



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:18 PM UTC

PDB ID : 2II7 / pdb_00002ii7
Title : Anabaena sensory rhodopsin transducer
Authors : Vogeley, L.
Deposited on : 2006-09-27
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:

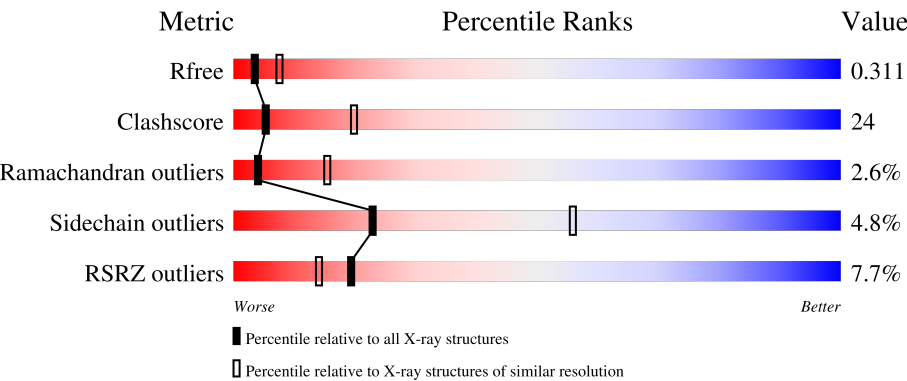
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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.



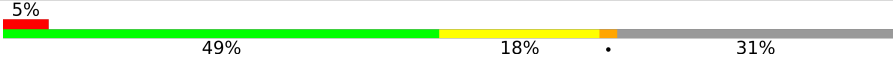
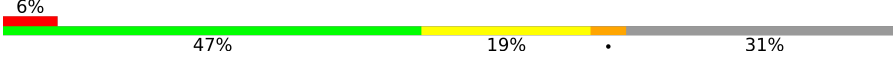
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	131	
1	B	131	
1	C	131	
1	D	131	
1	E	131	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	131	
1	G	131	
1	H	131	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6555 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 Anabaena sensory rhodopsin transducer protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	112	Total	C	N	O	S	0	0	0
			881	555	155	169	2			
1	B	115	Total	C	N	O	S	0	0	0
			905	571	159	173	2			
1	C	110	Total	C	N	O	S	0	0	0
			858	538	152	166	2			
1	D	103	Total	C	N	O	S	0	0	0
			809	511	142	154	2			
1	E	106	Total	C	N	O	S	0	0	0
			830	520	147	161	2			
1	F	91	Total	C	N	O	S	0	0	0
			714	453	123	136	2			
1	G	91	Total	C	N	O	S	0	0	0
			714	453	123	136	2			
1	H	106	Total	C	N	O	S	0	0	0
			829	520	146	161	2			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	125	HIS	-	expression tag	UNP Q8YSC3
A	126	HIS	-	expression tag	UNP Q8YSC3
A	127	HIS	-	expression tag	UNP Q8YSC3
A	128	HIS	-	expression tag	UNP Q8YSC3
A	129	HIS	-	expression tag	UNP Q8YSC3
A	130	HIS	-	expression tag	UNP Q8YSC3
B	125	HIS	-	expression tag	UNP Q8YSC3
B	126	HIS	-	expression tag	UNP Q8YSC3
B	127	HIS	-	expression tag	UNP Q8YSC3
B	128	HIS	-	expression tag	UNP Q8YSC3
B	129	HIS	-	expression tag	UNP Q8YSC3
B	130	HIS	-	expression tag	UNP Q8YSC3
C	125	HIS	-	expression tag	UNP Q8YSC3

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	126	HIS	-	expression tag	UNP Q8YSC3
C	127	HIS	-	expression tag	UNP Q8YSC3
C	128	HIS	-	expression tag	UNP Q8YSC3
C	129	HIS	-	expression tag	UNP Q8YSC3
C	130	HIS	-	expression tag	UNP Q8YSC3
D	125	HIS	-	expression tag	UNP Q8YSC3
D	126	HIS	-	expression tag	UNP Q8YSC3
D	127	HIS	-	expression tag	UNP Q8YSC3
D	128	HIS	-	expression tag	UNP Q8YSC3
D	129	HIS	-	expression tag	UNP Q8YSC3
D	130	HIS	-	expression tag	UNP Q8YSC3
E	125	HIS	-	expression tag	UNP Q8YSC3
E	126	HIS	-	expression tag	UNP Q8YSC3
E	127	HIS	-	expression tag	UNP Q8YSC3
E	128	HIS	-	expression tag	UNP Q8YSC3
E	129	HIS	-	expression tag	UNP Q8YSC3
E	130	HIS	-	expression tag	UNP Q8YSC3
F	125	HIS	-	expression tag	UNP Q8YSC3
F	126	HIS	-	expression tag	UNP Q8YSC3
F	127	HIS	-	expression tag	UNP Q8YSC3
F	128	HIS	-	expression tag	UNP Q8YSC3
F	129	HIS	-	expression tag	UNP Q8YSC3
F	130	HIS	-	expression tag	UNP Q8YSC3
G	125	HIS	-	expression tag	UNP Q8YSC3
G	126	HIS	-	expression tag	UNP Q8YSC3
G	127	HIS	-	expression tag	UNP Q8YSC3
G	128	HIS	-	expression tag	UNP Q8YSC3
G	129	HIS	-	expression tag	UNP Q8YSC3
G	130	HIS	-	expression tag	UNP Q8YSC3
H	125	HIS	-	expression tag	UNP Q8YSC3
H	126	HIS	-	expression tag	UNP Q8YSC3
H	127	HIS	-	expression tag	UNP Q8YSC3
H	128	HIS	-	expression tag	UNP Q8YSC3
H	129	HIS	-	expression tag	UNP Q8YSC3
H	130	HIS	-	expression tag	UNP Q8YSC3

