



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 08:08 AM UTC

PDB ID : 3TRH / pdb_00003trh
Title : Structure of a phosphoribosylaminoimidazole carboxylase catalytic subunit (purE) from *Coxiella burnetii*
Authors : Cheung, J.; Franklin, M.C.; Rudolph, M.; Cassidy, M.; Gary, E.; Burshteyn, F.; Love, J.
Deposited on : 2011-09-09
Resolution : 2.20 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
Mogul : 2022.3.0, CSD as543be (2022)
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

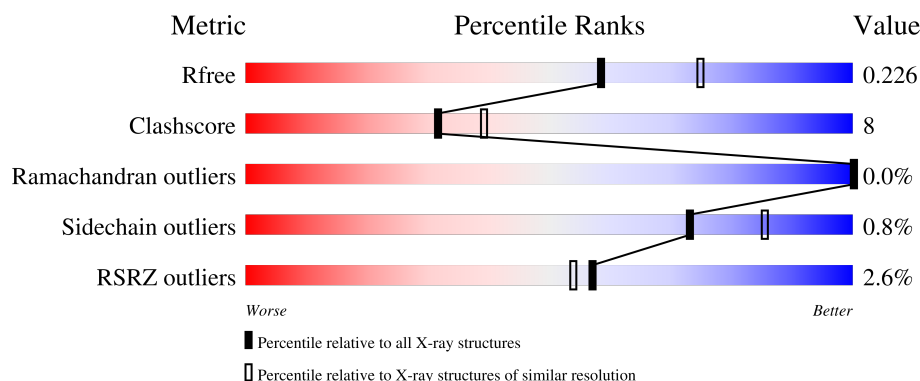
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	169	
1	B	169	
1	C	169	
1	D	169	
1	E	169	

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Mol	Chain	Length	Quality of chain
1	F	169	
1	G	169	
1	H	169	
1	I	169	
1	J	169	
1	K	169	
1	L	169	
1	M	169	
1	N	169	
1	O	169	
1	P	169	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 19690 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 Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	164	Total	C	N	O	S	Se	0	0	0
			1187	749	207	225	2	4			
1	B	164	Total	C	N	O	S	Se	0	0	0
			1187	749	207	225	2	4			
1	C	163	Total	C	N	O	S	Se	0	0	0
			1178	743	205	224	2	4			
1	D	164	Total	C	N	O	S	Se	0	0	0
			1187	749	207	225	2	4			
1	E	164	Total	C	N	O	S	Se	0	0	0
			1187	749	207	225	2	4			
1	F	163	Total	C	N	O	S	Se	0	0	0
			1178	743	205	224	2	4			
1	G	166	Total	C	N	O	S	Se	0	0	0
			1203	758	210	228	2	5			
1	H	163	Total	C	N	O	S	Se	0	0	0
			1178	743	205	224	2	4			
1	I	165	Total	C	N	O	S	Se	0	0	0
			1195	753	209	227	2	4			
1	J	164	Total	C	N	O	S	Se	0	0	0
			1187	749	207	225	2	4			
1	K	164	Total	C	N	O	S	Se	0	0	0
			1187	749	207	225	2	4			
1	L	163	Total	C	N	O	S	Se	0	0	0
			1178	743	205	224	2	4			
1	M	164	Total	C	N	O	S	Se	0	0	0
			1187	749	207	225	2	4			
1	N	164	Total	C	N	O	S	Se	0	0	0
			1187	749	207	225	2	4			
1	O	164	Total	C	N	O	S	Se	0	0	0
			1187	749	207	225	2	4			
1	P	164	Total	C	N	O	S	Se	0	0	0
			1187	749	207	225	2	4			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	57	Total O 57 57	0	0
2	B	66	Total O 66 66	0	0
2	C	72	Total O 72 72	0	0
2	D	67	Total O 67 67	0	0
2	E	44	Total O 44 44	0	0
2	F	57	Total O 57 57	0	0
2	G	53	Total O 53 53	0	0
2	H	43	Total O 43 43	0	0
2	I	59	Total O 59 59	0	0
2	J	49	Total O 49 49	0	0
2	K	40	Total O 40 40	0	0
2	L	28	Total O 28 28	0	0
2	M	21	Total O 21 21	0	0
2	N	22	Total O 22 22	0	0
2	O	17	Total O 17 17	0	0
2	P	15	Total O 15 15	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

