



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 12:07 am BST

PDB ID : 3VYI  
Title : Crystal Structure of a trimeric coiled-coil (I/I-type) assembly domain from the voltage-gated proton channel mutant  
Authors : Fujiwara, Y.; Takeshita, K.; Nakagawa, A.  
Deposited on : 2012-09-25  
Resolution : 2.31 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](mailto: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.

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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

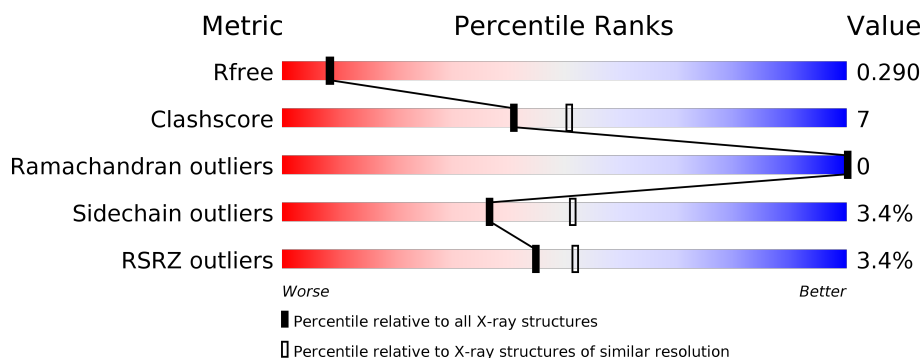
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.31 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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  $\geq 3$ , 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	51	<div> <div>75%</div> <div>16%</div> <div>10%</div> </div>
1	B	51	<div>2%</div> <div>71%</div> <div>20%</div> <div>10%</div>

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Mol	Chain	Length	Quality of chain
1	G	51	
1	H	51	
1	I	51	
1	J	51	
1	K	51	
1	L	51	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4529 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 Voltage-gated hydrogen channel 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	46	Total	C	N	O	0	0	0
			384	241	71	72			
1	B	46	Total	C	N	O	0	0	0
			384	241	71	72			
1	D	48	Total	C	N	O	0	0	0
			392	245	73	74			
1	C	48	Total	C	N	O	0	0	0
			396	247	73	76			
1	E	46	Total	C	N	O	0	0	0
			384	241	71	72			
1	F	49	Total	C	N	O	0	0	0
			400	249	74	77			
1	H	40	Total	C	N	O	0	0	0
			341	215	63	63			
1	G	40	Total	C	N	O	0	0	0
			338	212	62	64			
1	I	40	Total	C	N	O	0	0	0
			338	212	62	64			
1	J	42	Total	C	N	O	0	0	0
			356	223	66	67			
1	K	39	Total	C	N	O	0	0	0
			332	210	62	60			
1	L	38	Total	C	N	O	0	0	0
			324	203	60	61			

There are 132 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
A	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
A	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
A	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
A	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
A	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
A	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
A	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
A	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
A	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
B	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
B	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
B	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
B	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
B	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
B	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
B	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
B	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
B	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
B	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
B	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
D	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
D	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
D	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
D	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
D	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
D	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
D	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
D	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
D	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
D	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
D	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
C	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
C	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
C	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
C	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
C	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
C	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
C	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
C	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
C	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
C	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
C	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
E	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
E	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
E	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8

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Chain	Residue	Modelled	Actual	Comment	Reference
E	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
E	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
E	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
E	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
E	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
E	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
E	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
E	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
F	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
F	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
F	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
F	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
F	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
F	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
F	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
F	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
F	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
F	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
F	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
H	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
H	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
H	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
H	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
H	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
H	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
H	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
H	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
H	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
H	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
H	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
G	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
G	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
G	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
G	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
G	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
G	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
G	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
G	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
G	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
G	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
G	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
I	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8

