



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 02:08 PM GMT

PDB ID : 4LML
Title : GLIC double mutant I9'A T25'A
Authors : Grosman, C.; Gonzalez-Gutierrez, G.
Deposited on : 2013-07-10
Resolution : 3.80 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

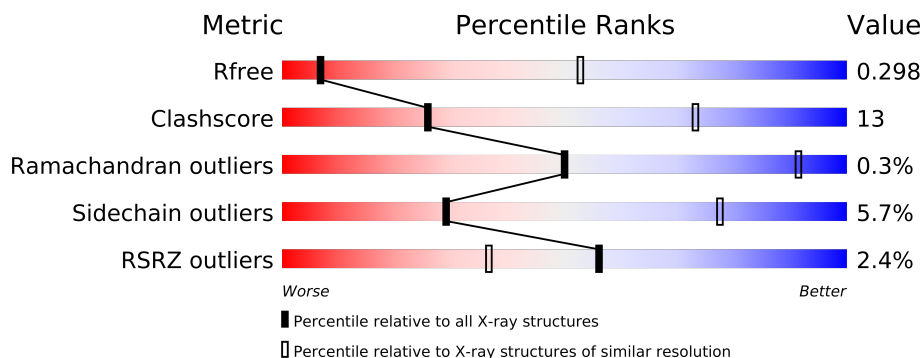
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.80 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}	66092	1162 (4.20-3.40)
Clashscore	79885	1100 (4.10-3.50)
Ramachandran outliers	78287	1050 (4.10-3.50)
Sidechain outliers	78261	1042 (4.10-3.50)
RSRZ outliers	66119	1163 (4.20-3.40)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	318	
1	B	318	
1	C	318	
1	D	318	
1	E	318	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 12580 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Proton-gated ion channel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	310	Total	C	N	O	S	0	0	0
			2516	1658	403	451	4			
1	B	310	Total	C	N	O	S	0	0	0
			2516	1658	403	451	4			
1	C	310	Total	C	N	O	S	0	0	0
			2516	1658	403	451	4			
1	D	310	Total	C	N	O	S	0	0	0
			2516	1658	403	451	4			
1	E	310	Total	C	N	O	S	0	0	0
			2516	1658	403	451	4			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP Q7NDN8
A	0	SER	-	EXPRESSION TAG	UNP Q7NDN8
A	232	ALA	ILE	ENGINEERED MUTATION	UNP Q7NDN8
A	248	ALA	THR	ENGINEERED MUTATION	UNP Q7NDN8
B	-1	GLY	-	EXPRESSION TAG	UNP Q7NDN8
B	0	SER	-	EXPRESSION TAG	UNP Q7NDN8
B	232	ALA	ILE	ENGINEERED MUTATION	UNP Q7NDN8
B	248	ALA	THR	ENGINEERED MUTATION	UNP Q7NDN8
C	-1	GLY	-	EXPRESSION TAG	UNP Q7NDN8
C	0	SER	-	EXPRESSION TAG	UNP Q7NDN8
C	232	ALA	ILE	ENGINEERED MUTATION	UNP Q7NDN8
C	248	ALA	THR	ENGINEERED MUTATION	UNP Q7NDN8
D	-1	GLY	-	EXPRESSION TAG	UNP Q7NDN8
D	0	SER	-	EXPRESSION TAG	UNP Q7NDN8
D	232	ALA	ILE	ENGINEERED MUTATION	UNP Q7NDN8
D	248	ALA	THR	ENGINEERED MUTATION	UNP Q7NDN8
E	-1	GLY	-	EXPRESSION TAG	UNP Q7NDN8
E	0	SER	-	EXPRESSION TAG	UNP Q7NDN8
E	232	ALA	ILE	ENGINEERED MUTATION	UNP Q7NDN8

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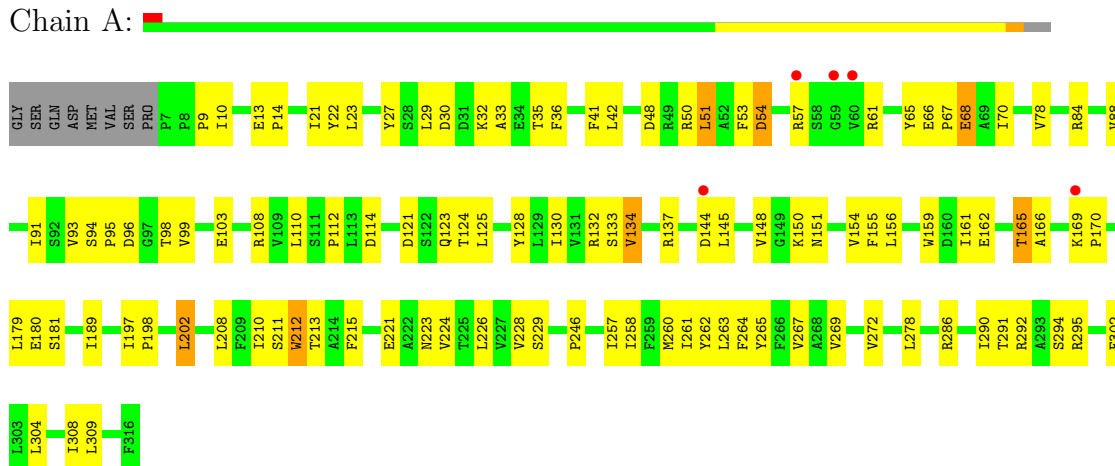
Chain	Residue	Modelled	Actual	Comment	Reference
E	248	ALA	THR	ENGINEERED MUTATION	UNP Q7NDN8

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 errors 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.