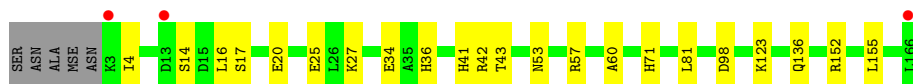
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain A: 



- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain B: 



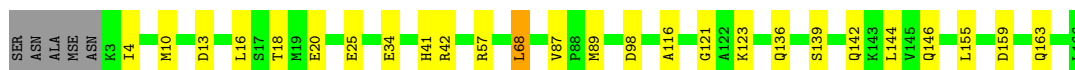
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain C: 



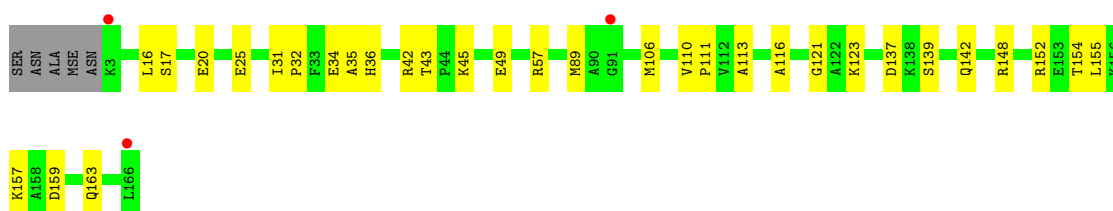
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain D: 



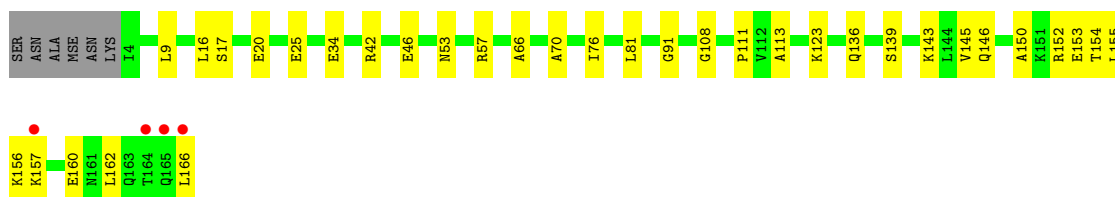
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain E: 



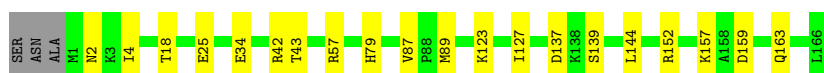
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain F:  2% 76% 20% .



- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain G:  86% 12% .



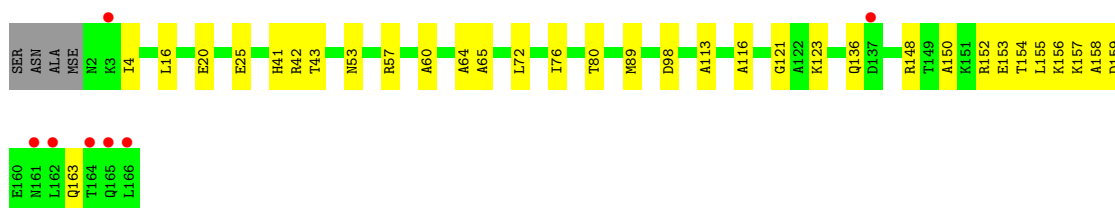
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain H:  4% 83% 13% .



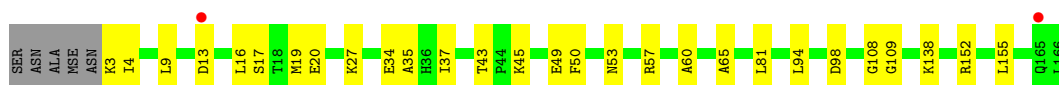
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain I:  4% 78% 20% .



- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain J:  80% 17% .