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total O 2 2	0	0
2	B	2	Total O 2 2	0	0

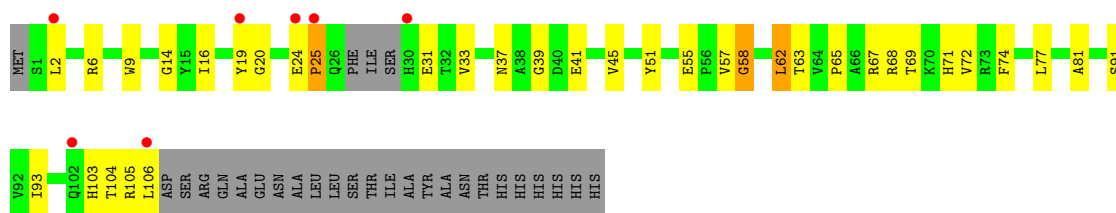
Continued on next page...

Continued from previous page...

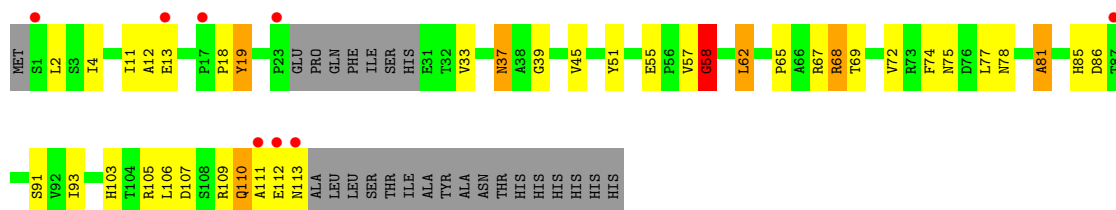
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	3	Total 3	O 3	0	0
2	E	5	Total 5	O 5	0	0
2	F	2	Total 2	O 2	0	0
2	H	1	Total 1	O 1	0	0

- Molecule 1: Anabaena sensory rhodopsin transducer protein

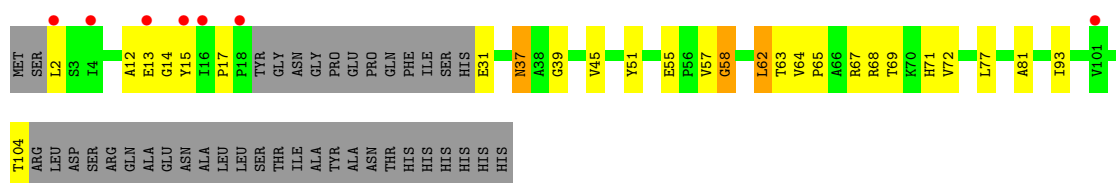




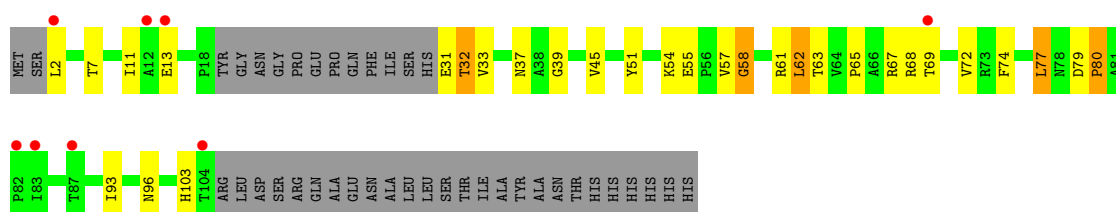
- Molecule 1: Anabaena sensory rhodopsin transducer protein



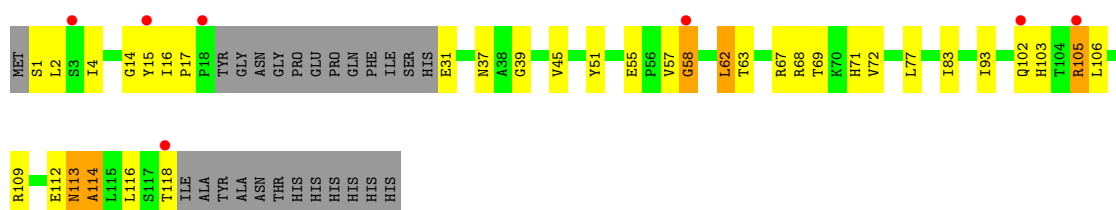
- Molecule 1: Anabaena sensory rhodopsin transducer protein



- Molecule 1: Anabaena sensory rhodopsin transducer protein



- Molecule 1: Anabaena sensory rhodopsin transducer protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.83Å 122.33Å 130.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 30.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.00-2.80) 99.7 (30.00-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.52 (at 2.61Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.259 , 0.315 0.257 , 0.311	Depositor DCC
R_{free} test set	2602 reflections (7.05%)	wwPDB-VP
Wilson B-factor (Å ²)	55.7	Xtriage
Anisotropy	0.137	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 56.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6555	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.44% 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	0.44	0/902	0.94	0/1233
1	B	0.50	0/930	1.05	4/1275 (0.3%)
1	C	0.52	0/878	1.00	2/1202 (0.2%)
1	D	0.51	0/832	0.98	4/1141 (0.4%)
1	E	0.61	0/851	1.05	4/1165 (0.3%)
1	F	0.50	0/733	0.96	3/1006 (0.3%)
1	G	0.44	0/733	0.95	1/1006 (0.1%)
1	H	0.47	0/848	0.99	1/1161 (0.1%)
All	All	0.50	0/6707	0.99	19/9189 (0.2%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	110	GLN	N-CA-C	-8.08	103.02	113.12
1	B	24	GLU	CA-C-N	7.57	127.52	120.03
1	B	24	GLU	C-N-CA	7.57	127.52	120.03
1	H	114	ALA	N-CA-C	-6.83	103.95	112.90
1	F	64	VAL	CA-C-N	6.42	126.44	119.89
1	F	64	VAL	C-N-CA	6.42	126.44	119.89
1	B	114	ALA	N-CA-C	-6.29	105.71	113.19
1	D	81	ALA	N-CA-C	6.13	113.74	108.22
1	E	58	GLY	CA-C-N	6.07	127.03	119.98
1	E	58	GLY	C-N-CA	6.07	127.03	119.98
1	D	91	SER	N-CA-C	5.59	117.68	109.24
1	D	16	ILE	CA-C-N	5.44	125.99	120.38
1	D	16	ILE	C-N-CA	5.44	125.99	120.38
1	C	58	GLY	CA-C-N	5.26	126.08	119.98
1	C	58	GLY	C-N-CA	5.26	126.08	119.98
1	E	91	SER	N-CA-C	5.13	117.18	109.23
1	G	32	THR	N-CA-C	5.09	117.12	109.23
1	E	81	ALA	N-CA-C	5.09	112.71	108.13
1	F	81	ALA	N-CA-C	5.05	112.77	108.22

