



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 11:24 am BST

PDB ID : 4LTQ
Title : Bacterial sodium channel in low calcium, P42 space group
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Nurva, S.; Loussouarn, G.; Minor, D.L.
Deposited on : 2013-07-23
Resolution : 5.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

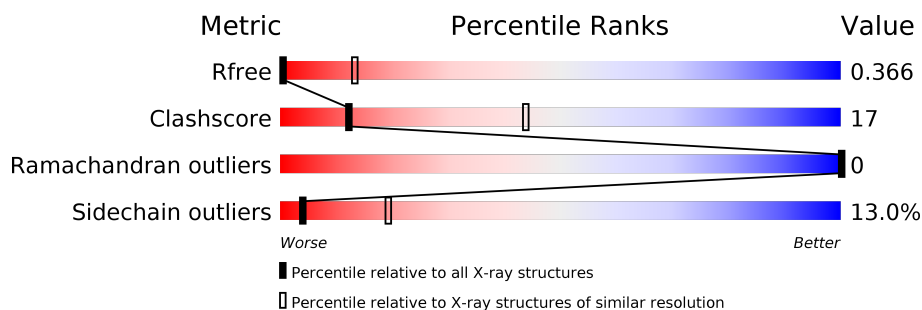
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 5.50 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}	130704	1019 (7.12-3.82)
Clashscore	141614	1010 (7.10-3.90)
Ramachandran outliers	138981	1014 (7.12-3.82)
Sidechain outliers	138945	1191 (7.20-3.80)

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 on 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	152	
1	B	152	
1	C	152	
1	D	152	
1	H	152	
1	I	152	
1	J	152	

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Mol	Chain	Length	Quality of chain
1	K	152	 A horizontal bar chart showing the quality of chain 1. The bar is divided into three segments: green (56%), yellow (30%), and orange (11%). The segments are labeled with their respective percentages: 56%, 30%, and 11%.

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 8086 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 Ion transport protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	136	Total	C	N	O	S	0	0	0
			1018	680	154	178	6			
1	B	136	Total	C	N	O	S	0	0	0
			1014	680	153	175	6			
1	C	136	Total	C	N	O	S	0	0	0
			1012	678	155	173	6			
1	D	136	Total	C	N	O	S	0	0	0
			999	671	153	169	6			
1	H	136	Total	C	N	O	S	0	0	0
			1018	680	154	178	6			
1	I	136	Total	C	N	O	S	0	0	0
			1014	680	153	175	6			
1	J	136	Total	C	N	O	S	0	0	0
			1012	678	155	173	6			
1	K	136	Total	C	N	O	S	0	0	0
			999	671	153	169	6			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	137	GLY	-	EXPRESSION TAG	UNP Q0ABW0
A	138	PRO	-	EXPRESSION TAG	UNP Q0ABW0
A	139	SER	-	EXPRESSION TAG	UNP Q0ABW0
A	140	SER	-	EXPRESSION TAG	UNP Q0ABW0
A	141	PRO	-	EXPRESSION TAG	UNP Q0ABW0
A	142	SER	-	EXPRESSION TAG	UNP Q0ABW0
B	137	GLY	-	EXPRESSION TAG	UNP Q0ABW0
B	138	PRO	-	EXPRESSION TAG	UNP Q0ABW0
B	139	SER	-	EXPRESSION TAG	UNP Q0ABW0
B	140	SER	-	EXPRESSION TAG	UNP Q0ABW0
B	141	PRO	-	EXPRESSION TAG	UNP Q0ABW0
B	142	SER	-	EXPRESSION TAG	UNP Q0ABW0
C	137	GLY	-	EXPRESSION TAG	UNP Q0ABW0