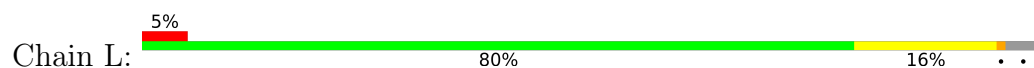
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

Chain K:  2% 78% 19% .

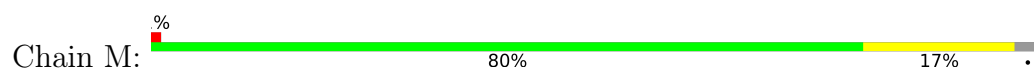




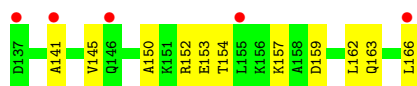
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit



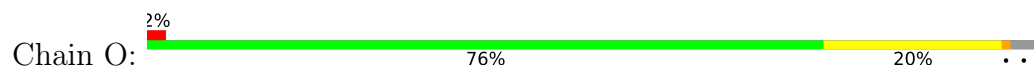
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit



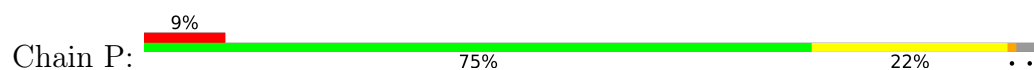
- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit

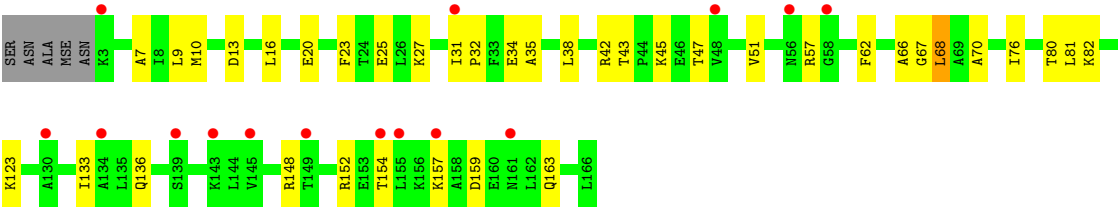


- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit



- Molecule 1: Phosphoribosylaminoimidazole carboxylase carboxyltransferase subunit





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.27Å 96.29Å 152.72Å 90.00° 91.95° 90.00°	Depositor
Resolution (Å)	40.80 – 2.20 40.80 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.7 (40.80-2.20) 95.6 (40.80-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.7 _650	Depositor
R, R_{free}	0.191 , 0.231 0.186 , 0.226	Depositor DCC
R_{free} test set	6139 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.343	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	19690	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% 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	0.37	0/1197	0.70	0/1612
1	B	0.37	0/1197	0.72	0/1612
1	C	0.35	0/1188	0.75	2/1601 (0.1%)
1	D	0.40	0/1197	0.75	1/1612 (0.1%)
1	E	0.35	0/1197	0.75	2/1612 (0.1%)
1	F	0.34	0/1188	0.71	0/1601
1	G	0.37	0/1213	0.74	0/1633
1	H	0.35	0/1188	0.72	0/1601
1	I	0.39	0/1205	0.72	0/1623
1	J	0.35	0/1197	0.70	0/1612
1	K	0.33	0/1197	0.72	0/1612
1	L	0.33	0/1188	0.72	0/1601
1	M	0.34	0/1197	0.71	0/1612
1	N	0.33	0/1197	0.71	0/1612
1	O	0.30	0/1197	0.76	1/1612 (0.1%)
1	P	0.32	0/1197	0.74	0/1612
All	All	0.35	0/19140	0.73	6/25780 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	31	ILE	CA-C-N	5.79	125.56	119.76
1	C	31	ILE	C-N-CA	5.79	125.56	119.76
1	E	106	MSE	CA-C-N	5.09	125.02	119.78
1	E	106	MSE	C-N-CA	5.09	125.02	119.78
1	D	68	LEU	CB-CA-C	-5.07	110.70	116.54
1	O	68	LEU	CB-CA-C	-5.05	110.77	116.63