There are no chirality outliers.

There are no planarity outliers.

5.2 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 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	881	0	862	58	0
1	B	905	0	872	61	0
1	C	858	0	838	53	0
1	D	809	0	780	37	0
1	E	830	0	799	42	0
1	F	714	0	692	31	0
1	G	714	0	692	33	0
1	H	829	0	810	42	0
2	A	2	0	0	1	0
2	B	2	0	0	1	0
2	D	3	0	0	1	0
2	E	5	0	0	0	0
2	F	2	0	0	1	0
2	H	1	0	0	0	0
All	All	6555	0	6345	313	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 24.

All (313) 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:A:2:LEU:HD12	1:A:2:LEU:H	1.33	0.92
1:H:2:LEU:HD23	1:H:4:ILE:HD11	1.54	0.90
1:A:30:HIS:CD2	1:A:73:ARG:HH12	1.91	0.87
1:C:31:GLU:HG2	1:C:109:ARG:NH2	1.90	0.86
1:C:31:GLU:HG2	1:C:109:ARG:HH21	1.41	0.86
1:H:31:GLU:HB3	1:H:103:HIS:HE1	1.45	0.81
1:C:110:GLN:NE2	1:C:113:ASN:HD22	1.77	0.80
1:C:29:SER:N	1:C:113:ASN:HD21	1.82	0.78
1:B:14:GLY:HA2	1:B:31:GLU:OE2	1.83	0.78