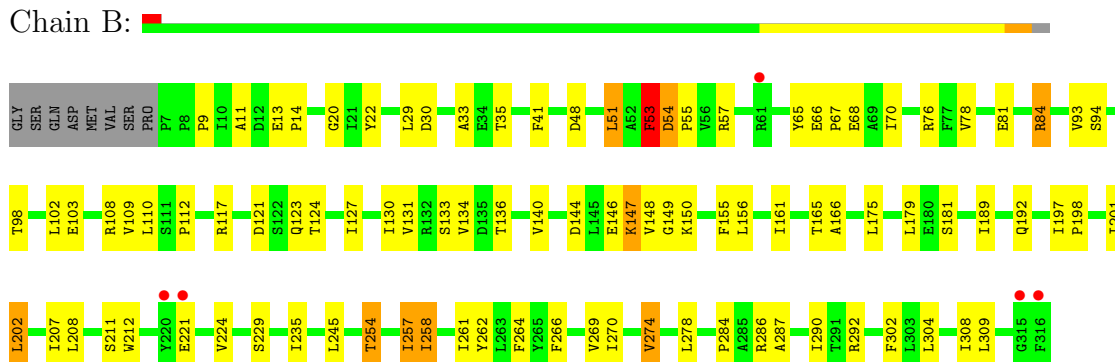
• Molecule 1: Proton-gated ion channel

Chain A:



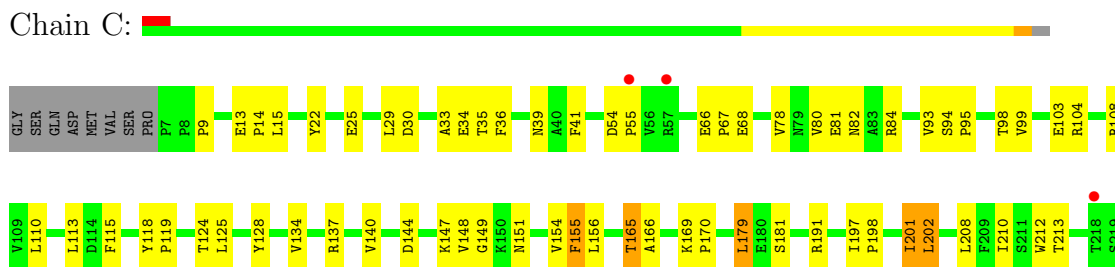
• Molecule 1: Proton-gated ion channel

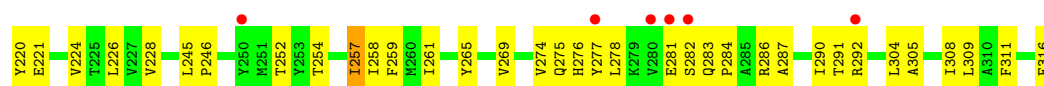
Chain B:



• Molecule 1: Proton-gated ion channel

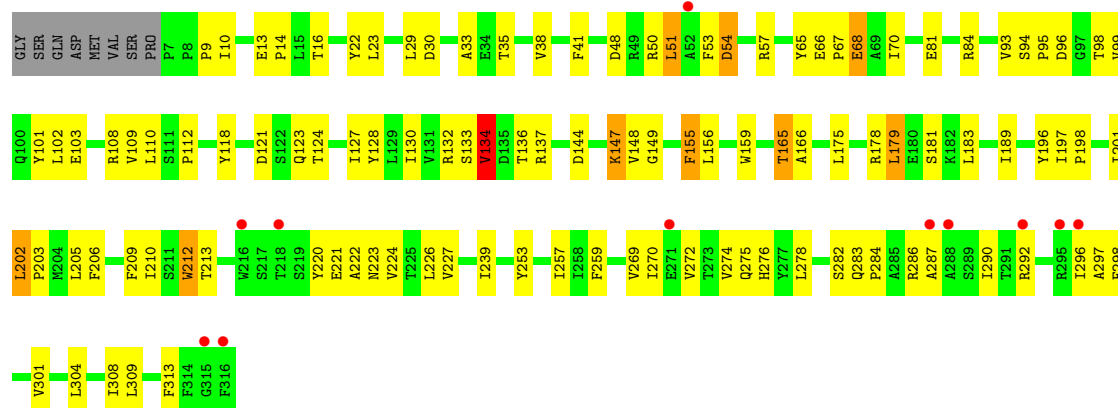
Chain C:





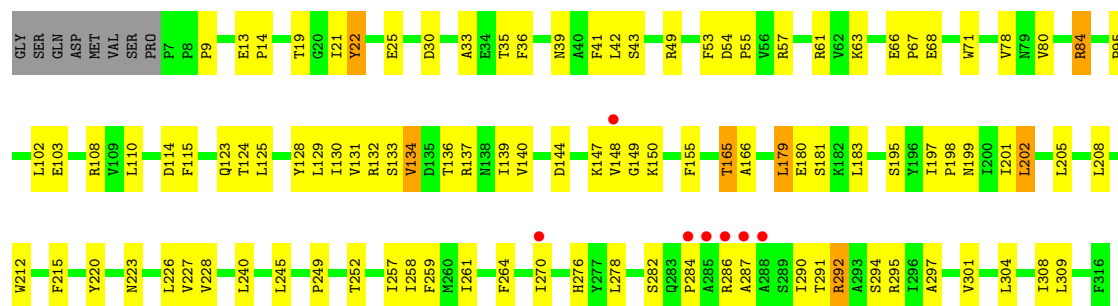
• Molecule 1: Proton-gated ion channel

Chain D:



• Molecule 1: Proton-gated ion channel

Chain E:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	122.80Å 128.00Å 190.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.93 – 3.80 39.43 – 3.79	Depositor EDS
% Data completeness (in resolution range)	72.2 (38.93-3.80) 63.9 (39.43-3.79)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 3.76Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.246 , 0.302 0.243 , 0.298	Depositor DCC
R_{free} test set	1801 reflections (8.35%)	DCC
Wilson B-factor (Å ²)	24.3	Xtriage
Anisotropy	1.738	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 19.8	EDS
Estimated twinning fraction	0.032 for k,h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 24257 reflections	Xtriage
F_o, F_c correlation	0.75	EDS
Total number of atoms	12580	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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	0/2584	0.60	0/3528
1	B	0.38	0/2584	0.59	0/3528
1	C	0.37	0/2584	0.58	0/3528
1	D	0.39	0/2584	0.59	0/3528
1	E	0.38	0/2584	0.59	1/3528 (0.0%)
All	All	0.38	0/12920	0.59	1/17640 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	149	GLY	N-CA-C	6.03	128.18	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	53	PHE	Peptide