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	1187	0	1241	17	0
1	B	1187	0	1241	19	0
1	C	1178	0	1228	14	0
1	D	1187	0	1241	19	0
1	E	1187	0	1241	26	0
1	F	1178	0	1228	27	0
1	G	1203	0	1259	21	0
1	H	1178	0	1228	16	0
1	I	1195	0	1247	30	0
1	J	1187	0	1241	20	0
1	K	1187	0	1241	28	0
1	L	1178	0	1228	28	0
1	M	1187	0	1241	26	0
1	N	1187	0	1241	33	0
1	O	1187	0	1241	31	0
1	P	1187	0	1241	25	0
2	A	57	0	0	2	0
2	B	66	0	0	7	0
2	C	72	0	0	1	0
2	D	67	0	0	2	0
2	E	44	0	0	5	0
2	F	57	0	0	4	0
2	G	53	0	0	3	0
2	H	43	0	0	0	0
2	I	59	0	0	7	0
2	J	49	0	0	2	0
2	K	40	0	0	1	0
2	L	28	0	0	2	0
2	M	21	0	0	1	0
2	N	22	0	0	2	0
2	O	17	0	0	4	0
2	P	15	0	0	0	0
All	All	19690	0	19828	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 8.

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:L:152:ARG:HG2	1:P:43:THR:HG22	1.45	0.98
1:N:152:ARG:HG2	1:O:43:THR:HG22	1.55	0.88
1:C:4:ILE:N	2:C:586:HOH:O	2.07	0.87
1:L:155:LEU:HD11	1:P:42:ARG:HD2	1.53	0.87
1:E:45:LYS:O	1:E:49:GLU:HG3	1.75	0.86
1:G:152:ARG:HG2	1:L:43:THR:HG22	1.57	0.85
1:I:136:GLN:HG2	2:I:617:HOH:O	1.74	0.84
1:G:89:MSE:HE2	1:G:89:MSE:HA	1.65	0.79
1:K:154:THR:OG1	2:K:647:HOH:O	2.02	0.78
1:B:43:THR:HG22	1:K:152:ARG:HG2	1.66	0.78
1:B:136:GLN:OE1	2:B:457:HOH:O	2.00	0.78
1:A:152:ARG:HG2	1:I:43:THR:HG22	1.67	0.77
1:F:16:LEU:O	1:F:20:GLU:HG2	1.85	0.76
2:B:457:HOH:O	1:D:136:GLN:HG2	1.85	0.76
1:L:136:GLN:HB3	1:O:136:GLN:CD	2.11	0.75
1:J:43:THR:HG22	1:O:152:ARG:HG2	1.69	0.75
1:P:25:GLU:HG3	1:P:123:LYS:HG3	1.70	0.73
1:A:106:MSE:HE1	1:D:41:HIS:CD2	2.24	0.73
1:B:36:HIS:NE2	2:B:441:HOH:O	2.21	0.73
1:L:136:GLN:NE2	2:L:698:HOH:O	2.19	0.72
1:E:43:THR:HG22	1:F:152:ARG:HG2	1.71	0.72
1:H:43:THR:HG22	1:J:152:ARG:HG2	1.72	0.72
1:L:81:LEU:HD22	1:O:144:LEU:HB3	1.71	0.72
1:O:89:MSE:HE2	1:O:89:MSE:HA	1.72	0.72
1:P:159:ASP:O	1:P:163:GLN:HG2	1.90	0.71
1:J:45:LYS:O	1:J:49:GLU:HG3	1.90	0.71
1:I:42:ARG:NE	2:I:483:HOH:O	2.15	0.71
1:L:42:ARG:NE	2:L:594:HOH:O	2.15	0.71
1:N:16:LEU:O	1:N:20:GLU:HG2	1.91	0.70
1:C:155:LEU:HD11	1:D:42:ARG:HD2	1.71	0.70
1:A:43:THR:HG22	1:B:152:ARG:HG2	1.74	0.70