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:HIS:CD2	1:A:73:ARG:NH1	2.53	0.76
1:B:106:LEU:HD13	1:B:110:GLN:HB2	1.68	0.74
1:F:67:ARG:HH11	1:F:67:ARG:HB2	1.53	0.73
1:E:103:HIS:CE1	1:E:105:ARG:HD3	2.22	0.73
1:E:11:ILE:HD13	1:E:103:HIS:HB2	1.71	0.73
1:F:71:HIS:HE1	2:F:132:HOH:O	1.70	0.73
1:H:113:ASN:C	1:H:113:ASN:HD22	1.97	0.73
1:A:105:ARG:HH22	1:F:104:THR:HA	1.55	0.71
1:D:71:HIS:NE2	1:G:13:GLU:OE1	2.23	0.71
1:B:62:LEU:HD11	1:B:72:VAL:HG21	1.73	0.70
1:F:67:ARG:HB2	1:F:67:ARG:NH1	2.07	0.70
1:B:30:HIS:HB2	1:B:110:GLN:NE2	2.08	0.69
1:H:105:ARG:HD2	1:H:106:LEU:H	1.56	0.68
1:A:2:LEU:HB3	1:A:4:ILE:HD11	1.75	0.68
1:C:110:GLN:HE21	1:C:110:GLN:HA	1.58	0.68
1:H:14:GLY:HA2	1:H:31:GLU:OE2	1.93	0.67
1:D:104:THR:HA	2:D:132:HOH:O	1.95	0.67
1:B:39:GLY:HA3	2:B:131:HOH:O	1.95	0.66
1:F:57:VAL:O	1:F:58:GLY:O	2.13	0.66
1:C:62:LEU:HD11	1:C:72:VAL:HG21	1.77	0.66
1:H:105:ARG:HD2	1:H:106:LEU:N	2.10	0.66
1:A:105:ARG:NH1	1:A:105:ARG:HB2	2.11	0.66
1:F:62:LEU:HD11	1:F:72:VAL:HG21	1.78	0.66
1:G:62:LEU:HD11	1:G:72:VAL:HG21	1.77	0.65
1:E:103:HIS:HE1	1:E:105:ARG:HD3	1.61	0.65
1:G:67:ARG:HH11	1:G:67:ARG:HB2	1.60	0.65
1:C:116:LEU:HD13	1:E:106:LEU:HD13	1.78	0.65
1:F:37:ASN:ND2	1:F:39:GLY:H	1.93	0.65
1:C:112:GLU:OE2	1:E:105:ARG:HB2	1.97	0.65
1:E:62:LEU:HD11	1:E:72:VAL:HG21	1.77	0.65
1:B:109:ARG:O	1:B:112:GLU:HG2	1.98	0.64
1:G:67:ARG:HB2	1:G:67:ARG:NH1	2.12	0.64
1:D:45:VAL:HB	1:D:62:LEU:HD22	1.79	0.64
1:C:15:TYR:O	1:C:31:GLU:HG3	1.98	0.64
1:C:57:VAL:O	1:C:58:GLY:O	2.15	0.64
1:E:2:LEU:HD21	1:E:4:ILE:HD11	1.80	0.64
1:B:103:HIS:HD2	1:B:104:THR:N	1.94	0.63
1:C:29:SER:N	1:C:109:ARG:HD3	2.12	0.63
1:A:67:ARG:HB2	1:A:67:ARG:NH1	2.13	0.63
1:B:18:PRO:HB3	1:B:86:ASP:HB2	1.81	0.63
1:A:107:ASP:OD1	1:F:104:THR:HG23	1.99	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:116:LEU:C	1:H:118:THR:H	2.06	0.62
1:A:30:HIS:HD2	1:A:73:ARG:NH1	1.97	0.61
1:E:93:ILE:N	1:E:93:ILE:HD12	2.13	0.61
1:H:62:LEU:HD11	1:H:72:VAL:HG21	1.81	0.61
1:G:37:ASN:HD22	1:G:39:GLY:H	1.49	0.61
1:D:62:LEU:C	1:D:62:LEU:HD23	2.26	0.61
1:A:105:ARG:CB	1:A:105:ARG:HH11	2.14	0.60
1:G:37:ASN:ND2	1:G:39:GLY:H	1.98	0.60
1:A:3:SER:HB2	1:G:7:THR:OG1	2.02	0.60
1:A:57:VAL:O	1:A:58:GLY:O	2.18	0.60
1:B:93:ILE:HD12	1:B:93:ILE:N	2.17	0.60
1:A:62:LEU:HD11	1:A:72:VAL:HG21	1.84	0.60
1:D:62:LEU:HD11	1:D:72:VAL:HG21	1.82	0.60
1:F:45:VAL:HB	1:F:62:LEU:HD22	1.83	0.60
1:F:93:ILE:N	1:F:93:ILE:HD12	2.16	0.60
1:F:37:ASN:HD22	1:F:39:GLY:H	1.47	0.59
1:D:67:ARG:NH1	1:D:67:ARG:HB2	2.17	0.59
1:A:68:ARG:NH2	1:B:55:GLU:OE2	2.35	0.59
1:C:118:THR:O	1:C:119:ILE:HD13	2.03	0.59
1:A:117:SER:O	1:A:119:ILE:N	2.36	0.59
1:A:67:ARG:HB2	1:A:67:ARG:HH11	1.68	0.59
1:F:13:GLU:OE1	1:H:71:HIS:NE2	2.35	0.58
1:E:18:PRO:HB3	1:E:86:ASP:HB2	1.85	0.58
1:A:27:PHE:H	1:A:27:PHE:HD1	1.50	0.58
1:D:68:ARG:HH21	1:G:55:GLU:N	2.01	0.58
1:F:15:TYR:CE2	1:F:17:PRO:HG3	2.38	0.58
1:F:37:ASN:HD22	1:F:37:ASN:C	2.11	0.58
1:H:2:LEU:CD2	1:H:4:ILE:HD11	2.31	0.58
1:B:57:VAL:O	1:B:58:GLY:O	2.21	0.58
1:G:31:GLU:HG3	1:G:74:PHE:HD2	1.68	0.58
1:D:103:HIS:CD2	1:D:104:THR:H	2.21	0.58
1:C:51:TYR:HE2	1:C:57:VAL:HG23	1.68	0.57
1:C:110:GLN:NE2	1:C:113:ASN:ND2	2.51	0.57
1:B:106:LEU:HD12	1:B:106:LEU:O	2.04	0.57
1:C:1:SER:O	1:C:2:LEU:HD22	2.03	0.57
1:C:62:LEU:C	1:C:62:LEU:HD23	2.29	0.57
1:E:57:VAL:O	1:E:58:GLY:O	2.23	0.57