5.2 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 hydrogens

added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2516	0	2529	71	0
1	B	2516	0	2529	72	1
1	C	2516	0	2529	76	0
1	D	2516	0	2529	78	1
1	E	2516	0	2529	72	0
All	All	12580	0	12645	340	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

All (340) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:121:ASP:OD1	1:B:123:GLN:NE2	2.04	0.90
1:B:117:ARG:HB2	1:B:121:ASP:HB3	1.60	0.83
1:C:22:TYR:HA	1:C:149:GLY:HA3	1.59	0.83
1:C:35:THR:OG1	1:C:108:ARG:NH2	2.12	0.82
1:D:51:LEU:HD12	1:D:93:VAL:HG11	1.63	0.81
1:C:84:ARG:NH1	1:C:103:GLU:OE1	2.15	0.78
1:B:29:LEU:HB2	1:B:156:LEU:HD11	1.68	0.73
1:A:29:LEU:HB2	1:A:156:LEU:HD11	1.69	0.72
1:A:51:LEU:HD12	1:A:93:VAL:HG11	1.72	0.72
1:D:54:ASP:HB2	1:D:57:ARG:HB2	1.71	0.71
1:D:84:ARG:NH1	1:D:103:GLU:OE1	2.15	0.71
1:D:22:TYR:HB3	1:D:41:PHE:HB2	1.73	0.70
1:B:22:TYR:HA	1:B:149:GLY:HA3	1.74	0.69
1:B:208:LEU:HD13	1:B:261:ILE:HG23	1.72	0.69
1:D:121:ASP:OD1	1:D:123:GLN:NE2	2.27	0.68
1:D:257:ILE:HG22	1:D:309:LEU:HD23	1.75	0.67
1:B:144:ASP:HA	1:B:147:LYS:HD2	1.76	0.66
1:E:278:LEU:HD21	1:E:286:ARG:HB3	1.77	0.66
1:A:22:TYR:HB3	1:A:41:PHE:HB2	1.77	0.66
1:D:175:LEU:HD11	1:E:22:TYR:CE1	2.32	0.65
1:C:22:TYR:HB3	1:C:41:PHE:HB2	1.79	0.65
1:E:208:LEU:HD13	1:E:261:ILE:HG23	1.77	0.65
1:B:53:PHE:HD2	1:B:53:PHE:H	1.45	0.65
1:D:94:SER:HB2	1:D:98:THR:HB	1.78	0.64
1:B:257:ILE:HG22	1:B:309:LEU:HD23	1.80	0.64
1:A:78:VAL:HB	1:A:128:TYR:HB2	1.80	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:269:VAL:HG11	1:D:210:ILE:HG23	1.80	0.63
1:E:257:ILE:HG22	1:E:309:LEU:HD23	1.81	0.63
1:C:35:THR:HG22	1:C:110:LEU:HG	1.80	0.63
1:D:165:THR:OG1	1:D:166:ALA:N	2.33	0.62
1:B:94:SER:HB2	1:B:98:THR:HB	1.80	0.62
1:D:112:PRO:HB3	1:E:155:PHE:CZ	2.35	0.62
1:D:278:LEU:HD21	1:D:286:ARG:HB3	1.82	0.61
1:A:210:ILE:HD13	1:C:269:VAL:HG21	1.83	0.61
1:E:165:THR:OG1	1:E:166:ALA:N	2.34	0.61
1:B:208:LEU:O	1:B:211:SER:HB3	2.01	0.60
1:E:22:TYR:HB3	1:E:41:PHE:HB2	1.83	0.60
1:C:208:LEU:HD13	1:C:261:ILE:HG23	1.83	0.60
1:B:54:ASP:HB2	1:B:57:ARG:HB2	1.84	0.59
1:A:30:ASP:HB3	1:A:33:ALA:HB3	1.84	0.59
1:A:66:GLU:HG3	1:A:67:PRO:HD2	1.85	0.59
1:B:197:ILE:HB	1:B:198:PRO:HD3	1.84	0.59
1:C:25:GLU:HB2	1:C:39:ASN:HB3	1.84	0.59
1:E:53:PHE:CE2	1:E:63:LYS:HB3	2.37	0.59
1:A:257:ILE:HG22	1:A:309:LEU:HD23	1.85	0.59
1:B:112:PRO:HB3	1:C:155:PHE:HZ	1.67	0.59
1:A:165:THR:OG1	1:A:166:ALA:N	2.34	0.59
1:E:25:GLU:HB2	1:E:39:ASN:HB3	1.84	0.58
1:B:165:THR:OG1	1:B:166:ALA:N	2.35	0.58
1:D:30:ASP:HB3	1:D:33:ALA:HB3	1.85	0.58
1:E:42:LEU:HB3	1:E:103:GLU:HG2	1.86	0.58
1:B:112:PRO:HB3	1:C:155:PHE:CZ	2.39	0.58
1:C:140:VAL:HG22	1:C:181:SER:HB3	1.86	0.58
1:D:29:LEU:HB2	1:D:156:LEU:HD11	1.85	0.57
1:A:278:LEU:HD21	1:A:286:ARG:HB3	1.86	0.57
1:B:53:PHE:HE1	1:B:55:PRO:HG3	1.68	0.57
1:B:22:TYR:HB3	1:B:41:PHE:HB2	1.86	0.57
1:D:175:LEU:HD11	1:E:22:TYR:HE1	1.68	0.57
1:D:134:VAL:O	1:D:178:ARG:NH2	2.38	0.57
1:A:144:ASP:O	1:A:148:VAL:HB	2.05	0.57
1:D:227:VAL:HG11	1:D:269:VAL:HG23	1.86	0.57