1:K:16:LEU:O	1:K:20:GLU:HG2	1.91	0.70
1:M:46:GLU:OE1	1:M:46:GLU:N	2.25	0.69
1:A:155:LEU:HD11	1:I:42:ARG:HD2	1.73	0.69
1:D:16:LEU:O	1:D:20:GLU:HG2	1.92	0.69
1:E:42:ARG:HD2	1:F:155:LEU:HD11	1.75	0.69
1:L:45:LYS:O	1:L:49:GLU:HG3	1.93	0.69
1:D:89:MSE:HE2	1:D:89:MSE:HA	1.75	0.68
1:H:166:LEU:HD11	1:N:35:ALA:HB3	1.76	0.68
1:H:143:LYS:O	1:H:146:GLN:HG2	1.95	0.67
1:B:20:GLU:OE1	2:B:317:HOH:O	2.11	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:137:ASP:OD1	1:G:139:SER:HB3	1.95	0.67
1:L:34:GLU:OE2	1:L:57:ARG:NH2	2.27	0.67
1:D:155:LEU:HD11	1:F:42:ARG:HD2	1.75	0.66
1:P:23:PHE:O	1:P:27:LYS:HG3	1.96	0.66
1:I:25:GLU:HG3	1:I:123:LYS:HG3	1.76	0.66
1:G:43:THR:HG22	1:M:152:ARG:CG	2.26	0.65
1:E:36:HIS:HE1	2:E:599:HOH:O	1.80	0.65
1:G:43:THR:HG22	1:M:152:ARG:HG3	1.79	0.65
1:J:16:LEU:O	1:J:20:GLU:HG2	1.98	0.64
1:E:34:GLU:OE2	1:E:57:ARG:NH2	2.30	0.64
1:M:45:LYS:O	1:M:49:GLU:HG3	1.97	0.64
1:O:68:LEU:HA	1:O:98:ASP:OD2	1.98	0.64
1:L:166:LEU:HD11	1:P:35:ALA:HB3	1.79	0.64
1:K:25:GLU:HG3	1:K:123:LYS:HG3	1.81	0.63
1:M:89:MSE:HA	1:M:89:MSE:HE2	1.80	0.63
1:B:81:LEU:HD22	1:D:144:LEU:HB3	1.80	0.63
1:L:136:GLN:HB3	1:O:136:GLN:OE1	1.97	0.62
1:N:154:THR:HA	1:N:157:LYS:HE3	1.81	0.62
1:D:18:THR:HG21	1:D:89:MSE:HE1	1.81	0.62
1:L:111:PRO:HB3	1:O:124:ASN:HB3	1.81	0.62
1:N:116:ALA:O	1:N:121:GLY:HA3	1.99	0.62
1:H:155:LEU:HD11	1:N:42:ARG:HD2	1.82	0.61
1:D:142:GLN:NE2	2:D:668:HOH:O	2.32	0.61
1:G:144:LEU:HB3	1:J:81:LEU:HD22	1.81	0.61
1:J:3:LYS:HG3	1:J:4:ILE:H	1.66	0.61
1:K:90:ALA:HB2	1:K:117:ILE:HG22	1.81	0.61
1:I:154:THR:HA	1:I:157:LYS:HE3	1.83	0.61
1:O:137:ASP:OD1	1:O:139:SER:HB3	2.01	0.61
1:K:89:MSE:HA	1:K:89:MSE:HE2	1.83	0.60
1:C:42:ARG:HD2	1:E:155:LEU:HD11	1.82	0.60
1:K:154:THR:O	1:K:157:LYS:HE3	2.02	0.60
1:G:42:ARG:HD2	1:M:155:LEU:HD11	1.84	0.59
1:M:45:LYS:HE3	1:N:79:HIS:CD2	2.37	0.59
1:K:41:HIS:NE2	1:K:71:HIS:CD2	2.70	0.59
1:B:34:GLU:OE2	1:B:57:ARG:NH2	2.37	0.58
1:M:106:MSE:HE1	1:N:41:HIS:CD2	2.39	0.58
1:N:89:MSE:HE2	1:N:89:MSE:HA	1.85	0.57
1:M:18:THR:HG21	1:M:89:MSE:HE1	1.86	0.57
1:O:90:ALA:HB2	1:O:117:ILE:HG22	1.87	0.57