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Chain	Residue	Modelled	Actual	Comment	Reference
I	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
I	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
I	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
I	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
I	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
I	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
I	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
I	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
I	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
I	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
J	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
J	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
J	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
J	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
J	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
J	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
J	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
J	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
J	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
J	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
J	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
K	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
K	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
K	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
K	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
K	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
K	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
K	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
K	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
K	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
K	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
K	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
L	219	GLY	-	EXPRESSION TAG	UNP Q3U2S8
L	227	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
L	230	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
L	231	ILE	ASN	ENGINEERED MUTATION	UNP Q3U2S8
L	232	ASN	ILE	ENGINEERED MUTATION	UNP Q3U2S8
L	234	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
L	241	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8
L	245	ILE	CYS	ENGINEERED MUTATION	UNP Q3U2S8
L	248	ILE	LYS	ENGINEERED MUTATION	UNP Q3U2S8
L	255	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8

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Chain	Residue	Modelled	Actual	Comment	Reference
L	259	ILE	LEU	ENGINEERED MUTATION	UNP Q3U2S8

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	14	Total O 14 14	0	0
2	B	22	Total O 22 22	0	0
2	D	12	Total O 12 12	0	0
2	C	15	Total O 15 15	0	0
2	E	26	Total O 26 26	0	0
2	F	16	Total O 16 16	0	0
2	H	10	Total O 10 10	0	0
2	G	13	Total O 13 13	0	0
2	I	11	Total O 11 11	0	0
2	J	3	Total O 3 3	0	0
2	K	8	Total O 8 8	0	0
2	L	10	Total O 10 10	0	0



### 3 Residue-property plots [i](#)

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 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: Voltage-gated hydrogen channel 1

Chain A: 



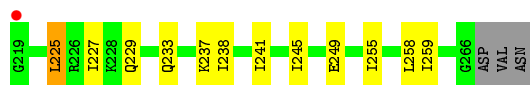
- Molecule 1: Voltage-gated hydrogen channel 1

Chain B: 



- Molecule 1: Voltage-gated hydrogen channel 1

Chain D: 



- Molecule 1: Voltage-gated hydrogen channel 1

Chain C: 

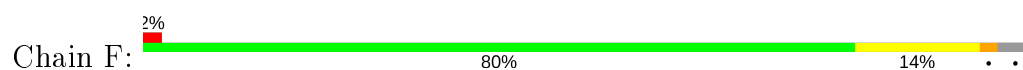


- Molecule 1: Voltage-gated hydrogen channel 1

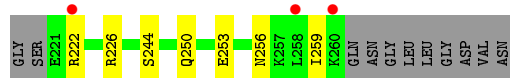
Chain E: 



- Molecule 1: Voltage-gated hydrogen channel 1



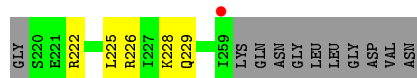
- Molecule 1: Voltage-gated hydrogen channel 1



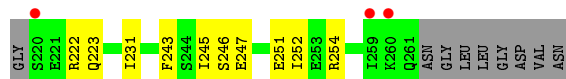
- Molecule 1: Voltage-gated hydrogen channel 1



- Molecule 1: Voltage-gated hydrogen channel 1



- Molecule 1: Voltage-gated hydrogen channel 1



- Molecule 1: Voltage-gated hydrogen channel 1



- Molecule 1: Voltage-gated hydrogen channel 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.27Å 84.01Å 88.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.11 – 2.31 39.11 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.7 (39.11-2.31) 98.4 (39.11-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.58 (at 2.31Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, $R_{free}$	0.232 , 0.295 0.230 , 0.290	Depositor DCC
$R_{free}$ test set	1398 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.6	Xtriage
Anisotropy	0.227	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 47.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4529	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.47 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.9742e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.49	0/385	0.57	0/513
1	B	0.47	0/385	0.52	0/513
1	C	0.43	0/397	0.55	0/529
1	D	0.48	0/393	0.56	0/523
1	E	0.48	0/385	0.57	0/513
1	F	0.44	0/401	0.52	0/534
1	G	0.51	0/339	0.60	0/452
1	H	0.41	0/342	0.47	0/455
1	I	0.44	0/339	0.47	0/452
1	J	0.47	0/357	0.52	0/475
1	K	0.41	0/333	0.45	0/443
1	L	0.45	0/325	0.53	0/433
All	All	0.46	0/4381	0.53	0/5835