1:C:197:ILE:O	1:C:202:LEU:HB2	2.04	0.57
1:B:51:LEU:HD23	1:B:51:LEU:H	1.69	0.57
1:C:29:LEU:HB2	1:C:156:LEU:HD11	1.87	0.57
1:D:147:LYS:HG2	1:D:147:LYS:O	2.03	0.57
1:E:66:GLU:HG3	1:E:67:PRO:HD2	1.87	0.56
1:B:53:PHE:CE1	1:B:55:PRO:HG3	2.40	0.56
1:E:132:ARG:HA	1:E:180:GLU:HG2	1.86	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:54:ASP:HB2	1:A:57:ARG:HB2	1.88	0.56
1:E:140:VAL:HG23	1:E:183:LEU:HG	1.87	0.56
1:B:11:ALA:HB3	1:B:136:THR:HG21	1.88	0.56
1:A:202:LEU:HD12	1:C:259:PHE:CZ	2.41	0.56
1:D:22:TYR:HA	1:D:149:GLY:HA3	1.88	0.55
1:D:197:ILE:HA	1:D:201:ILE:HB	1.88	0.55
1:A:114:ASP:O	1:A:123:GLN:NE2	2.39	0.55
1:A:132:ARG:HA	1:A:180:GLU:HG2	1.89	0.55
1:E:78:VAL:HB	1:E:128:TYR:HB2	1.88	0.55
1:A:32:LYS:HA	1:A:246:PRO:HD3	1.89	0.55
1:E:84:ARG:NH2	1:E:103:GLU:OE1	2.19	0.55
1:E:61:ARG:O	1:E:95:PRO:HG3	2.07	0.55
1:A:112:PRO:HB3	1:D:155:PHE:CZ	2.43	0.54
1:B:144:ASP:O	1:B:148:VAL:HG13	2.07	0.54
1:C:278:LEU:HD21	1:C:286:ARG:HB3	1.90	0.54
1:B:66:GLU:HG3	1:B:67:PRO:HD2	1.90	0.54
1:C:66:GLU:HG3	1:C:67:PRO:HD2	1.88	0.54
1:D:9:PRO:HB3	1:D:48:ASP:CG	2.29	0.53
1:D:304:LEU:O	1:D:308:ILE:HG12	2.08	0.53
1:E:202:LEU:HD22	1:E:205:LEU:HD12	1.90	0.53
1:E:133:SER:HB3	1:E:136:THR:O	2.08	0.53
1:E:30:ASP:HB3	1:E:33:ALA:HB3	1.89	0.53
1:D:137:ARG:HB3	1:D:179:LEU:HD23	1.89	0.53
1:A:262:TYR:CE1	1:D:203:PRO:HB3	2.44	0.53
1:C:55:PRO:HD3	1:C:95:PRO:HB3	1.91	0.53
1:A:133:SER:HB3	1:A:137:ARG:HA	1.90	0.53
1:A:96:ASP:N	1:A:96:ASP:OD1	2.42	0.53
1:E:197:ILE:HB	1:E:198:PRO:HD3	1.91	0.53
1:E:304:LEU:O	1:E:308:ILE:HG12	2.09	0.52
1:A:161:ILE:HA	1:A:189:ILE:HG22	1.89	0.52
1:E:57:ARG:HA	1:E:57:ARG:HE	1.75	0.52
1:E:144:ASP:O	1:E:148:VAL:HG13	2.09	0.52
1:D:197:ILE:HB	1:D:198:PRO:HD3	1.91	0.52
1:D:130:ILE:HA	1:D:181:SER:O	2.09	0.52
1:B:109:VAL:HG11	1:B:127:ILE:HG12	1.92	0.52
1:E:286:ARG:O	1:E:290:ILE:HG13	2.10	0.52
1:D:156:LEU:HD23	1:D:189:ILE:HG21	1.90	0.52
1:C:30:ASP:HB3	1:C:33:ALA:HB3	1.92	0.52
1:D:286:ARG:O	1:D:290:ILE:HG13	2.09	0.51
1:C:197:ILE:HA	1:C:201:ILE:HB	1.92	0.51
1:A:223:ASN:HB3	1:A:272:VAL:HG11	1.92	0.51
1:C:197:ILE:HB	1:C:198:PRO:HD3	1.91	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:93:VAL:HG22	1:C:99:VAL:HG22	1.91	0.51
1:B:147:LYS:O	1:B:147:LYS:HG2	2.11	0.51
1:C:147:LYS:C	1:C:149:GLY:H	2.14	0.51
1:A:264:PHE:HA	1:A:267:VAL:HG22	1.93	0.51
1:C:137:ARG:HB3	1:C:179:LEU:HD21	1.93	0.51
1:E:292:ARG:HD2	1:E:295:ARG:HD2	1.91	0.51
1:D:10:ILE:HD12	1:D:50:ARG:HD3	1.92	0.51
1:B:284:PRO:HA	1:B:287:ALA:HB3	1.93	0.51
1:A:65:TYR:CG	1:A:70:ILE:HD11	2.45	0.51
1:B:20:GLY:HA3	1:B:146:GLU:HB3	1.92	0.51
1:E:197:ILE:HA	1:E:201:ILE:HB	1.93	0.50
1:E:215:PHE:HB3	1:E:295:ARG:HG2	1.94	0.50
1:D:144:ASP:O	1:D:148:VAL:HG13	2.12	0.50
1:C:118:TYR:O	1:C:254:THR:OG1	2.28	0.50
1:B:269:VAL:HG11	1:C:210:ILE:HG23	1.94	0.50
1:D:41:PHE:CD1	1:D:102:LEU:HD11	2.46	0.50
1:D:133:SER:HB3	1:D:136:THR:O	2.12	0.50
1:D:147:LYS:C	1:D:149:GLY:H	2.15	0.49
1:C:286:ARG:O	1:C:290:ILE:HG13	2.11	0.49
1:C:144:ASP:O	1:C:148:VAL:HG13	2.13	0.49
1:E:41:PHE:CD1	1:E:102:LEU:HD11	2.47	0.49