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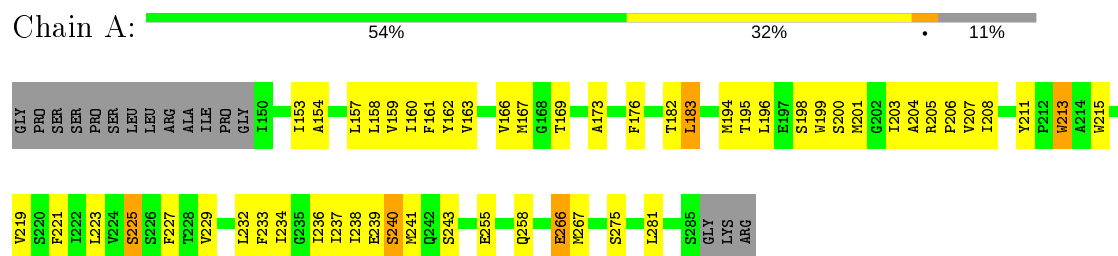
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Chain	Residue	Modelled	Actual	Comment	Reference
C	138	PRO	-	EXPRESSION TAG	UNP Q0ABW0
C	139	SER	-	EXPRESSION TAG	UNP Q0ABW0
C	140	SER	-	EXPRESSION TAG	UNP Q0ABW0
C	141	PRO	-	EXPRESSION TAG	UNP Q0ABW0
C	142	SER	-	EXPRESSION TAG	UNP Q0ABW0
D	137	GLY	-	EXPRESSION TAG	UNP Q0ABW0
D	138	PRO	-	EXPRESSION TAG	UNP Q0ABW0
D	139	SER	-	EXPRESSION TAG	UNP Q0ABW0
D	140	SER	-	EXPRESSION TAG	UNP Q0ABW0
D	141	PRO	-	EXPRESSION TAG	UNP Q0ABW0
D	142	SER	-	EXPRESSION TAG	UNP Q0ABW0
H	137	GLY	-	EXPRESSION TAG	UNP Q0ABW0
H	138	PRO	-	EXPRESSION TAG	UNP Q0ABW0
H	139	SER	-	EXPRESSION TAG	UNP Q0ABW0
H	140	SER	-	EXPRESSION TAG	UNP Q0ABW0
H	141	PRO	-	EXPRESSION TAG	UNP Q0ABW0
H	142	SER	-	EXPRESSION TAG	UNP Q0ABW0
I	137	GLY	-	EXPRESSION TAG	UNP Q0ABW0
I	138	PRO	-	EXPRESSION TAG	UNP Q0ABW0
I	139	SER	-	EXPRESSION TAG	UNP Q0ABW0
I	140	SER	-	EXPRESSION TAG	UNP Q0ABW0
I	141	PRO	-	EXPRESSION TAG	UNP Q0ABW0
I	142	SER	-	EXPRESSION TAG	UNP Q0ABW0
J	137	GLY	-	EXPRESSION TAG	UNP Q0ABW0
J	138	PRO	-	EXPRESSION TAG	UNP Q0ABW0
J	139	SER	-	EXPRESSION TAG	UNP Q0ABW0
J	140	SER	-	EXPRESSION TAG	UNP Q0ABW0
J	141	PRO	-	EXPRESSION TAG	UNP Q0ABW0
J	142	SER	-	EXPRESSION TAG	UNP Q0ABW0
K	137	GLY	-	EXPRESSION TAG	UNP Q0ABW0
K	138	PRO	-	EXPRESSION TAG	UNP Q0ABW0
K	139	SER	-	EXPRESSION TAG	UNP Q0ABW0
K	140	SER	-	EXPRESSION TAG	UNP Q0ABW0
K	141	PRO	-	EXPRESSION TAG	UNP Q0ABW0
K	142	SER	-	EXPRESSION TAG	UNP Q0ABW0

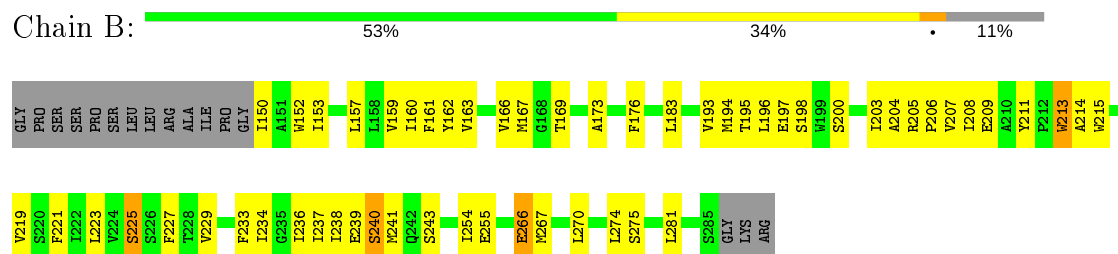
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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. 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. 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: Ion transport protein