1:A:51:TYR:C	1:F:69:THR:HG22	2.30	0.57
1:F:62:LEU:C	1:F:62:LEU:HD23	2.29	0.57
1:H:31:GLU:HB3	1:H:103:HIS:CE1	2.35	0.57
1:B:19:TYR:CG	1:B:20:GLY:N	2.73	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:102:GLN:HB3	1:E:12:ALA:HB3	1.86	0.57
1:D:103:HIS:CD2	1:D:104:THR:N	2.73	0.56
1:B:105:ARG:HG3	1:B:105:ARG:HH11	1.68	0.56
1:C:110:GLN:HA	1:C:113:ASN:HD22	1.70	0.56
1:B:111:ALA:C	1:B:113:ASN:H	2.13	0.56
1:B:103:HIS:CD2	1:B:104:THR:N	2.73	0.56
1:A:37:ASN:ND2	1:A:39:GLY:H	2.03	0.56
1:A:93:ILE:N	1:A:93:ILE:HD12	2.21	0.56
1:F:14:GLY:HA2	1:F:31:GLU:OE1	2.06	0.56
1:A:62:LEU:C	1:A:62:LEU:HD23	2.30	0.55
1:C:62:LEU:HD23	1:C:63:THR:N	2.21	0.55
1:E:110:GLN:C	1:E:110:GLN:CD	2.74	0.55
1:G:65:PRO:HB2	1:G:68:ARG:HG3	1.89	0.55
1:B:113:ASN:C	1:B:115:LEU:H	2.13	0.55
1:C:29:SER:N	1:C:113:ASN:ND2	2.53	0.55
1:E:67:ARG:NH1	1:E:67:ARG:HB2	2.22	0.55
1:A:2:LEU:H	1:A:2:LEU:CD1	2.11	0.55
1:G:45:VAL:HB	1:G:62:LEU:HD22	1.89	0.55
1:H:112:GLU:O	1:H:116:LEU:HD13	2.07	0.55
1:H:57:VAL:O	1:H:58:GLY:O	2.25	0.55
1:C:71:HIS:NE2	1:E:13:GLU:OE1	2.36	0.55
1:E:67:ARG:HB2	1:E:67:ARG:HH11	1.71	0.55
1:A:69:THR:HG22	1:B:51:TYR:C	2.32	0.54
1:E:19:TYR:OH	1:E:75:ASN:ND2	2.37	0.54
1:F:65:PRO:HB2	1:F:68:ARG:HG3	1.90	0.54
1:G:62:LEU:HD23	1:G:62:LEU:C	2.31	0.54
1:A:45:VAL:HB	1:A:62:LEU:HD22	1.90	0.54
1:A:2:LEU:HB3	1:A:4:ILE:CD1	2.37	0.54
1:A:109:ARG:HG2	1:A:109:ARG:HH11	1.72	0.54
1:C:1:SER:C	1:C:2:LEU:HD22	2.33	0.53
1:D:14:GLY:HA2	1:D:31:GLU:OE2	2.08	0.53
1:D:67:ARG:HB2	1:D:67:ARG:HH11	1.72	0.53
1:A:102:GLN:HB3	1:B:12:ALA:HB3	1.89	0.53
1:D:105:ARG:O	1:D:106:LEU:C	2.51	0.53
1:B:113:ASN:C	1:B:115:LEU:N	2.67	0.53
1:E:62:LEU:C	1:E:62:LEU:HD23	2.33	0.53
1:D:51:TYR:C	1:E:69:THR:HG22	2.33	0.53
1:C:45:VAL:HB	1:C:62:LEU:HD22	1.91	0.53
1:D:93:ILE:HD12	1:D:93:ILE:N	2.24	0.53
1:A:71:HIS:NE2	1:B:13:GLU:OE1	2.42	0.53
1:A:28:ILE:HG23	1:A:113:ASN:OD1	2.09	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:ASN:HD22	1:A:39:GLY:H	1.56	0.52
1:E:37:ASN:ND2	1:E:39:GLY:H	2.08	0.52
1:C:116:LEU:CD1	1:E:106:LEU:HD13	2.40	0.52
1:B:68:ARG:CG	1:B:68:ARG:HH11	2.23	0.52
1:G:93:ILE:HD12	1:G:93:ILE:N	2.25	0.52
1:A:37:ASN:HD22	1:A:37:ASN:C	2.18	0.51
1:C:110:GLN:NE2	1:C:110:GLN:HA	2.23	0.51
1:E:107:ASP:O	1:E:110:GLN:HB3	2.09	0.51
1:C:67:ARG:HB2	1:C:67:ARG:NH1	2.25	0.51
1:A:51:TYR:HE2	1:A:57:VAL:HG23	1.76	0.51
1:A:62:LEU:HD23	1:A:63:THR:N	2.25	0.51
1:B:106:LEU:HD12	1:B:106:LEU:C	2.35	0.51
1:H:113:ASN:C	1:H:113:ASN:ND2	2.65	0.51
1:C:69:THR:HG22	1:E:51:TYR:C	2.35	0.51
1:D:37:ASN:C	1:D:37:ASN:HD22	2.19	0.51
1:C:93:ILE:N	1:C:93:ILE:HD12	2.26	0.51
1:D:33:VAL:HG23	1:D:74:PHE:CE2	2.46	0.51
1:C:67:ARG:HB2	1:C:67:ARG:HH11	1.76	0.50
1:C:15:TYR:HB3	1:C:106:LEU:HD21	1.94	0.50
1:C:29:SER:O	1:C:30:HIS:HB2	2.11	0.50
1:E:110:GLN:C	1:E:112:GLU:N	2.61	0.50
1:G:57:VAL:O	1:G:58:GLY:O	2.29	0.50
1:E:37:ASN:HD22	1:E:37:ASN:C	2.18	0.50
1:E:65:PRO:HB2	1:E:68:ARG:HG3	1.93	0.50
1:B:108:SER:HB3	1:H:15:TYR:CD2	2.47	0.50
1:C:37:ASN:ND2	1:C:39:GLY:H	2.08	0.50
1:D:65:PRO:HB2	1:D:68:ARG:HG3	1.93	0.50
1:B:115:LEU:O	1:B:115:LEU:HD12	2.11	0.50
1:D:68:ARG:HD3	1:G:51:TYR:O	2.11	0.50
1:E:37:ASN:HD22	1:E:39:GLY:H	1.59	0.50
1:H:37:ASN:ND2	1:H:39:GLY:H	2.10	0.50
1:H:62:LEU:HD23	1:H:63:THR:N	2.27	0.50
1:A:61:ARG:HH11	1:A:61:ARG:HG3	1.77	0.49
1:C:117:SER:O	1:C:119:ILE:N	2.44	0.49
1:D:2:LEU:HG	1:D:2:LEU:O	2.12	0.49