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	384	0	409	6	0
1	B	384	0	409	12	0
1	C	396	0	416	6	0
1	D	392	0	415	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	384	0	409	13	0
1	F	400	0	419	6	0
1	G	338	0	357	5	0
1	H	341	0	365	3	0
1	I	338	0	357	2	0
1	J	356	0	378	9	0
1	K	332	0	359	5	0
1	L	324	0	341	19	0
2	A	14	0	0	0	0
2	B	22	0	0	0	0
2	C	15	0	0	0	0
2	D	12	0	0	0	0
2	E	26	0	0	0	0
2	F	16	0	0	0	0
2	G	13	0	0	0	0
2	H	10	0	0	0	0
2	I	11	0	0	1	0
2	J	3	0	0	0	0
2	K	8	0	0	1	0
2	L	10	0	0	0	0
All	All	4529	0	4634	64	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 7.

All (64) 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:B:229:GLN:HG3	1:L:233:GLN:HE21	1.39	0.88
1:B:229:GLN:HG3	1:L:233:GLN:NE2	2.06	0.70
1:B:229:GLN:HG2	1:E:240:HIS:CE1	2.30	0.67
1:A:256:ASN:OD1	1:B:254:ARG:NH2	2.29	0.60
1:E:247:GLU:HG2	1:L:225:LEU:HD23	1.83	0.60
1:E:255:ILE:O	1:E:259:ILE:HG12	2.05	0.57
1:F:266:GLY:O	1:F:267:ASP:HB2	2.04	0.56
1:I:225:LEU:HD23	1:I:228:LYS:HE2	1.87	0.56
1:I:226:ARG:NH1	2:I:305:HOH:O	2.38	0.56
1:F:246:SER:O	1:F:250:GLN:HG2	2.05	0.56
1:L:224:ILE:HG22	1:L:228:LYS:HE2	1.88	0.55
1:E:251:GLU:HG2	1:L:222:ARG:HH22	1.72	0.54
1:E:243:PHE:CG	1:L:229:GLN:HG3	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:228:LYS:NZ	1:L:250:GLN:OE1	2.43	0.52
1:D:225:LEU:O	1:D:229:GLN:HG3	2.11	0.51
1:E:243:PHE:CD2	1:L:229:GLN:HG3	2.46	0.50
1:D:258:LEU:HD22	1:F:259:ILE:HD13	1.94	0.49
1:B:243:PHE:CD2	1:G:229:GLN:HG3	2.47	0.49
1:L:251:GLU:OE2	1:L:254:ARG:NH2	2.46	0.49
1:L:222:ARG:HD2	1:L:226:ARG:CZ	2.42	0.49
1:A:227:ILE:O	1:A:231:ILE:HG12	2.12	0.49
1:G:250:GLN:HA	1:G:253:GLU:HG2	1.95	0.48
1:J:231:ILE:HG23	1:K:234:ILE:HD11	1.94	0.48
1:K:245:ILE:O	1:K:249:GLU:HG3	2.13	0.48
1:A:265:LEU:HD13	1:B:258:LEU:HD22	1.93	0.48
1:J:222:ARG:HD2	1:J:222:ARG:HA	1.58	0.48
1:H:256:ASN:O	1:H:259:ILE:HG13	2.14	0.48
1:K:252:ILE:HG21	1:L:251:GLU:HG2	1.96	0.47
1:B:265:LEU:HB2	1:C:262:ASN:HD21	1.80	0.47
1:D:237:LYS:HB2	1:D:237:LYS:HE3	1.65	0.46
1:D:255:ILE:O	1:D:259:ILE:HG12	2.15	0.46
1:B:226:ARG:NH2	1:D:249:GLU:OE2	2.41	0.46
1:L:230:ASN:O	1:L:234:ILE:HG12	2.16	0.45
1:L:223:GLN:CD	1:L:223:GLN:H	2.19	0.45
1:K:227:ILE:O	1:K:231:ILE:HG12	2.16	0.45
1:H:222:ARG:O	1:H:226:ARG:HG3	2.17	0.45
1:E:239:GLN:NE2	1:L:232:ASN:OD1	2.49	0.44
1:B:238:ILE:HD11	1:C:234:ILE:HG23	2.00	0.44