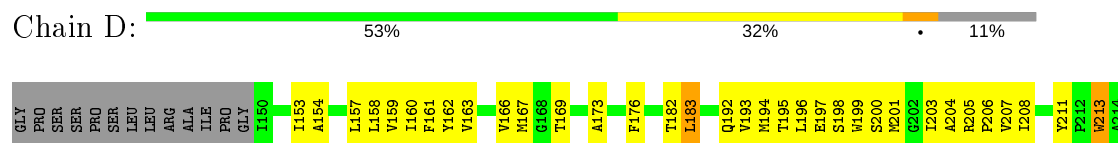
- Molecule 1: Ion transport protein



- Molecule 1: Ion transport protein



- Molecule 1: Ion transport protein





- Molecule 1: Ion transport protein

Chain H: 55% 31% 11%



- Molecule 1: Ion transport protein

Chain I: 56% 30% 11%



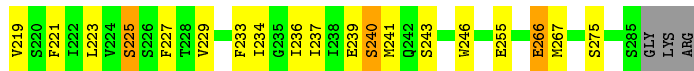
- Molecule 1: Ion transport protein

Chain J: 55% 29% 5% 11%



- Molecule 1: Ion transport protein

Chain K: 56% 30% 11%



4 Data and refinement statistics

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α , β , γ	181.14Å 181.14Å 94.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 5.50 128.08 – 5.50	Depositor EDS
% Data completeness (in resolution range)	96.9 (15.00-5.50) 96.6 (128.08-5.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 5.42Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.266 , 0.294 0.352 , 0.366	Depositor DCC
R_{free} test set	449 reflections (4.59%)	wwPDB-VP
Wilson B-factor (Å ²)	377.2	Xtriage
Anisotropy	0.060	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 452.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.30$, $\langle L^2 \rangle = 0.14$	Xtriage
Estimated twinning fraction	0.326 for h,-k,-l	Xtriage
Reported twinning fraction	0.598 for H, K, L 0.402 for -H, K, -L	Depositor
Outliers	0 of 9776 reflections	Xtriage
F_o, F_c correlation	0.76	EDS
Total number of atoms	8086	wwPDB-VP
Average B, all atoms (Å ²)	292.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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	0.53	0/1045	0.62	0/1431
1	B	0.54	0/1041	0.62	0/1426
1	C	0.54	0/1040	0.62	0/1425
1	D	0.53	0/1026	0.61	0/1407
1	H	0.53	0/1045	0.62	0/1431
1	I	0.51	0/1041	0.60	0/1426
1	J	0.52	0/1040	0.60	0/1425
1	K	0.53	0/1026	0.61	0/1407
All	All	0.53	0/8304	0.61	0/11378

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	1018	0	946	44	0
1	B	1014	0	947	44	0
1	C	1012	0	939	38	0
1	D	999	0	926	45	0
1	H	1018	0	946	41	0
1	I	1014	0	947	40	0
1	J	1012	0	939	44	0
1	K	999	0	926	42	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	8086	0	7516	265	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 17.