1:I:157:LYS:HG2	1:I:158:ALA:N	2.18	0.56
1:N:45:LYS:O	1:N:49:GLU:HG3	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:52:GLU:OE2	2:O:671:HOH:O	2.17	0.56
1:C:34:GLU:OE2	1:C:57:ARG:NH2	2.35	0.56
1:D:34:GLU:OE2	1:D:57:ARG:NH2	2.38	0.56
1:I:150:ALA:O	1:I:153:GLU:HB2	2.04	0.56
1:D:25:GLU:HG3	1:D:123:LYS:HG3	1.88	0.56
1:K:18:THR:HG21	1:K:89:MSE:HE1	1.86	0.56
1:G:152:ARG:NH2	2:G:470:HOH:O	2.39	0.56
1:P:76:ILE:O	1:P:80:THR:HG22	2.06	0.56
1:C:116:ALA:O	1:C:121:GLY:HA3	2.06	0.55
1:L:143:LYS:O	1:L:146:GLN:HG2	2.06	0.55
1:A:30:GLY:O	2:A:170:HOH:O	2.18	0.55
1:F:143:LYS:HA	1:F:146:GLN:HG2	1.88	0.55
1:M:4:ILE:HG23	1:M:60:ALA:HB2	1.87	0.55
1:E:16:LEU:HD11	1:F:162:LEU:HD13	1.88	0.55
1:M:34:GLU:OE2	1:M:57:ARG:NH2	2.39	0.55
1:P:133:ILE:O	1:P:136:GLN:HG2	2.07	0.55
1:I:53:ASN:O	1:I:57:ARG:HG3	2.06	0.54
1:K:34:GLU:OE2	1:K:57:ARG:NH2	2.39	0.54
1:P:148:ARG:O	1:P:152:ARG:HG3	2.07	0.54
1:I:155:LEU:HD11	1:K:42:ARG:HD2	1.90	0.54
1:M:137:ASP:OD1	1:M:139:SER:HB3	2.07	0.54
1:A:154:THR:HA	1:A:157:LYS:HE3	1.89	0.54
1:L:152:ARG:CG	1:P:43:THR:HG22	2.30	0.54
1:I:16:LEU:O	1:I:20:GLU:HG2	2.08	0.53
1:F:145:VAL:HG22	1:K:81:LEU:HD12	1.89	0.53
1:J:27:LYS:HD2	2:J:560:HOH:O	2.08	0.53
1:L:34:GLU:CD	1:L:57:ARG:HH22	2.16	0.53
1:J:94:LEU:HD12	1:J:98:ASP:HB3	1.91	0.53
1:L:16:LEU:O	1:L:20:GLU:HG2	2.09	0.53
1:N:159:ASP:O	1:N:163:GLN:HG2	2.08	0.53
1:M:47:THR:O	1:M:51:VAL:HG23	2.09	0.52
1:D:34:GLU:CD	1:D:57:ARG:HH22	2.17	0.52
1:M:9:LEU:HD22	1:M:76:ILE:HD11	1.91	0.52
1:P:38:LEU:HD13	1:P:47:THR:HA	1.90	0.52
1:I:152:ARG:CG	1:K:43:THR:HG22	2.40	0.52
1:N:76:ILE:O	1:N:80:THR:HG22	2.10	0.52
1:N:166:LEU:HD11	1:O:35:ALA:HB3	1.90	0.52
1:F:9:LEU:HD22	1:F:76:ILE:HD11	1.92	0.52
1:I:25:GLU:HG2	2:I:600:HOH:O	2.10	0.52
1:I:152:ARG:HG3	1:K:43:THR:HG22	1.92	0.52
1:K:41:HIS:NE2	1:K:71:HIS:HD2	2.06	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:53:ASN:OD1	1:J:57:ARG:HD3	2.10	0.51
1:O:19:MSE:SE	1:O:65:ALA:HB2	2.60	0.51
1:P:13:ASP:O	1:P:16:LEU:HB2	2.10	0.51
1:F:139:SER:O	1:F:143:LYS:HG3	2.10	0.51
1:M:16:LEU:O	1:M:20:GLU:HG2	2.11	0.51
