHE:O	1:K:107:LYS:HA	2.16	0.46
2:A:201:HOH:O	1:C:139[A]:SER:HB2	2.15	0.46
1:B:119:LEU:O	1:B:123:VAL:HG13	2.14	0.46
1:C:79:PHE:CE1	1:C:81:LEU:HB2	2.50	0.46
1:I:115:SER:HB3	1:I:171:PHE:CE1	2.51	0.46
1:C:146:VAL:HG23	1:C:182:ILE:HD11	1.97	0.46
1:H:95:VAL:HG21	1:H:138:LEU:HB3	1.97	0.46
1:C:164:ASN:ND2	1:C:171:PHE:HB2	2.32	0.45
1:I:138:LEU:HD23	1:I:155:LEU:HD11	1.97	0.45
1:K:122:GLU:HG2	1:K:126:LYS:HD2	1.98	0.45
1:L:119:LEU:HD22	1:L:170:GLY:HA3	1.98	0.45
1:D:114:LEU:HB3	1:D:119:LEU:HD23	1.98	0.45
1:J:131:LYS:O	1:J:135:ILE:HG13	2.17	0.45
1:K:107:LYS:HB3	1:K:181:ILE:HG13	1.99	0.45
1:A:137:ILE:HD13	1:A:158:GLU:HG2	1.99	0.44
1:I:118:LYS:HD2	1:I:169:ASP:HB3	1.99	0.44
1:L:86:LEU:HD12	1:L:113:GLU:HA	2.00	0.44
1:L:160:LYS:HG3	1:L:172:ILE:HB	2.00	0.44
1:K:83:LEU:HA	1:K:113:GLU:OE2	2.17	0.44
1:H:118:LYS:HG2	1:H:169:ASP:HB3	2.00	0.43
1:F:124:LYS:HB2	1:F:124:LYS:HE3	1.60	0.43
1:H:145:GLU:O	1:H:151:GLY:HA3	2.19	0.43
1:J:86:LEU:HD23	1:J:86:LEU:HA	1.78	0.43
1:J:93:PHE:O	1:J:107:LYS:HA	2.19	0.43
1:L:86:LEU:CD1	1:L:113:GLU:HG3	2.48	0.43
1:I:109:SER:HB2	1:I:179[B]:ASP:HB2	2.00	0.43
1:L:141:LYS:HD3	1:L:141:LYS:HA	1.74	0.43
1:A:149:ASN:CG	1:C:136:GLU:HG2	2.44	0.43
1:J:115:SER:HB3	1:J:171:PHE:O	2.18	0.43
1:G:123:VAL:HB	1:G:130:ILE:HD11	2.00	0.42
1:L:127:ASP:OD1	2:L:201:HOH:O	2.21	0.42
1:I:82:VAL:N	2:I:202:HOH:O	2.52	0.42
1:C:142:SER:OG	1:C:145:GLU:HG3	2.20	0.42
1:I:123:VAL:HG13	1:I:130:ILE:HD13	2.02	0.42
1:C:93:PHE:O	1:C:107:LYS:HA	2.20	0.42
1:C:119:LEU:HD22	1:C:170:GLY:HA3	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:83:LEU:HD12	1:J:83:LEU:HA	1.88	0.42
1:C:124:LYS:HE3	1:C:124:LYS:HB2	1.48	0.42
1:A:98:VAL:O	1:A:142:SER:HA	2.21	0.41
1:E:113:GLU:OE1	1:E:174:ASN:ND2	2.46	0.41
1:L:155:LEU:HD23	1:L:177:PHE:CZ	2.55	0.41
1:A:142:SER:OG	1:A:145:GLU:HG3	2.20	0.41
1:H:86:LEU:HD23	1:H:86:LEU:HA	1.92	0.41
1:K:163:LEU:O	1:K:167:LEU:HG	2.20	0.41
1:C:144:GLU:H	1:C:144:GLU:HG3	1.60	0.41
1:F:104:ARG:HD3	1:F:143:VAL:HG21	2.02	0.41
1:F:123:VAL:HG12	1:F:167:LEU:HD22	2.03	0.41
1:C:147:VAL:HG23	1:C:148:THR:HG23	2.03	0.41
1:H:152:LYS:O	1:H:156:LYS:HG3	2.21	0.41
1:I:163:LEU:O	1:I:167:LEU:HG	2.21	0.41
1:A:148:THR:O	1:A:152:LYS:HG3	2.20	0.41
1:C:148:THR:OG1	1:C:150:LYS:HG3	2.21	0.40
1:E:127:ASP:OD2	1:E:131:LYS:HE3	2.21	0.40
1:J:163:LEU:HD23	1:J:163:LEU:HA	1.90	0.40
1:L:143:VAL:HG23	1:L:144:GLU:N	2.36	0.40
1:L:122:GLU:OE1	1:L:169:ASP:HB2	2.22	0.40
1:E:109:SER:HB2	1:E:179[A]:ASP:HB3	2.03	0.40
1:L:109:SER:O	1:L:178:THR:HG23	2.21	0.40
1:L:167:LEU:HA	1:L:167:LEU:HD23	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 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 X-ray entries followed by that with respect to entries of similar resolution.