All (265) 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:234:ILE:HD11	1:D:237:ILE:HA	1.50	0.92
1:J:245:HIS:CE1	1:K:246:TRP:CB	2.54	0.91
1:H:234:ILE:HD11	1:K:237:ILE:HA	1.52	0.90
1:J:195:THR:HB	1:K:196:LEU:HD13	1.54	0.89
1:C:195:THR:HB	1:D:196:LEU:HD13	1.56	0.87
1:I:195:THR:HB	1:J:196:LEU:HD13	1.57	0.87
1:A:195:THR:HB	1:B:196:LEU:HD13	1.58	0.86
1:H:195:THR:HB	1:I:196:LEU:HD13	1.58	0.85
1:C:205:ARG:O	1:C:209:GLU:HG2	1.81	0.80
1:H:196:LEU:HD13	1:K:195:THR:HB	1.62	0.80
1:A:223:LEU:O	1:A:227:PHE:HB2	1.83	0.79
1:J:213:TRP:HD1	1:J:213:TRP:H	1.31	0.79
1:H:223:LEU:O	1:H:227:PHE:HB2	1.85	0.77
1:K:213:TRP:HD1	1:K:213:TRP:H	1.31	0.77
1:B:195:THR:HB	1:C:196:LEU:HD13	1.65	0.76
1:I:236:ILE:O	1:I:240:SER:HB2	1.86	0.76
1:B:236:ILE:O	1:B:240:SER:HB2	1.86	0.75
1:D:223:LEU:O	1:D:227:PHE:HB2	1.87	0.75
1:K:236:ILE:O	1:K:240:SER:HB2	1.86	0.74
1:K:223:LEU:O	1:K:227:PHE:HB2	1.87	0.74
1:A:213:TRP:H	1:A:213:TRP:HD1	1.34	0.74
1:A:205:ARG:HB2	1:A:206:PRO:HD3	1.68	0.74
1:D:213:TRP:H	1:D:213:TRP:HD1	1.34	0.74
1:A:196:LEU:HD13	1:D:195:THR:HB	1.69	0.74
1:I:223:LEU:O	1:I:227:PHE:HB2	1.88	0.74
1:B:213:TRP:HD1	1:B:213:TRP:H	1.35	0.73
1:J:223:LEU:O	1:J:227:PHE:HB2	1.88	0.73
1:B:223:LEU:O	1:B:227:PHE:HB2	1.88	0.73
1:C:223:LEU:O	1:C:227:PHE:HB2	1.89	0.73
1:C:213:TRP:H	1:C:213:TRP:HD1	1.37	0.73
1:D:236:ILE:O	1:D:240:SER:HB2	1.87	0.73
1:I:213:TRP:H	1:I:213:TRP:HD1	1.36	0.73
1:B:281:LEU:HD13	1:C:282:GLU:HG2	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:205:ARG:HB2	1:H:206:PRO:HD3	1.71	0.72
1:I:205:ARG:HB2	1:I:206:PRO:HD3	1.71	0.72
1:B:205:ARG:HB2	1:B:206:PRO:HD3	1.72	0.72
1:D:205:ARG:HB2	1:D:206:PRO:HD3	1.71	0.72
1:C:236:ILE:O	1:C:240:SER:HB2	1.89	0.71
1:J:245:HIS:HE1	1:K:246:TRP:CB	2.02	0.71
1:A:233:PHE:CE2	1:D:233:PHE:HZ	2.08	0.71
1:H:236:ILE:O	1:H:240:SER:HB2	1.91	0.71
1:H:213:TRP:H	1:H:213:TRP:HD1	1.35	0.71
1:A:236:ILE:O	1:A:240:SER:HB2	1.91	0.71
1:J:236:ILE:O	1:J:240:SER:HB2	1.91	0.71
1:J:205:ARG:O	1:J:209:GLU:HG2	1.90	0.70
1:J:205:ARG:HB2	1:J:206:PRO:HD3	1.73	0.70
1:D:200:SER:HA	1:D:204:ALA:HB3	1.73	0.70
1:C:205:ARG:HB2	1:C:206:PRO:HD3	1.74	0.70
1:K:205:ARG:HB2	1:K:206:PRO:HD3	1.74	0.70
1:C:245:HIS:NE2	1:D:246:TRP:CB	2.55	0.69
1:B:200:SER:HA	1:B:204:ALA:HB3	1.74	0.69
1:K:200:SER:HA	1:K:204:ALA:HB3	1.75	0.69
1:I:200:SER:HA	1:I:204:ALA:HB3	1.75	0.68
1:I:205:ARG:O	1:I:209:GLU:HG3	1.94	0.68
1:J:200:SER:HA	1:J:204:ALA:HB3	1.75	0.67
1:A:200:SER:HA	1:A:204:ALA:HB3	1.76	0.67
1:C:200:SER:HA	1:C:204:ALA:HB3	1.76	0.66
1:H:200:SER:HA	1:H:204:ALA:HB3	1.77	0.66
1:A:200:SER:HB3	1:D:192:GLN:HG3	1.79	0.65
1:H:233:PHE:CE2	1:K:233:PHE:HZ	2.14	0.64
1:C:211:TYR:HB3	1:C:213:TRP:CD1	2.32	0.64
1:H:176:PHE:CZ	1:H:207:VAL:HA	2.33	0.64
1:H:211:TYR:HB3	1:H:213:TRP:CD1	2.32	0.64
1:I:211:TYR:HB3	1:I:213:TRP:CD1	2.33	0.64
1:A:234:ILE:HD11	1:D:237:ILE:CA	2.26	0.64
1:C:176:PHE:CZ	1:C:207:VAL:HA	2.34	0.63
1:A:176:PHE:CZ	1:A:207:VAL:HA	2.33	0.63
1:A:211:TYR:HB3	1:A:213:TRP:CD1	2.34	0.63
1:A:215:TRP:O	1:A:219:VAL:HG23	1.98	0.63
1:B:194:MET:HG3	1:B:225:SER:OG	1.98	0.62
1:D:194:MET:HG3	1:D:225:SER:OG	1.99	0.62