1:A:65:TYR:O	1:A:91:ILE:HB	2.13	0.49
1:B:35:THR:OG1	1:B:108:ARG:NH2	2.46	0.49
1:C:278:LEU:HD23	1:C:287:ALA:HA	1.95	0.49
1:E:215:PHE:CZ	1:E:294:SER:HB3	2.48	0.49
1:A:291:THR:O	1:A:295:ARG:HG3	2.12	0.49
1:D:66:GLU:HG3	1:D:67:PRO:HD2	1.95	0.49
1:D:93:VAL:HG22	1:D:99:VAL:HG13	1.94	0.48
1:D:94:SER:N	1:D:98:THR:O	2.39	0.48
1:B:13:GLU:HB3	1:B:14:PRO:HD2	1.94	0.48
1:D:275:GLN:HG3	1:D:287:ALA:HB1	1.94	0.48
1:C:80:VAL:HG13	1:C:108:ARG:O	2.14	0.48
1:D:35:THR:OG1	1:D:108:ARG:NH2	2.45	0.48
1:B:266:PHE:HA	1:C:210:ILE:HD11	1.96	0.48
1:B:175:LEU:HD11	1:C:22:TYR:CE1	2.48	0.48
1:D:13:GLU:HB3	1:D:14:PRO:HD2	1.96	0.48
1:E:36:PHE:CD1	1:E:125:LEU:HD13	2.49	0.48
1:A:304:LEU:O	1:A:308:ILE:HG12	2.14	0.48
1:E:264:PHE:HZ	1:E:301:VAL:HG12	1.78	0.48
1:B:304:LEU:O	1:B:308:ILE:HG12	2.13	0.48
1:E:9:PRO:HD3	1:E:71:TRP:CE3	2.49	0.48
1:D:259:PHE:CZ	1:E:202:LEU:HD12	2.48	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:13:GLU:HB3	1:C:14:PRO:HD2	1.96	0.48
1:D:239:ILE:HG12	1:E:240:LEU:HD22	1.96	0.47
1:D:65:TYR:CG	1:D:70:ILE:HD11	2.48	0.47
1:D:253:TYR:HD1	1:D:313:PHE:CD1	2.33	0.47
1:B:147:LYS:C	1:B:149:GLY:H	2.16	0.47
1:C:257:ILE:HG22	1:C:309:LEU:HD23	1.96	0.47
1:B:35:THR:HG22	1:B:110:LEU:HG	1.96	0.47
1:E:80:VAL:HG13	1:E:108:ARG:O	2.14	0.47
1:E:130:ILE:HA	1:E:181:SER:O	2.14	0.47
1:B:270:ILE:HA	1:B:270:ILE:HD13	1.79	0.47
1:C:284:PRO:HA	1:C:287:ALA:HB3	1.96	0.47
1:A:263:LEU:HD12	1:D:206:PHE:CE1	2.50	0.47
1:B:197:ILE:HG23	1:B:202:LEU:HG	1.96	0.47
1:E:131:VAL:HG11	1:E:140:VAL:HG13	1.96	0.47
1:A:53:PHE:CD2	1:A:96:ASP:HA	2.50	0.47
1:A:197:ILE:HB	1:A:198:PRO:HD3	1.97	0.47
1:E:195:SER:O	1:E:199:ASN:ND2	2.48	0.47
1:E:137:ARG:HB3	1:E:179:LEU:HD23	1.97	0.47
1:D:101:TYR:OH	1:D:103:GLU:OE2	2.25	0.47
1:B:65:TYR:CG	1:B:70:ILE:HD11	2.50	0.47
1:B:76:ARG:HH11	1:C:104:ARG:HH11	1.63	0.47
1:C:311:PHE:CZ	1:C:316:PHE:HB2	2.49	0.47
1:B:150:LYS:HE2	1:B:150:LYS:HB3	1.68	0.46
1:B:278:LEU:HD21	1:B:286:ARG:HB3	1.97	0.46
1:D:283:GLN:N	1:D:284:PRO:HD3	2.31	0.46
1:E:19:THR:HA	1:E:43:SER:O	2.14	0.46
1:E:35:THR:HG22	1:E:110:LEU:HG	1.97	0.46
1:D:81:GLU:HG3	1:D:108:ARG:HG3	1.98	0.46
1:B:192:GLN:OE1	1:E:249:PRO:HG3	2.16	0.46
1:E:147:LYS:HG2	1:E:147:LYS:O	2.16	0.46
1:E:21:ILE:HD11	1:E:129:LEU:HD11	1.97	0.46
1:A:13:GLU:HB3	1:A:14:PRO:HD2	1.98	0.46
1:C:304:LEU:O	1:C:308:ILE:HG12	2.15	0.46
1:C:147:LYS:O	1:C:147:LYS:HG2	2.16	0.46
1:C:118:TYR:HD1	1:C:191:ARG:NH2	2.13	0.46
1:A:21:ILE:HA	1:A:41:PHE:O	2.16	0.46
1:B:202:LEU:HD12	1:E:259:PHE:HZ	1.81	0.46
1:E:78:VAL:HG22	1:E:130:ILE:HG12	1.98	0.46
1:A:212:TRP:HB2	1:A:215:PHE:CE1	2.51	0.46
1:B:30:ASP:HB3	1:B:33:ALA:HB3	1.98	0.46
1:E:114:ASP:O	1:E:123:GLN:NE2	2.47	0.46
1:C:147:LYS:HE3	1:C:166:ALA:H	1.80	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:93:VAL:HG22	1:A:99:VAL:HG13	1.97	0.45
1:E:131:VAL:CG1	1:E:140:VAL:HG13	2.47	0.45
1:B:133:SER:HB3	1:B:136:THR:O	2.16	0.45
1:B:274:VAL:O	1:B:278:LEU:HB2	2.16	0.45
1:C:309:LEU:HD12	1:C:309:LEU:HA	1.71	0.45
1:E:55:PRO:HD3	1:E:95:PRO:HB3	1.97	0.45
1:E:284:PRO:HA	1:E:287:ALA:HB3	1.98	0.45