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	101/113 (89%)	98 (97%)	3 (3%)	0	100	100
1	B	101/113 (89%)	92 (91%)	9 (9%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	108/113 (96%)	106 (98%)	2 (2%)	0	100	100
1	D	99/113 (88%)	95 (96%)	4 (4%)	0	100	100
1	E	101/113 (89%)	96 (95%)	5 (5%)	0	100	100
1	F	107/113 (95%)	104 (97%)	3 (3%)	0	100	100
1	G	103/113 (91%)	100 (97%)	3 (3%)	0	100	100
1	H	101/113 (89%)	93 (92%)	8 (8%)	0	100	100
1	I	102/113 (90%)	94 (92%)	8 (8%)	0	100	100
1	J	107/113 (95%)	103 (96%)	4 (4%)	0	100	100
1	K	95/113 (84%)	90 (95%)	5 (5%)	0	100	100
1	L	100/113 (88%)	93 (93%)	7 (7%)	0	100	100
All	All	1225/1356 (90%)	1164 (95%)	61 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 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 X-ray entries followed by that with respect to entries of similar resolution.

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	92/103 (89%)	88 (96%)	4 (4%)	26	60
1	B	91/103 (88%)	91 (100%)	0	100	100
1	C	97/103 (94%)	92 (95%)	5 (5%)	21	53
1	D	91/103 (88%)	91 (100%)	0	100	100
1	E	93/103 (90%)	93 (100%)	0	100	100
1	F	98/103 (95%)	98 (100%)	0	100	100
1	G	94/103 (91%)	93 (99%)	1 (1%)	65	87
1	H	94/103 (91%)	93 (99%)	1 (1%)	65	87
1	I	92/103 (89%)	90 (98%)	2 (2%)	45	78
1	J	96/103 (93%)	96 (100%)	0	100	100
1	K	88/103 (85%)	87 (99%)	1 (1%)	65	87

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	91/103 (88%)	90 (99%)	1 (1%)	65	87
All	All	1117/1236 (90%)	1102 (99%)	15 (1%)	70	86

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	136[A]	GLU
1	A	136[B]	GLU
1	A	139[A]	SER
1	A	139[B]	SER
1	C	83	LEU
1	C	139[A]	SER
1	C	139[B]	SER
1	C	179[A]	ASP
1	C	179[B]	ASP
1	G	123	VAL
1	H	144	GLU
1	I	136[A]	GLU
1	I	136[B]	GLU
1	K	148	THR
1	L	178	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	121	ASN
1	H	100	GLN
1	K	149	ASN
1	K	162	HIS
1	L	174	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	101/113 (89%)	-0.78	1 (0%) 79 72	11, 24, 52, 111	2 (1%)
1	B	101/113 (89%)	-0.64	2 (1%) 65 56	11, 29, 55, 122	2 (1%)
1	C	108/113 (95%)	-0.88	1 (0%) 81 74	10, 22, 48, 85	2 (1%)
1	D	101/113 (89%)	-0.76	1 (0%) 79 72	13, 26, 52, 91	0
1	E	100/113 (88%)	-0.56	1 (1%) 79 72	12, 30, 67, 116	3 (3%)
1	F	108/113 (95%)	-0.66	0 100 100	12, 28, 57, 102	1 (0%)
1	G	103/113 (91%)	-0.66	2 (1%) 66 57	10, 25, 55, 90	2 (1%)
1	H	101/113 (89%)	-0.62	1 (0%) 79 72	13, 32, 54, 104	2 (1%)
1	I	102/113 (90%)	-0.56	1 (0%) 79 72	12, 33, 56, 102	2 (1%)
1	J	108/113 (95%)	-0.33	2 (1%) 66 57	12, 36, 63, 102	1 (0%)
1	K	98/113 (86%)	-0.07	1 (1%) 79 72	17, 47, 68, 75	1 (1%)
1	L	101/113 (89%)	-0.04	0 100 100	19, 52, 73, 92	1 (0%)
All	All	1232/1356 (90%)	-0.55	13 (1%) 78 70	10, 31, 67, 122	19 (1%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	101	ASN	3.5
1	J	77	ASP	3.2
1	I	100	GLN	3.1
1	G	101	ASN	2.8
1	J	100	GLN	2.8
1	A	83	LEU	2.6
1	B	102	GLY	2.3
1	K	84	GLY	2.3
1	E	121[A]	ASN	2.2
1	C	101	ASN	2.1
1	D	83	LEU	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	G	100	GLN	2.0
1	H	101	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.