1:K:176:PHE:CZ	1:K:207:VAL:HA	2.35	0.61
1:J:211:TYR:HB3	1:J:213:TRP:CD1	2.35	0.61
1:H:234:ILE:HD11	1:K:237:ILE:CA	2.28	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:238:ILE:HG12	1:K:241:MET:HG2	1.81	0.61
1:D:211:TYR:HB3	1:D:213:TRP:CD1	2.36	0.61
1:D:176:PHE:CZ	1:D:207:VAL:HA	2.35	0.60
1:H:200:SER:HB3	1:K:192:GLN:HG3	1.82	0.60
1:A:194:MET:HG3	1:A:225:SER:OG	2.01	0.60
1:J:245:HIS:NE2	1:K:246:TRP:CB	2.63	0.60
1:I:281:LEU:HD13	1:J:282:GLU:HG2	1.83	0.60
1:H:215:TRP:O	1:H:219:VAL:HG23	2.01	0.60
1:B:176:PHE:CZ	1:B:207:VAL:HA	2.36	0.60
1:B:266:GLU:HG3	1:B:267:MET:N	2.16	0.59
1:B:225:SER:O	1:B:229:VAL:HG23	2.02	0.59
1:K:211:TYR:HB3	1:K:213:TRP:CD1	2.39	0.58
1:K:194:MET:HG3	1:K:225:SER:OG	2.04	0.58
1:B:211:TYR:HB3	1:B:213:TRP:CD1	2.38	0.58
1:J:194:MET:HG3	1:J:225:SER:OG	2.04	0.58
1:C:194:MET:HG3	1:C:225:SER:OG	2.05	0.57
1:H:194:MET:HG3	1:H:225:SER:OG	2.03	0.57
1:A:236:ILE:HG22	1:B:234:ILE:HD13	1.87	0.57
1:A:198:SER:HA	1:D:197:GLU:OE1	2.05	0.57
1:C:225:SER:O	1:C:229:VAL:HG23	2.04	0.57
1:I:215:TRP:O	1:I:219:VAL:HG23	2.04	0.57
1:B:215:TRP:O	1:B:219:VAL:HG23	2.06	0.55
1:K:215:TRP:O	1:K:219:VAL:HG23	2.06	0.55
1:A:238:ILE:HG12	1:D:241:MET:HG2	1.88	0.55
1:A:266:GLU:HG3	1:A:267:MET:N	2.21	0.55
1:D:266:GLU:HG3	1:D:267:MET:N	2.21	0.55
1:I:194:MET:HG3	1:I:225:SER:OG	2.06	0.55
1:J:203:ILE:O	1:J:207:VAL:HG23	2.06	0.55
1:I:270:LEU:HD21	1:J:272:ARG:CG	2.37	0.55
1:C:237:ILE:HA	1:D:234:ILE:HD11	1.89	0.55
1:J:176:PHE:CZ	1:J:207:VAL:HA	2.42	0.55
1:C:215:TRP:O	1:C:219:VAL:HG23	2.07	0.55
1:I:203:ILE:O	1:I:207:VAL:HG23	2.08	0.54
1:B:197:GLU:OE2	1:C:200:SER:HB3	2.07	0.54
1:C:266:GLU:HG3	1:C:267:MET:N	2.20	0.54
1:J:197:GLU:OE2	1:K:200:SER:HB3	2.07	0.54
1:I:270:LEU:HD21	1:J:272:ARG:HG2	1.90	0.54
1:J:225:SER:O	1:J:229:VAL:HG23	2.08	0.54
1:B:281:LEU:CD1	1:C:282:GLU:HG2	2.36	0.54
1:K:225:SER:O	1:K:229:VAL:HG23	2.07	0.54
1:I:176:PHE:CZ	1:I:207:VAL:HA	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:ARG:O	1:B:209:GLU:HG3	2.07	0.54
1:H:237:ILE:HA	1:I:234:ILE:HD11	1.89	0.54
1:K:266:GLU:HG3	1:K:267:MET:N	2.22	0.53
1:C:197:GLU:HG3	1:D:199:TRP:CD1	2.44	0.52
1:H:162:TYR:O	1:H:166:VAL:HG23	2.09	0.52
1:J:213:TRP:N	1:J:213:TRP:CD1	2.76	0.52
1:H:225:SER:O	1:H:229:VAL:HG23	2.09	0.52
1:I:225:SER:O	1:I:229:VAL:HG23	2.09	0.52
1:J:215:TRP:O	1:J:219:VAL:HG23	2.09	0.52
1:A:237:ILE:O	1:A:241:MET:HB2	2.09	0.52
1:I:266:GLU:HG3	1:I:267:MET:N	2.25	0.52
1:I:237:ILE:HA	1:J:234:ILE:HD11	1.92	0.52
1:C:237:ILE:O	1:C:241:MET:HB2	2.10	0.52
1:D:215:TRP:O	1:D:219:VAL:HG23	2.10	0.52
1:I:197:GLU:HG3	1:J:199:TRP:CD1	2.45	0.52
1:A:166:VAL:HG22	1:A:183:LEU:HD21	1.92	0.51
1:B:197:GLU:HG3	1:C:199:TRP:CD1	2.46	0.51
1:I:197:GLU:OE2	1:J:200:SER:HB3	2.10	0.51
1:H:198:SER:HA	1:K:197:GLU:OE1	2.10	0.51
1:K:162:TYR:O	1:K:166:VAL:HG23	2.11	0.51
1:B:237:ILE:HA	1:C:234:ILE:HD11	1.93	0.50
1:B:240:SER:HB3	1:C:238:ILE:HD11	1.92	0.50
1:A:233:PHE:CE2	1:D:233:PHE:CZ	2.94	0.50
1:B:221:PHE:O	1:B:225:SER:HB3	2.12	0.50
1:A:225:SER:O	1:A:229:VAL:HG23	2.11	0.50
1:A:159:VAL:O	1:A:163:VAL:HG23	2.12	0.50
1:J:237:ILE:HA	1:K:234:ILE:HD11	1.93	0.50