1:F:37:ASN:ND2	1:F:37:ASN:C	2.70	0.49
1:A:30:HIS:HA	1:A:109:ARG:HD2	1.95	0.49
1:C:51:TYR:C	1:G:69:THR:HG22	2.37	0.49
1:D:62:LEU:HD23	1:D:63:THR:N	2.28	0.49
1:E:45:VAL:HB	1:E:62:LEU:HD22	1.94	0.49
1:F:51:TYR:C	1:H:69:THR:HG22	2.38	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:PRO:HB2	1:B:68:ARG:HG3	1.93	0.49
1:B:51:TYR:HE2	1:B:57:VAL:HG23	1.78	0.48
1:H:62:LEU:HD23	1:H:62:LEU:C	2.37	0.48
1:H:67:ARG:HB2	1:H:67:ARG:NH1	2.27	0.48
1:A:68:ARG:HH21	1:B:55:GLU:HA	1.78	0.48
1:H:2:LEU:N	1:H:2:LEU:HD12	2.27	0.48
1:C:2:LEU:HD12	1:C:4:ILE:HD11	1.94	0.48
1:B:37:ASN:ND2	1:B:39:GLY:H	2.11	0.48
1:F:55:GLU:N	1:H:68:ARG:HH21	2.12	0.47
1:D:19:TYR:CE2	1:D:24:GLU:HG3	2.49	0.47
1:H:93:ILE:N	1:H:93:ILE:HD12	2.30	0.47
1:D:37:ASN:HD22	1:D:39:GLY:H	1.63	0.47
1:B:67:ARG:NH1	1:B:67:ARG:HB2	2.29	0.47
1:B:105:ARG:HG3	1:B:105:ARG:NH1	2.30	0.47
1:B:106:LEU:HD12	1:B:107:ASP:O	2.15	0.47
1:E:68:ARG:HH11	1:E:68:ARG:CG	2.28	0.47
1:B:68:ARG:NH2	1:H:55:GLU:OE2	2.48	0.47
1:F:67:ARG:HH11	1:F:67:ARG:CB	2.25	0.47
1:E:109:ARG:HG3	1:E:109:ARG:HH11	1.79	0.47
1:B:32:THR:HG22	1:B:33:VAL:N	2.29	0.46
1:C:6:ARG:HG2	1:C:9:TRP:CE2	2.50	0.46
1:D:55:GLU:OE2	1:E:68:ARG:NH2	2.48	0.46
1:B:109:ARG:HG3	1:B:109:ARG:HH11	1.81	0.46
1:D:68:ARG:HH21	1:G:55:GLU:CA	2.28	0.46
1:A:105:ARG:NH1	1:A:105:ARG:CB	2.75	0.46
1:C:29:SER:O	1:C:30:HIS:CB	2.64	0.46
1:D:68:ARG:HH21	1:G:54:LYS:C	2.22	0.46
1:G:62:LEU:HD23	1:G:63:THR:N	2.31	0.46
1:B:109:ARG:HG3	1:B:109:ARG:NH1	2.31	0.46
1:D:57:VAL:O	1:D:58:GLY:O	2.33	0.46
1:B:62:LEU:HD23	1:B:62:LEU:C	2.41	0.46
1:H:105:ARG:HG2	1:H:105:ARG:HH11	1.80	0.46
1:E:110:GLN:O	1:E:111:ALA:C	2.57	0.45
1:B:35:ILE:HG22	1:B:36:LEU:N	2.31	0.45
1:A:117:SER:O	1:A:118:THR:C	2.59	0.45
1:C:68:ARG:NH2	1:E:55:GLU:OE2	2.50	0.45
1:C:79:ASP:HA	1:C:80:PRO:C	2.42	0.45
1:E:109:ARG:HG2	1:E:109:ARG:O	2.16	0.45
1:D:67:ARG:NH1	1:G:55:GLU:OE2	2.50	0.45
1:B:41:GLU:OE2	1:D:41:GLU:OE2	2.34	0.45
1:B:67:ARG:HB2	1:B:67:ARG:HH11	1.82	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:51:TYR:HE2	1:F:57:VAL:HG23	1.81	0.45
1:D:103:HIS:HD2	1:D:104:THR:H	1.62	0.45
1:B:69:THR:HG22	1:H:51:TYR:C	2.42	0.45
1:B:103:HIS:CD2	1:B:103:HIS:C	2.95	0.45
1:C:112:GLU:O	1:C:116:LEU:HB2	2.16	0.45
1:A:2:LEU:HD23	1:A:4:ILE:HD11	1.98	0.44
1:B:17:PRO:HB2	1:B:28:ILE:HD11	1.98	0.44
1:E:112:GLU:O	1:E:113:ASN:CG	2.60	0.44
1:C:79:ASP:OD1	1:C:80:PRO:HA	2.17	0.44
1:E:111:ALA:C	1:E:113:ASN:H	2.25	0.44
1:A:13:GLU:OE2	1:A:106:LEU:HD23	2.17	0.44
1:B:111:ALA:C	1:B:113:ASN:N	2.75	0.44
1:C:68:ARG:HH11	1:C:68:ARG:CG	2.30	0.44
1:D:37:ASN:ND2	1:D:39:GLY:H	2.16	0.44
1:D:69:THR:HG22	1:G:51:TYR:C	2.43	0.44
1:A:37:ASN:ND2	1:A:37:ASN:C	2.74	0.44
1:A:65:PRO:HB2	1:A:68:ARG:HG3	2.00	0.44
1:B:25:PRO:O	1:B:30:HIS:HE1	2.01	0.44
1:A:55:GLU:OE2	1:F:68:ARG:NH2	2.51	0.44
1:A:83:ILE:HD12	1:A:83:ILE:N	2.33	0.44
1:C:30:HIS:CD2	1:C:73:ARG:HH22	2.35	0.44
1:A:105:ARG:NH2	1:A:107:ASP:OD2	2.51	0.44
1:B:17:PRO:HD3	1:B:29:SER:HB2	2.00	0.44
1:B:74:PHE:O	1:B:77:LEU:HB2	2.18	0.44
1:H:51:TYR:HE2	1:H:57:VAL:HG23	1.82	0.44
1:H:116:LEU:C	1:H:118:THR:N	2.73	0.44
1:A:118:THR:O	1:A:119:ILE:C	2.60	0.44
1:B:37:ASN:HD22	1:B:39:GLY:H	1.65	0.44
1:A:109:ARG:HG2	1:A:109:ARG:NH1	2.33	0.44
1:B:68:ARG:HH21	1:H:55:GLU:N	2.16	0.44
1:A:68:ARG:HH21	1:B:55:GLU:CA	2.29	0.43
1:F:15:TYR:HE2	1:F:17:PRO:HG3	1.83	0.43