1:O:79:HIS:CD2	1:P:45:LYS:HE3	2.46	0.51
1:F:66:ALA:HB3	1:F:70:ALA:HB1	1.92	0.50
1:O:9:LEU:N	1:O:9:LEU:HD12	2.25	0.50
1:D:68:LEU:HA	1:D:98:ASP:OD2	2.11	0.50
1:O:56:ASN:ND2	2:O:695:HOH:O	2.35	0.50
1:B:27:LYS:NZ	2:B:346:HOH:O	2.42	0.50
1:G:18:THR:HG21	1:G:89:MSE:HE1	1.94	0.50
1:F:17:SER:O	2:F:183:HOH:O	2.20	0.49
1:K:20:GLU:OE1	1:K:20:GLU:HA	2.10	0.49
1:N:46:GLU:H	1:N:46:GLU:CD	2.18	0.49
1:A:41:HIS:NE2	1:A:71:HIS:CD2	2.80	0.49
1:H:141:ALA:O	1:H:145:VAL:HG23	2.12	0.49
1:J:37:ILE:HD11	1:O:162:LEU:HD22	1.94	0.49
1:G:2:ASN:ND2	2:G:550:HOH:O	2.45	0.49
1:I:163:GLN:O	2:I:430:HOH:O	2.18	0.49
1:J:4:ILE:HG23	1:J:60:ALA:HB2	1.93	0.49
1:E:16:LEU:O	1:E:20:GLU:HG2	2.12	0.49
1:F:25:GLU:HG2	2:F:557:HOH:O	2.13	0.48
1:L:10:MSE:SE	1:L:16:LEU:HD23	2.63	0.48
1:A:152:ARG:NH2	2:A:515:HOH:O	2.45	0.48
1:E:35:ALA:HB3	1:F:166:LEU:HD11	1.95	0.48
1:E:137:ASP:OD1	1:E:139:SER:HB3	2.13	0.48
1:O:16:LEU:O	1:O:20:GLU:HG2	2.13	0.48
1:I:116:ALA:O	1:I:121:GLY:HA3	2.13	0.48
1:P:7:ALA:HB3	1:P:62:PHE:CD2	2.49	0.48
1:P:9:LEU:HD12	1:P:9:LEU:N	2.28	0.48
1:F:34:GLU:OE2	1:F:57:ARG:NH2	2.47	0.48
1:E:36:HIS:CE1	2:E:599:HOH:O	2.61	0.48
1:E:123:LYS:NZ	2:E:485:HOH:O	2.46	0.48
1:F:154:THR:HA	1:F:157:LYS:HE3	1.95	0.48
1:L:137:ASP:C	1:L:137:ASP:OD1	2.56	0.48
1:K:25:GLU:CG	1:K:123:LYS:HG3	2.43	0.48
1:O:149:THR:HA	1:O:152:ARG:NH1	2.29	0.48
1:N:134:ALA:C	1:N:136:GLN:H	2.21	0.47
1:P:66:ALA:HB3	1:P:70:ALA:HB1	1.95	0.47
1:F:91:GLY:O	2:F:589:HOH:O	2.20	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:34:GLU:OE2	1:P:57:ARG:NH2	2.47	0.47
1:L:98:ASP:OD1	1:L:98:ASP:N	2.47	0.47
1:N:150:ALA:O	1:N:153:GLU:HB2	2.15	0.47
1:P:51:VAL:HG21	1:P:76:ILE:HG12	1.94	0.47
1:D:116:ALA:O	1:D:121:GLY:HA3	2.14	0.47
1:F:81:LEU:HD22	1:K:144:LEU:HB3	1.97	0.47
1:H:124:ASN:HB3	1:M:111:PRO:HB3	1.97	0.47
1:A:149:THR:O	1:A:153:GLU:HG2	2.14	0.47
1:D:87:VAL:HG12	1:D:89:MSE:HE3	1.97	0.47
1:D:159:ASP:O	1:D:163:GLN:HG2	2.15	0.47
1:K:66:ALA:HB3	1:K:70:ALA:HB1	1.97	0.47
1:O:148:ARG:NH2	2:O:535:HOH:O	2.17	0.47
1:G:43:THR:HG22	1:M:152:ARG:HG2	1.95	0.47