1:D:225:SER:O	1:D:229:VAL:HG23	2.11	0.50
1:C:197:GLU:OE2	1:D:200:SER:HB3	2.11	0.50
1:D:208:ILE:HA	1:D:211:TYR:O	2.12	0.49
1:D:213:TRP:N	1:D:213:TRP:CD1	2.79	0.49
1:J:197:GLU:HG3	1:K:199:TRP:CD1	2.47	0.49
1:H:237:ILE:O	1:H:241:MET:HB2	2.11	0.49
1:D:203:ILE:O	1:D:207:VAL:HG23	2.12	0.49
1:A:213:TRP:CD1	1:A:213:TRP:N	2.79	0.49
1:J:266:GLU:HG3	1:J:267:MET:N	2.27	0.49
1:I:213:TRP:N	1:I:213:TRP:CD1	2.80	0.49
1:B:162:TYR:O	1:B:166:VAL:HG23	2.13	0.48
1:C:198:SER:HB2	1:D:201:MET:HG3	1.94	0.48
1:H:233:PHE:CE2	1:K:233:PHE:CZ	3.00	0.48
1:C:203:ILE:O	1:C:207:VAL:HG23	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:166:VAL:HG22	1:D:183:LEU:HD21	1.95	0.48
1:A:201:MET:HG2	1:D:192:GLN:NE2	2.27	0.48
1:C:213:TRP:N	1:C:213:TRP:CD1	2.81	0.48
1:A:157:LEU:HA	1:A:160:ILE:HD12	1.94	0.48
1:A:237:ILE:HA	1:B:234:ILE:HD11	1.96	0.48
1:I:162:TYR:O	1:I:166:VAL:HG23	2.13	0.48
1:D:193:VAL:O	1:D:196:LEU:HD23	2.13	0.48
1:H:239:GLU:O	1:H:243:SER:HB3	2.14	0.48
1:K:203:ILE:O	1:K:207:VAL:HG23	2.14	0.48
1:J:198:SER:HB2	1:K:201:MET:HG3	1.96	0.48
1:H:157:LEU:HA	1:H:160:ILE:HD12	1.97	0.47
1:D:157:LEU:HA	1:D:160:ILE:HD12	1.95	0.47
1:C:197:GLU:OE1	1:D:198:SER:HA	2.14	0.47
1:K:159:VAL:O	1:K:163:VAL:HG23	2.15	0.47
1:A:162:TYR:O	1:A:166:VAL:HG23	2.14	0.47
1:I:169:THR:O	1:I:173:ALA:HB2	2.14	0.47
1:J:162:TYR:O	1:J:166:VAL:HG23	2.15	0.47
1:I:240:SER:HB3	1:J:238:ILE:HD11	1.97	0.47
1:H:201:MET:HG3	1:K:198:SER:HB2	1.95	0.47
1:J:237:ILE:O	1:J:241:MET:HB2	2.14	0.47
1:H:169:THR:O	1:H:173:ALA:HB2	2.14	0.47
1:A:219:VAL:O	1:A:223:LEU:HG	2.15	0.47
1:B:208:ILE:HA	1:B:211:TYR:O	2.15	0.47
1:D:162:TYR:O	1:D:166:VAL:HG23	2.15	0.46
1:H:266:GLU:HG3	1:H:267:MET:N	2.30	0.46
1:B:166:VAL:HG22	1:B:183:LEU:HD21	1.98	0.46
1:C:169:THR:O	1:C:173:ALA:HB2	2.16	0.46
1:H:213:TRP:CD1	1:H:213:TRP:N	2.79	0.46
1:I:208:ILE:HD11	1:I:215:TRP:HA	1.97	0.46
1:A:203:ILE:O	1:A:207:VAL:HG23	2.16	0.46
1:A:258:GLN:O	1:A:258:GLN:HG2	2.16	0.46
1:I:159:VAL:O	1:I:163:VAL:HG23	2.16	0.46
1:C:166:VAL:HG22	1:C:183:LEU:HD21	1.98	0.46
1:H:240:SER:HB3	1:I:238:ILE:HD11	1.97	0.46
1:B:270:LEU:HD21	1:C:272:ARG:CG	2.46	0.45
1:I:208:ILE:HA	1:I:211:TYR:O	2.16	0.45
1:K:193:VAL:O	1:K:196:LEU:HD23	2.17	0.45
1:I:198:SER:HB2	1:J:201:MET:HG3	1.97	0.45
1:J:169:THR:O	1:J:173:ALA:HB2	2.17	0.45
1:I:263:GLU:OE2	1:J:261:HIS:CE1	2.69	0.45
1:J:208:ILE:HA	1:J:211:TYR:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:SER:HB2	1:C:201:MET:HG3	1.98	0.44
1:I:166:VAL:HG22	1:I:183:LEU:HD21	2.00	0.44
1:J:221:PHE:O	1:J:225:SER:HB3	2.17	0.44
1:K:213:TRP:CD1	1:K:213:TRP:N	2.77	0.44
1:D:154:ALA:O	1:D:158:LEU:HD12	2.18	0.44
1:H:203:ILE:O	1:H:207:VAL:HG23	2.16	0.44
1:H:236:ILE:HG22	1:I:234:ILE:HD13	1.99	0.44
1:D:159:VAL:O	1:D:163:VAL:HG23	2.17	0.44
1:H:233:PHE:HZ	1:I:233:PHE:CE2	2.35	0.44
1:D:239:GLU:O	1:D:243:SER:HB3	2.18	0.44
1:H:221:PHE:O	1:H:225:SER:HB3	2.18	0.44
1:J:166:VAL:HG22	1:J:183:LEU:HD21	1.98	0.44
1:H:208:ILE:HA	1:H:211:TYR:O	2.17	0.44
1:K:237:ILE:O	1:K:241:MET:HB2	2.17	0.44
1:D:169:THR:O	1:D:173:ALA:HB2	2.18	0.44
1:A:199:TRP:H	1:D:197:GLU:HG3	1.83	0.43
1:H:197:GLU:OE1	1:I:198:SER:HA	2.18	0.43