1:B:9:PRO:HB3	1:B:48:ASP:OD1	2.16	0.45
1:A:208:LEU:HD13	1:A:261:ILE:HG23	1.98	0.45
1:B:78:VAL:HG22	1:B:130:ILE:HG12	1.99	0.45
1:D:53:PHE:HB3	1:D:95:PRO:C	2.36	0.45
1:A:151:ASN:O	1:A:154:VAL:HG22	2.16	0.45
1:C:118:TYR:CD1	1:C:119:PRO:HA	2.51	0.45
1:D:212:TRP:HE3	1:D:298:PHE:HD1	1.64	0.45
1:B:254:THR:O	1:B:258:ILE:HB	2.17	0.45
1:A:9:PRO:HB3	1:A:48:ASP:OD1	2.16	0.45
1:C:221:GLU:O	1:C:224:VAL:HG22	2.17	0.45
1:C:36:PHE:CD1	1:C:125:LEU:HD13	2.51	0.45
1:B:202:LEU:HD12	1:E:259:PHE:CZ	2.52	0.45
1:C:197:ILE:HG23	1:C:202:LEU:HD23	1.99	0.45
1:B:13:GLU:HB3	1:B:14:PRO:CD	2.47	0.45
1:D:159:TRP:CE3	1:D:189:ILE:HD12	2.52	0.44
1:C:80:VAL:HG12	1:C:82:ASN:O	2.18	0.44
1:B:41:PHE:CD1	1:B:102:LEU:HD11	2.52	0.44
1:E:291:THR:O	1:E:295:ARG:HG3	2.17	0.44
1:B:131:VAL:HG11	1:B:140:VAL:HG13	1.99	0.44
1:C:165:THR:OG1	1:C:166:ALA:N	2.50	0.44
1:D:35:THR:HG22	1:D:110:LEU:HG	1.98	0.44
1:E:297:ALA:O	1:E:301:VAL:HG23	2.17	0.44
1:A:68:GLU:CD	1:A:68:GLU:H	2.20	0.44
1:C:277:TYR:O	1:C:281:GLU:HG2	2.18	0.44
1:D:175:LEU:HD12	1:D:175:LEU:HA	1.68	0.44
1:A:260:MET:O	1:A:263:LEU:HB2	2.16	0.44
1:A:67:PRO:HG2	1:A:88:VAL:HG21	2.00	0.44
1:B:81:GLU:HG3	1:B:108:ARG:HG3	2.00	0.44
1:E:264:PHE:CZ	1:E:301:VAL:HG12	2.52	0.44
1:B:207:ILE:HG13	1:B:208:LEU:N	2.33	0.44
1:C:29:LEU:HD12	1:C:125:LEU:HD11	2.00	0.44
1:B:286:ARG:O	1:B:290:ILE:HG13	2.18	0.44
1:D:212:TRP:CE3	1:D:298:PHE:HD1	2.36	0.44
1:E:223:ASN:O	1:E:227:VAL:HG23	2.18	0.44
1:E:13:GLU:HB3	1:E:14:PRO:HD2	1.99	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:51:LEU:HD23	1:D:51:LEU:N	2.33	0.44
1:A:9:PRO:HB3	1:A:48:ASP:CG	2.37	0.44
1:A:221:GLU:OE1	1:D:222:ALA:HA	2.18	0.44
1:C:274:VAL:O	1:C:278:LEU:HB2	2.18	0.43
1:A:121:ASP:OD1	1:A:123:GLN:NE2	2.49	0.43
1:B:130:ILE:HA	1:B:181:SER:O	2.18	0.43
1:D:96:ASP:N	1:D:96:ASP:OD1	2.51	0.43
1:B:197:ILE:O	1:B:202:LEU:HB2	2.18	0.43
1:A:215:PHE:CE1	1:A:294:SER:HB3	2.54	0.43
1:A:261:ILE:O	1:A:265:TYR:HD1	2.01	0.43
1:A:130:ILE:HA	1:A:181:SER:O	2.18	0.43
1:B:229:SER:HB3	1:E:228:VAL:HG11	2.00	0.43
1:C:169:LYS:HA	1:C:170:PRO:HD2	1.86	0.43
1:E:220:TYR:CD1	1:E:276:HIS:HB2	2.53	0.43
1:A:286:ARG:O	1:A:290:ILE:HG13	2.18	0.43
1:B:76:ARG:NH2	1:B:130:ILE:HD12	2.34	0.43
1:C:94:SER:HB2	1:C:98:THR:HB	2.01	0.43
1:C:78:VAL:HB	1:C:128:TYR:HB2	2.00	0.43
1:A:10:ILE:HD12	1:A:50:ARG:HD3	2.00	0.43
1:B:51:LEU:HD12	1:B:93:VAL:HG11	2.00	0.43
1:C:34:GLU:OE2	1:C:113:LEU:N	2.39	0.43
1:A:53:PHE:HB3	1:A:95:PRO:CA	2.48	0.43
1:A:159:TRP:CE3	1:A:189:ILE:HD12	2.53	0.43
1:E:215:PHE:CE1	1:E:294:SER:HB3	2.54	0.43
1:A:224:VAL:O	1:A:228:VAL:HB	2.19	0.43
1:D:223:ASN:HB3	1:D:272:VAL:HG11	2.01	0.43
1:B:269:VAL:HG21	1:C:210:ILE:HD13	2.01	0.43
1:A:36:PHE:CD1	1:A:125:LEU:HD13	2.54	0.43
1:A:94:SER:N	1:A:98:THR:O	2.44	0.43
1:B:309:LEU:HA	1:B:309:LEU:HD12	1.81	0.43
1:A:202:LEU:HD12	1:C:259:PHE:HZ	1.82	0.43
1:A:137:ARG:HD3	1:A:137:ARG:HA	1.73	0.43
1:C:22:TYR:CD1	1:C:149:GLY:HA2	2.53	0.42
1:B:156:LEU:HD12	1:B:156:LEU:HA	1.78	0.42
1:D:259:PHE:HZ	1:E:202:LEU:HD12	1.83	0.42
1:C:137:ARG:HD3	1:C:137:ARG:HA	1.89	0.42