1:A:105:ARG:O	1:A:106:LEU:C	2.61	0.43
1:C:109:ARG:NH1	1:C:110:GLN:OE1	2.51	0.43
1:F:62:LEU:HD23	1:F:63:THR:N	2.34	0.43
1:A:33:VAL:HG23	1:A:74:PHE:CE2	2.54	0.43
1:A:36:LEU:HD12	1:A:37:ASN:H	1.83	0.43
1:G:79:ASP:HA	1:G:80:PRO:C	2.44	0.43
1:B:67:ARG:NH1	1:H:55:GLU:OE2	2.52	0.43
1:B:86:ASP:OD1	1:B:86:ASP:O	2.37	0.43
1:H:1:SER:HB3	1:H:2:LEU:HD12	2.00	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:115:LEU:O	1:B:115:LEU:CD1	2.67	0.43
1:C:17:PRO:HA	1:C:18:PRO:HD3	1.94	0.43
1:E:110:GLN:C	1:E:110:GLN:NE2	2.76	0.43
1:H:83:ILE:HD12	1:H:83:ILE:N	2.34	0.43
1:A:61:ARG:HG3	1:A:61:ARG:NH1	2.34	0.43
1:G:74:PHE:O	1:G:77:LEU:HB2	2.19	0.43
1:C:30:HIS:CG	1:C:73:ARG:HH12	2.35	0.43
1:D:68:ARG:HH21	1:G:55:GLU:HA	1.84	0.43
1:E:37:ASN:ND2	1:E:37:ASN:C	2.76	0.43
1:G:11:ILE:HG21	1:G:103:HIS:CD2	2.53	0.43
1:B:62:LEU:HD23	1:B:63:THR:N	2.33	0.43
1:C:112:GLU:C	1:C:114:ALA:H	2.25	0.43
1:F:12:ALA:HB3	1:H:102:GLN:HB3	2.01	0.42
1:G:51:TYR:HE2	1:G:57:VAL:HG23	1.84	0.42
1:B:37:ASN:HD22	1:B:37:ASN:C	2.26	0.42
1:E:33:VAL:HG23	1:E:74:PHE:CE2	2.54	0.42
1:G:61:ARG:HG3	1:G:61:ARG:HH11	1.84	0.42
1:A:2:LEU:HB3	1:A:4:ILE:CG1	2.49	0.42
1:E:110:GLN:O	1:E:110:GLN:NE2	2.53	0.42
1:D:37:ASN:C	1:D:37:ASN:ND2	2.77	0.42
1:H:113:ASN:HD22	1:H:114:ALA:N	2.16	0.42
1:E:85:HIS:O	1:E:86:ASP:C	2.63	0.42
1:A:2:LEU:HD13	1:A:38:ALA:HB1	2.01	0.42
1:C:13:GLU:OE2	1:C:106:LEU:HD23	2.18	0.42
1:D:68:ARG:CG	1:D:68:ARG:HH11	2.33	0.42
1:E:109:ARG:HG3	1:E:109:ARG:NH1	2.35	0.42
1:A:105:ARG:HG2	2:A:132:HOH:O	2.18	0.42
1:C:83:ILE:HD12	1:C:83:ILE:N	2.35	0.42
1:F:68:ARG:HH11	1:F:68:ARG:CG	2.33	0.42
1:H:31:GLU:CB	1:H:103:HIS:HE1	2.24	0.42
1:B:29:SER:C	1:B:31:GLU:H	2.28	0.41
1:B:45:VAL:HB	1:B:62:LEU:HD22	2.02	0.41
1:A:97:VAL:HG12	1:G:96:ASN:O	2.20	0.41
1:B:31:GLU:HG3	1:B:103:HIS:NE2	2.35	0.41
1:G:32:THR:HG22	1:G:33:VAL:N	2.34	0.41
1:A:55:GLU:N	1:F:68:ARG:HH21	2.19	0.41
1:G:37:ASN:HD22	1:G:37:ASN:C	2.29	0.41
1:H:16:ILE:HA	1:H:17:PRO:HD3	1.84	0.41
1:C:51:TYR:HE2	1:C:57:VAL:CG2	2.31	0.41
1:F:55:GLU:CA	1:H:68:ARG:HH21	2.34	0.41
1:B:37:ASN:ND2	1:B:37:ASN:C	2.77	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:45:VAL:HB	1:H:62:LEU:HD22	2.02	0.41
1:F:93:ILE:N	1:F:93:ILE:CD1	2.84	0.41
1:H:109:ARG:HB2	1:H:109:ARG:NH1	2.35	0.41
1:B:24:GLU:HA	1:B:25:PRO:HD3	1.72	0.41
1:C:32:THR:HG22	1:C:33:VAL:N	2.36	0.41
1:C:55:GLU:OE2	1:G:67:ARG:NH1	2.54	0.41
1:D:24:GLU:HA	1:D:25:PRO:HD2	1.87	0.41
1:E:78:ASN:O	1:E:81:ALA:HA	2.21	0.41
1:H:67:ARG:HB2	1:H:67:ARG:HH11	1.86	0.41
1:H:37:ASN:HD22	1:H:39:GLY:H	1.69	0.40
1:A:27:PHE:CD1	1:A:27:PHE:N	2.88	0.40
1:B:68:ARG:HH11	1:B:68:ARG:HG3	1.86	0.40
1:C:55:GLU:N	1:G:68:ARG:HH21	2.19	0.40
1:D:51:TYR:HE2	1:D:57:VAL:HG23	1.86	0.40
1:G:68:ARG:HH11	1:G:68:ARG:CG	2.34	0.40
1:H:68:ARG:HH11	1:H:68:ARG:CG	2.34	0.40
1:D:6:ARG:HG2	1:D:9:TRP:CE2	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 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 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	108/131 (82%)	94 (87%)	10 (9%)	4 (4%)	2	9
1	B	113/131 (86%)	99 (88%)	12 (11%)	2 (2%)	6	23
1	C	106/131 (81%)	97 (92%)	3 (3%)	6 (6%)	1	4
1	D	99/131 (76%)	89 (90%)	7 (7%)	3 (3%)	3	12
1	E	102/131 (78%)	91 (89%)	9 (9%)	2 (2%)	6	21
1	F	87/131 (66%)	80 (92%)	6 (7%)	1 (1%)	11	36
1	G	87/131 (66%)	81 (93%)	4 (5%)	2 (2%)	5	18