1:B:274:LEU:CD1	1:C:275:SER:HA	2.48	0.43
1:I:274:LEU:CD1	1:J:275:SER:HA	2.47	0.43
1:J:274:LEU:HA	1:J:274:LEU:HD12	1.90	0.43
1:A:169:THR:O	1:A:173:ALA:HB2	2.18	0.43
1:C:162:TYR:O	1:C:166:VAL:HG23	2.17	0.43
1:A:208:ILE:HD11	1:A:215:TRP:HA	2.01	0.43
1:B:157:LEU:HA	1:B:160:ILE:HD12	2.00	0.43
1:B:208:ILE:HD11	1:B:215:TRP:HA	2.00	0.43
1:D:221:PHE:O	1:D:225:SER:HB3	2.18	0.43
1:K:239:GLU:O	1:K:243:SER:HB3	2.18	0.43
1:B:203:ILE:O	1:B:207:VAL:HG23	2.18	0.43
1:B:274:LEU:HD12	1:B:274:LEU:HA	1.91	0.43
1:C:157:LEU:HA	1:C:160:ILE:HD12	2.00	0.43
1:D:237:ILE:O	1:D:241:MET:HB2	2.19	0.43
1:J:197:GLU:OE1	1:K:198:SER:HA	2.18	0.42
1:I:237:ILE:O	1:I:241:MET:HB2	2.19	0.42
1:A:233:PHE:CD2	1:D:233:PHE:HZ	2.37	0.42
1:K:157:LEU:HA	1:K:160:ILE:HD12	2.00	0.42
1:H:199:TRP:H	1:K:197:GLU:HG3	1.84	0.42
1:B:213:TRP:CD1	1:B:213:TRP:N	2.80	0.42
1:C:159:VAL:O	1:C:163:VAL:HG23	2.19	0.42
1:H:201:MET:HG2	1:K:192:GLN:NE2	2.34	0.42
1:J:239:GLU:O	1:J:243:SER:HB3	2.20	0.42
1:J:157:LEU:HA	1:J:160:ILE:HD12	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:PHE:HZ	1:B:233:PHE:CE2	2.37	0.42
1:B:237:ILE:O	1:B:241:MET:HB2	2.19	0.42
1:H:159:VAL:O	1:H:163:VAL:HG23	2.20	0.42
1:A:239:GLU:O	1:A:243:SER:HB3	2.20	0.41
1:B:270:LEU:HD21	1:C:272:ARG:HG3	2.01	0.41
1:A:201:MET:HG3	1:D:198:SER:HB2	2.02	0.41
1:H:175:SER:HG	1:H:211:TYR:HH	1.65	0.41
1:B:239:GLU:O	1:B:243:SER:HB3	2.21	0.41
1:K:169:THR:O	1:K:173:ALA:HB2	2.19	0.41
1:A:208:ILE:HA	1:A:211:TYR:O	2.21	0.41
1:B:159:VAL:O	1:B:163:VAL:HG23	2.19	0.41
1:K:221:PHE:O	1:K:225:SER:HB3	2.21	0.41
1:H:166:VAL:HG22	1:H:183:LEU:HD21	2.01	0.41
1:A:221:PHE:O	1:A:225:SER:HB3	2.20	0.41
1:K:166:VAL:HG22	1:K:183:LEU:HD21	2.03	0.41
1:B:169:THR:O	1:B:173:ALA:HB2	2.19	0.41
1:B:207:VAL:HG12	1:B:214:ALA:CB	2.51	0.41
1:A:240:SER:HB3	1:B:238:ILE:HD11	2.02	0.41
1:B:150:ILE:N	1:B:152:TRP:HD1	2.19	0.41
1:I:239:GLU:O	1:I:243:SER:HB3	2.21	0.41
1:H:234:ILE:HA	1:H:237:ILE:HD12	2.04	0.40
1:I:270:LEU:HD21	1:J:272:ARG:HG3	2.03	0.40
1:A:232:LEU:O	1:A:236:ILE:HG13	2.20	0.40
1:B:193:VAL:O	1:B:196:LEU:HD23	2.21	0.40
1:D:208:ILE:HD11	1:D:215:TRP:HA	2.02	0.40
1:A:154:ALA:O	1:A:158:LEU:HD12	2.21	0.40
1:D:232:LEU:O	1:D:236:ILE:HG13	2.22	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	134/152 (88%)	129 (96%)	5 (4%)	0	100	100
1	B	134/152 (88%)	130 (97%)	4 (3%)	0	100	100
1	C	134/152 (88%)	128 (96%)	6 (4%)	0	100	100
1	D	134/152 (88%)	129 (96%)	5 (4%)	0	100	100
1	H	134/152 (88%)	129 (96%)	5 (4%)	0	100	100
1	I	134/152 (88%)	130 (97%)	4 (3%)	0	100	100
1	J	134/152 (88%)	128 (96%)	6 (4%)	0	100	100
1	K	134/152 (88%)	129 (96%)	5 (4%)	0	100	100
All	All	1072/1216 (88%)	1032 (96%)	40 (4%)	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	95/131 (72%)	83 (87%)	12 (13%)	4	20
1	B	94/131 (72%)	84 (89%)	10 (11%)	6	25
1	C	93/131 (71%)	81 (87%)	12 (13%)	4	19
1	D	90/131 (69%)	78 (87%)	12 (13%)	4	19
1	H	95/131 (72%)	81 (85%)	14 (15%)	3	16
1	I	94/131 (72%)	81 (86%)	13 (14%)	3	18
1	J	93/131 (71%)	81 (87%)	12 (13%)	4	19
1	K	90/131 (69%)	78 (87%)	12 (13%)	4	19
All	All	744/1048 (71%)	647 (87%)	97 (13%)	4	19