1:B:161:ILE:HA	1:B:189:ILE:HG22	2.00	0.42
1:D:202:LEU:HD22	1:D:205:LEU:HD12	1.99	0.42
1:E:139:ILE:HA	1:E:181:SER:OG	2.19	0.42
1:A:208:LEU:O	1:A:211:SER:HB3	2.18	0.42
1:B:53:PHE:N	1:B:53:PHE:CD2	2.82	0.42
1:C:282:SER:C	1:C:284:PRO:HD3	2.40	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:282:SER:C	1:E:284:PRO:HD3	2.40	0.42
1:C:151:ASN:O	1:C:154:VAL:HG22	2.20	0.42
1:A:221:GLU:O	1:A:224:VAL:HG22	2.20	0.42
1:B:229:SER:CB	1:E:228:VAL:HG11	2.50	0.42
1:A:42:LEU:HB3	1:A:103:GLU:HG2	2.01	0.42
1:C:144:ASP:HA	1:C:147:LYS:HD2	2.01	0.42
1:D:13:GLU:HB3	1:D:14:PRO:CD	2.49	0.42
1:A:229:SER:HB3	1:C:228:VAL:HG11	2.02	0.42
1:B:221:GLU:O	1:B:224:VAL:HG22	2.20	0.42
1:C:261:ILE:O	1:C:265:TYR:HD1	2.02	0.42
1:D:282:SER:C	1:D:284:PRO:HD3	2.40	0.42
1:D:109:VAL:HG11	1:D:127:ILE:HG12	2.02	0.42
1:C:245:LEU:HD12	1:C:245:LEU:HA	1.74	0.42
1:A:61:ARG:O	1:A:95:PRO:HG3	2.20	0.42
1:C:13:GLU:HB3	1:C:14:PRO:CD	2.49	0.42
1:D:297:ALA:O	1:D:301:VAL:HG23	2.20	0.42
1:D:220:TYR:CD1	1:D:276:HIS:HB2	2.55	0.42
1:A:27:TYR:OH	1:C:81:GLU:O	2.36	0.41
1:D:84:ARG:HH11	1:D:84:ARG:HD2	1.75	0.41
1:B:197:ILE:HA	1:B:201:ILE:HB	2.01	0.41
1:D:155:PHE:HA	1:D:155:PHE:HD2	1.65	0.41
1:A:35:THR:HG22	1:A:110:LEU:HG	2.02	0.41
1:C:220:TYR:CD1	1:C:276:HIS:HB2	2.55	0.41
1:C:275:GLN:OE1	1:C:291:THR:OG1	2.28	0.41
1:E:270:ILE:HD13	1:E:270:ILE:HA	1.86	0.41
1:E:49:ARG:H	1:E:49:ARG:HG2	1.64	0.41
1:D:128:TYR:HA	1:D:183:LEU:O	2.20	0.41
1:B:264:PHE:HE2	1:B:302:PHE:HB2	1.85	0.41
1:B:84:ARG:NH2	1:B:103:GLU:OE1	2.44	0.41
1:C:115:PHE:O	1:C:252:THR:HG22	2.20	0.41
1:E:133:SER:HB3	1:E:137:ARG:HA	2.03	0.41
1:A:35:THR:OG1	1:A:108:ARG:NH2	2.46	0.41
1:B:245:LEU:HA	1:B:245:LEU:HD12	1.82	0.41
1:E:245:LEU:HD12	1:E:245:LEU:HA	1.87	0.41
1:A:264:PHE:CE2	1:A:302:PHE:HD1	2.39	0.41
1:A:27:TYR:HB3	1:C:110:LEU:HD11	2.03	0.41
1:C:147:LYS:HE2	1:C:165:THR:HA	2.02	0.41
1:E:13:GLU:HB3	1:E:14:PRO:CD	2.51	0.41
1:D:296:ILE:HD13	1:D:296:ILE:HA	1.91	0.41
1:C:9:PRO:HG2	1:C:15:LEU:HD22	2.03	0.41
1:A:169:LYS:HA	1:A:170:PRO:HD2	1.98	0.41
1:D:23:LEU:HD22	1:D:38:VAL:HG21	2.03	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:257:ILE:O	1:C:261:ILE:HG13	2.21	0.40
1:D:132:ARG:NH2	1:D:178:ARG:HB2	2.36	0.40
1:A:53:PHE:HD2	1:A:96:ASP:HA	1.84	0.40
1:D:209:PHE:HA	1:D:212:TRP:NE1	2.37	0.40
1:E:130:ILE:HD12	1:E:180:GLU:OE1	2.22	0.40
1:D:221:GLU:O	1:D:224:VAL:HG22	2.21	0.40
1:D:118:TYR:HH	1:D:196:TYR:HE2	1.68	0.40
1:C:305:ALA:O	1:C:309:LEU:HB2	2.21	0.40
1:D:9:PRO:HB3	1:D:48:ASP:OD1	2.21	0.40
1:C:245:LEU:HG	1:C:246:PRO:HD2	2.02	0.40
1:B:235:ILE:HG12	1:B:262:TYR:OH	2.20	0.40
1:D:270:ILE:HD13	1:D:270:ILE:HA	1.82	0.40
1:E:22:TYR:C	1:E:22:TYR:CD2	2.94	0.40
1:C:283:GLN:N	1:C:284:PRO:HD3	2.36	0.40
1:D:53:PHE:HB3	1:D:95:PRO:O	2.21	0.40
1:A:23:LEU:HB2	1:A:150:LYS:HB2	2.02	0.40
1:E:115:PHE:O	1:E:252:THR:HG22	2.20	0.40
1:A:202:LEU:HD22	1:A:202:LEU:HA	1.94	0.40
1:A:215:PHE:HB3	1:A:295:ARG:HG2	2.02	0.40
1:D:68:GLU:H	1:D:68:GLU:HG2	1.53	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:147:LYS:NZ	1:D:98:THR:OG1[3.554]	2.11	0.09