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	102/131 (78%)	93 (91%)	8 (8%)	1 (1%)	12	38
All	All	804/1048 (77%)	724 (90%)	59 (7%)	21 (3%)	4	15

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	31	GLU
1	A	58	GLY
1	A	118	THR
1	B	58	GLY
1	C	58	GLY
1	D	58	GLY
1	E	58	GLY
1	F	58	GLY
1	G	58	GLY
1	H	58	GLY
1	B	29	SER
1	C	118	THR
1	E	19	TYR
1	C	30	HIS
1	D	20	GLY
1	C	31	GLU
1	A	106	LEU
1	D	25	PRO
1	C	119	ILE
1	G	80	PRO
1	C	80	PRO

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	99/114 (87%)	91 (92%)	8 (8%)	11	33
1	B	100/114 (88%)	94 (94%)	6 (6%)	17	47
1	C	96/114 (84%)	93 (97%)	3 (3%)	35	70

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	90/114 (79%)	88 (98%)	2 (2%)	45	78
1	E	92/114 (81%)	87 (95%)	5 (5%)	20	51
1	F	80/114 (70%)	76 (95%)	4 (5%)	22	54
1	G	80/114 (70%)	77 (96%)	3 (4%)	29	64
1	H	93/114 (82%)	89 (96%)	4 (4%)	26	60
All	All	730/912 (80%)	695 (95%)	35 (5%)	23	56

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	3	SER
1	A	30	HIS
1	A	37	ASN
1	A	62	LEU
1	A	77	LEU
1	A	105	ARG
1	A	116	LEU
1	B	28	ILE
1	B	62	LEU
1	B	68	ARG
1	B	77	LEU
1	B	106	LEU
1	B	116	LEU
1	C	62	LEU
1	C	68	ARG
1	C	77	LEU
1	D	62	LEU
1	D	77	LEU
1	E	37	ASN
1	E	62	LEU
1	E	68	ARG
1	E	77	LEU
1	E	110	GLN
1	F	2	LEU
1	F	37	ASN
1	F	62	LEU
1	F	77	LEU
1	G	2	LEU
1	G	62	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	G	77	LEU
1	H	62	LEU
1	H	77	LEU
1	H	105	ARG
1	H	113	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	85	HIS
1	A	94	GLN
1	A	102	GLN
1	B	37	ASN
1	B	85	HIS
1	B	94	GLN
1	B	102	GLN
1	B	110	GLN
1	B	113	ASN
1	C	37	ASN
1	C	85	HIS
1	C	110	GLN
1	D	37	ASN
1	D	85	HIS
1	D	94	GLN
1	D	102	GLN
1	E	37	ASN
1	E	85	HIS
1	E	94	GLN
1	E	102	GLN
1	E	110	GLN
1	F	37	ASN
1	F	85	HIS
1	F	94	GLN
1	F	102	GLN
1	G	37	ASN
1	G	85	HIS
1	G	94	GLN
1	G	102	GLN
1	H	37	ASN
1	H	85	HIS
1	H	94	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	H	113	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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	112/131 (85%)	0.68	6 (5%) 31 24	40, 69, 91, 97	0
1	B	115/131 (87%)	0.54	9 (7%) 19 14	32, 56, 95, 113	0
1	C	110/131 (83%)	0.76	12 (10%) 10 8	20, 62, 97, 113	0
1	D	103/131 (78%)	0.39	7 (6%) 23 17	24, 50, 96, 120	0
1	E	106/131 (80%)	0.25	8 (7%) 20 15	17, 39, 101, 121	0
1	F	91/131 (69%)	0.54	7 (7%) 19 14	28, 53, 87, 108	0
1	G	91/131 (69%)	0.90	8 (8%) 15 11	43, 69, 102, 113	0
1	H	106/131 (80%)	0.36	7 (6%) 24 18	22, 56, 100, 113	0
All	All	834/1048 (79%)	0.55	64 (7%) 19 14	17, 59, 98, 121	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	119	ILE	5.5
1	C	117	SER	4.9
1	C	113	ASN	4.9
1	D	106	LEU	4.7
1	C	116	LEU	4.5
1	G	2	LEU	4.2
1	F	18	PRO	4.1
1	H	58	GLY	3.8
1	B	116	LEU	3.7
1	G	104	THR	3.6
1	G	13	GLU	3.5
1	C	16	ILE	3.5
1	B	2	LEU	3.5
1	D	30	HIS	3.3
1	A	13	GLU	3.2
1	F	4	ILE	3.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	120	ALA	3.1
1	G	12	ALA	3.1
1	C	15	TYR	3.0
1	B	114	ALA	3.0
1	H	118	THR	3.0
1	E	111	ALA	3.0
1	E	112	GLU	2.9
1	F	2	LEU	2.9
1	F	15	TYR	2.9
1	B	88	ASP	2.8
1	B	58	GLY	2.8
1	C	1	SER	2.8
1	D	25	PRO	2.7
1	C	110	GLN	2.7
1	H	15	TYR	2.6
1	A	62	LEU	2.6
1	C	13	GLU	2.6
1	D	24	GLU	2.6
1	E	1	SER	2.6
1	C	112	GLU	2.6
1	D	2	LEU	2.4
1	G	82	PRO	2.4
1	A	15	TYR	2.4
1	B	83	ILE	2.4
1	G	83	ILE	2.4
1	E	23	PRO	2.4
1	C	56	PRO	2.4
1	B	57	VAL	2.4
1	B	87	THR	2.4
1	G	87	THR	2.3
1	F	101	VAL	2.3
1	E	87	THR	2.3
1	B	19	TYR	2.3
1	E	17	PRO	2.3
1	E	13	GLU	2.3
1	A	117	SER	2.2
1	E	113	ASN	2.2
1	F	13	GLU	2.2
1	G	69	THR	2.2
1	H	102	GLN	2.1
1	H	3	SER	2.1
1	D	102	GLN	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	H	18	PRO	2.1
1	C	2	LEU	2.1
1	F	16	ILE	2.1
1	A	65	PRO	2.1
1	D	19	TYR	2.0
1	H	105	ARG	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.