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	ILE
1	A	161	PHE

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Mol	Chain	Res	Type
1	A	167	MET
1	A	182	THR
1	A	183	LEU
1	A	213	TRP
1	A	225	SER
1	A	240	SER
1	A	255	GLU
1	A	266	GLU
1	A	275	SER
1	A	281	LEU
1	B	153	ILE
1	B	161	PHE
1	B	167	MET
1	B	213	TRP
1	B	225	SER
1	B	240	SER
1	B	254	ILE
1	B	255	GLU
1	B	266	GLU
1	B	275	SER
1	C	153	ILE
1	C	161	PHE
1	C	166	VAL
1	C	167	MET
1	C	182	THR
1	C	213	TRP
1	C	225	SER
1	C	240	SER
1	C	250	ASP
1	C	255	GLU
1	C	266	GLU
1	C	275	SER
1	D	153	ILE
1	D	161	PHE
1	D	167	MET
1	D	182	THR
1	D	183	LEU
1	D	213	TRP
1	D	215	TRP
1	D	225	SER
1	D	240	SER
1	D	255	GLU

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Mol	Chain	Res	Type
1	D	266	GLU
1	D	275	SER
1	H	153	ILE
1	H	161	PHE
1	H	167	MET
1	H	182	THR
1	H	183	LEU
1	H	213	TRP
1	H	215	TRP
1	H	225	SER
1	H	240	SER
1	H	255	GLU
1	H	266	GLU
1	H	275	SER
1	H	279	ASP
1	H	282	GLU
1	I	153	ILE
1	I	161	PHE
1	I	167	MET
1	I	182	THR
1	I	183	LEU
1	I	213	TRP
1	I	215	TRP
1	I	225	SER
1	I	240	SER
1	I	250	ASP
1	I	255	GLU
1	I	266	GLU
1	I	275	SER
1	J	153	ILE
1	J	161	PHE
1	J	167	MET
1	J	183	LEU
1	J	209	GLU
1	J	213	TRP
1	J	215	TRP
1	J	225	SER
1	J	240	SER
1	J	255	GLU
1	J	266	GLU
1	J	275	SER
1	K	153	ILE

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Mol	Chain	Res	Type
1	K	161	PHE
1	K	167	MET
1	K	182	THR
1	K	183	LEU
1	K	213	TRP
1	K	215	TRP
1	K	225	SER
1	K	240	SER
1	K	255	GLU
1	K	266	GLU
1	K	275	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	J	245	HIS

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 carbohydrates 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

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.