1:E:234:ILE:O	1:E:238:ILE:HG12	2.18	0.44
1:C:226:ARG:HD3	2:K:303:HOH:O	2.17	0.44
1:D:258:LEU:HD12	1:F:267:ASP:N	2.33	0.44
1:C:257:LYS:HB2	1:C:257:LYS:HE3	1.84	0.44
1:J:223:GLN:HG3	1:L:224:ILE:HG21	2.00	0.44
1:D:241:ILE:O	1:D:245:ILE:HG12	2.17	0.43
1:J:251:GLU:HG2	1:L:252:ILE:HG21	1.99	0.43
1:J:243:PHE:O	1:J:247:GLU:HG2	2.18	0.43
1:G:241:ILE:O	1:G:245:ILE:HG12	2.19	0.43
1:J:251:GLU:OE1	1:J:254:ARG:HD3	2.19	0.43
1:L:221:GLU:HG2	1:L:224:ILE:HG13	2.01	0.42
1:A:255:ILE:O	1:A:259:ILE:HG12	2.20	0.42
1:D:227:ILE:HD11	1:F:224:ILE:HG23	2.02	0.42
1:D:238:ILE:HD11	1:E:238:ILE:HD11	2.01	0.42
1:A:245:ILE:HD11	1:B:241:ILE:HG23	2.02	0.41
1:J:251:GLU:HA	1:J:254:ARG:HG2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:252:ILE:HG23	1:K:255:ILE:HD11	2.02	0.41
1:B:243:PHE:CG	1:G:229:GLN:HG3	2.54	0.41
1:F:240:HIS:NE2	1:H:250:GLN:OE1	2.52	0.41
1:A:249:GLU:OE2	1:E:226:ARG:NH2	2.50	0.41
1:C:224:ILE:HA	1:C:224:ILE:HD13	1.87	0.41
1:J:245:ILE:HA	1:J:245:ILE:HD13	1.89	0.41
1:E:243:PHE:CE1	1:L:225:LEU:HG	2.55	0.41
1:B:229:GLN:HG2	1:E:240:HIS:NE2	2.36	0.41
1:C:222:ARG:HG3	1:C:226:ARG:NH2	2.37	0.41
1:E:243:PHE:CD1	1:L:229:GLN:HG3	2.55	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	44/51 (86%)	44 (100%)	0	0	100	100
1	B	44/51 (86%)	44 (100%)	0	0	100	100
1	C	46/51 (90%)	46 (100%)	0	0	100	100
1	D	46/51 (90%)	45 (98%)	1 (2%)	0	100	100
1	E	44/51 (86%)	44 (100%)	0	0	100	100
1	F	47/51 (92%)	47 (100%)	0	0	100	100
1	G	38/51 (74%)	38 (100%)	0	0	100	100
1	H	38/51 (74%)	38 (100%)	0	0	100	100
1	I	38/51 (74%)	37 (97%)	1 (3%)	0	100	100
1	J	40/51 (78%)	39 (98%)	1 (2%)	0	100	100
1	K	37/51 (72%)	36 (97%)	1 (3%)	0	100	100
1	L	36/51 (71%)	35 (97%)	1 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	498/612 (81%)	493 (99%)	5 (1%)	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	44/47 (94%)	44 (100%)	0	100	100
1	B	44/47 (94%)	42 (96%)	2 (4%)	27	39
1	C	45/47 (96%)	42 (93%)	3 (7%)	16	21
1	D	44/47 (94%)	42 (96%)	2 (4%)	27	39
1	E	44/47 (94%)	44 (100%)	0	100	100
1	F	45/47 (96%)	43 (96%)	2 (4%)	28	39
1	G	39/47 (83%)	38 (97%)	1 (3%)	46	63
1	H	39/47 (83%)	37 (95%)	2 (5%)	24	33
1	I	39/47 (83%)	37 (95%)	2 (5%)	24	33
1	J	41/47 (87%)	40 (98%)	1 (2%)	49	66
1	K	38/47 (81%)	38 (100%)	0	100	100
1	L	37/47 (79%)	35 (95%)	2 (5%)	22	30
All	All	499/564 (88%)	482 (97%)	17 (3%)	37	51