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	308/318 (97%)	291 (94%)	16 (5%)	1 (0%)	50 91
1	B	308/318 (97%)	290 (94%)	18 (6%)	0	100 100
1	C	308/318 (97%)	286 (93%)	21 (7%)	1 (0%)	50 91

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	308/318 (97%)	289 (94%)	18 (6%)	1 (0%)	50	91
1	E	308/318 (97%)	289 (94%)	18 (6%)	1 (0%)	50	91
All	All	1540/1590 (97%)	1445 (94%)	91 (6%)	4 (0%)	50	91

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	134	VAL
1	C	201	ILE
1	D	134	VAL
1	E	134	VAL

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	276/283 (98%)	259 (94%)	17 (6%)	26	75
1	B	276/283 (98%)	259 (94%)	17 (6%)	26	75
1	C	276/283 (98%)	262 (95%)	14 (5%)	33	81
1	D	276/283 (98%)	260 (94%)	16 (6%)	28	76
1	E	276/283 (98%)	262 (95%)	14 (5%)	33	81
All	All	1380/1415 (98%)	1302 (94%)	78 (6%)	29	77

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	LEU
1	A	54	ASP
1	A	68	GLU
1	A	84	ARG
1	A	124	THR
1	A	134	VAL
1	A	145	LEU
1	A	155	PHE

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Mol	Chain	Res	Type
1	A	162	GLU
1	A	165	THR
1	A	179	LEU
1	A	202	LEU
1	A	212	TRP
1	A	213	THR
1	A	226	LEU
1	A	258	ILE
1	A	292	ARG
1	B	51	LEU
1	B	53	PHE
1	B	54	ASP
1	B	68	GLU
1	B	84	ARG
1	B	124	THR
1	B	134	VAL
1	B	147	LYS
1	B	155	PHE
1	B	179	LEU
1	B	202	LEU
1	B	212	TRP
1	B	254	THR
1	B	257	ILE
1	B	258	ILE
1	B	274	VAL
1	B	292	ARG
1	C	54	ASP
1	C	68	GLU
1	C	124	THR
1	C	134	VAL
1	C	155	PHE
1	C	165	THR
1	C	179	LEU
1	C	202	LEU
1	C	212	TRP
1	C	213	THR
1	C	226	LEU
1	C	257	ILE
1	C	258	ILE
1	C	292	ARG
1	D	16	THR
1	D	51	LEU

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Mol	Chain	Res	Type
1	D	54	ASP
1	D	68	GLU
1	D	124	THR
1	D	134	VAL
1	D	147	LYS
1	D	155	PHE
1	D	165	THR
1	D	179	LEU
1	D	202	LEU
1	D	212	TRP
1	D	213	THR
1	D	226	LEU
1	D	274	VAL
1	D	292	ARG
1	E	22	TYR
1	E	54	ASP
1	E	68	GLU
1	E	84	ARG
1	E	124	THR
1	E	134	VAL
1	E	150	LYS
1	E	165	THR
1	E	179	LEU
1	E	202	LEU
1	E	212	TRP
1	E	226	LEU
1	E	258	ILE
1	E	292	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

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, 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	310/318 (97%)	0.21	5 (1%) 68 47	13, 48, 116, 160	0
1	B	310/318 (97%)	0.27	5 (1%) 68 47	12, 45, 103, 128	0
1	C	310/318 (97%)	0.33	9 (2%) 49 32	18, 50, 98, 154	0
1	D	310/318 (97%)	0.36	11 (3%) 42 28	14, 45, 107, 130	0
1	E	310/318 (97%)	0.30	7 (2%) 57 38	13, 45, 117, 139	0
All	All	1550/1590 (97%)	0.30	37 (2%) 56 36	12, 46, 108, 160	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	292	ARG	6.3
1	E	286	ARG	5.5
1	C	281	GLU	4.7
1	D	288	ALA	4.5
1	B	220	TYR	3.7
1	C	57	ARG	3.7
1	A	59	GLY	3.5
1	E	287	ALA	3.4
1	B	221	GLU	3.3
1	E	288	ALA	3.3
1	D	316	PHE	3.3
1	A	57	ARG	3.3
1	D	52	ALA	3.0
1	E	148	VAL	3.0
1	B	315	GLY	2.9
1	D	296	ILE	2.8
1	D	216	TRP	2.7
1	E	285	ALA	2.7
1	D	271	GLU	2.7
1	C	277	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	316	PHE	2.5
1	C	218	THR	2.4
1	C	55	PRO	2.4
1	A	144	ASP	2.3
1	E	270	ILE	2.3
1	A	169	LYS	2.3
1	E	284	PRO	2.2
1	D	315	GLY	2.2
1	A	60	VAL	2.2
1	C	292	ARG	2.1
1	B	61	ARG	2.1
1	C	282	SER	2.0
1	D	295	ARG	2.0
1	D	287	ALA	2.0
1	C	280	VAL	2.0
1	D	218	THR	2.0
1	C	250	TYR	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.