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	233	GLN
1	B	257	LYS
1	D	225	LEU
1	D	233	GLN
1	C	222	ARG
1	C	225	LEU

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Mol	Chain	Res	Type
1	C	258	LEU
1	F	222	ARG
1	F	267	ASP
1	H	244	SER
1	H	253	GLU
1	G	257	LYS
1	I	222	ARG
1	I	229	GLN
1	J	246	SER
1	L	246	SER
1	L	257	LYS

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

Mol	Chain	Res	Type
1	B	229	GLN

### 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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	46/51 (90%)	0.20	0 <span>100</span> <span>100</span>	32, 44, 56, 61	0
1	B	46/51 (90%)	0.36	1 (2%) <span>62</span> <span>69</span>	22, 40, 54, 72	0
1	C	48/51 (94%)	0.17	0 <span>100</span> <span>100</span>	28, 43, 57, 70	0
1	D	48/51 (94%)	0.30	1 (2%) <span>63</span> <span>70</span>	31, 43, 59, 68	0
1	E	46/51 (90%)	0.23	1 (2%) <span>62</span> <span>69</span>	24, 39, 57, 68	0
1	F	49/51 (96%)	0.39	1 (2%) <span>65</span> <span>71</span>	29, 43, 59, 65	0
1	G	40/51 (78%)	0.51	3 (7%) <span>14</span> <span>19</span>	28, 42, 65, 68	0
1	H	40/51 (78%)	0.34	3 (7%) <span>14</span> <span>19</span>	32, 46, 70, 75	0
1	I	40/51 (78%)	0.35	1 (2%) <span>57</span> <span>64</span>	38, 50, 65, 67	0
1	J	42/51 (82%)	0.48	3 (7%) <span>16</span> <span>21</span>	41, 54, 69, 81	0
1	K	39/51 (76%)	0.44	4 (10%) <span>6</span> <span>9</span>	31, 47, 72, 81	0
1	L	38/51 (74%)	0.27	0 <span>100</span> <span>100</span>	31, 44, 66, 71	0
All	All	522/612 (85%)	0.33	18 (3%) <span>45</span> <span>52</span>	22, 45, 67, 81	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	219	GLY	5.9
1	B	265	LEU	4.1
1	K	258	LEU	3.7
1	I	259	ILE	3.5
1	H	222	ARG	3.5
1	H	258	LEU	3.2
1	K	222	ARG	3.2
1	G	258	LEU	3.1
1	G	257	LYS	3.1
1	J	260	LYS	3.0
1	J	259	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	G	221	GLU	2.7
1	K	225	LEU	2.6
1	H	260	LYS	2.6
1	K	223	GLN	2.4
1	E	260	LYS	2.4
1	F	225	LEU	2.1
1	J	220	SER	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 carbohydrates 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.