1:J:35:ALA:HB3	1:O:166:LEU:HD11	1.97	0.46
1:A:45:LYS:O	1:A:49:GLU:HG3	2.15	0.46
1:G:34:GLU:CD	1:G:57:ARG:HH22	2.23	0.46
1:O:31:ILE:HA	1:O:32:PRO:HD3	1.71	0.46
1:C:31:ILE:HA	1:C:32:PRO:HD3	1.78	0.46
1:I:98:ASP:OD1	1:I:98:ASP:N	2.48	0.46
1:H:89:MSE:HA	1:H:89:MSE:HE2	1.97	0.46
1:L:90:ALA:HB2	1:L:117:ILE:CG2	2.45	0.46
1:N:152:ARG:CG	1:O:43:THR:HG22	2.38	0.46
1:O:9:LEU:HD22	1:O:76:ILE:HD11	1.97	0.46
1:E:89:MSE:HE2	1:E:89:MSE:HA	1.97	0.46
1:A:154:THR:HG23	1:A:157:LYS:CE	2.46	0.46
1:C:8:ILE:O	1:C:35:ALA:HA	2.16	0.45
1:C:151:LYS:HE3	1:D:42:ARG:HD3	1.98	0.45
1:E:31:ILE:HA	1:E:32:PRO:HD3	1.84	0.45
1:E:34:GLU:CD	1:E:57:ARG:HH22	2.24	0.45
1:H:42:ARG:HD2	1:J:155:LEU:HD11	1.96	0.45
1:E:154:THR:O	1:E:157:LYS:HE3	2.16	0.45
1:I:25:GLU:CG	1:I:123:LYS:HG3	2.45	0.45
1:J:138:LYS:HG3	2:J:495:HOH:O	2.15	0.45
1:N:18:THR:HG21	1:N:89:MSE:HE1	1.99	0.45
1:G:34:GLU:OE2	1:G:57:ARG:NH2	2.49	0.45
1:I:76:ILE:O	1:I:80:THR:HG22	2.16	0.45
1:H:116:ALA:O	1:H:121:GLY:HA3	2.16	0.45
1:L:154:THR:O	1:L:157:LYS:HB3	2.17	0.45
1:N:29:LEU:HD13	1:N:130:ALA:HB2	1.98	0.45
1:J:34:GLU:OE2	1:J:57:ARG:NH2	2.49	0.45
1:N:162:LEU:O	1:N:166:LEU:HG	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:GLU:CD	1:B:57:ARG:HH22	2.24	0.45
1:P:154:THR:O	1:P:157:LYS:HE3	2.16	0.45
1:B:41:HIS:NE2	1:B:71:HIS:CD2	2.84	0.45
1:I:159:ASP:O	1:I:163:GLN:HG2	2.17	0.45
1:P:31:ILE:HA	1:P:32:PRO:HD3	1.84	0.45
1:L:19:MSE:SE	1:L:65:ALA:HB2	2.67	0.45
1:L:154:THR:O	1:L:157:LYS:HE3	2.15	0.45
1:M:34:GLU:CD	1:M:57:ARG:HH22	2.25	0.44
1:O:7:ALA:HB3	1:O:62:PHE:CD2	2.53	0.44
1:E:110:VAL:HG13	1:I:148:ARG:NH2	2.32	0.44
1:C:108:GLY:O	1:I:42:ARG:NH2	2.45	0.44
1:G:25:GLU:HG3	1:G:123:LYS:HG3	1.99	0.44
1:G:89:MSE:HA	1:G:89:MSE:CE	2.43	0.44
1:N:90:ALA:HB2	1:N:117:ILE:CG2	2.47	0.44
1:J:108:GLY:O	1:L:42:ARG:NH2	2.48	0.44
1:K:143:LYS:HA	1:K:146:GLN:HG2	1.98	0.44
1:M:16:LEU:O	1:M:17:SER:C	2.60	0.44
1:E:111:PRO:O	2:E:173:HOH:O	2.21	0.44
1:M:8:ILE:O	1:M:35:ALA:HA	2.18	0.44
1:O:89:MSE:O	2:O:453:HOH:O	2.20	0.44
1:H:34:GLU:OE2	1:H:57:ARG:NH2	2.51	0.44
1:A:34:GLU:OE2	1:A:57:ARG:NH2	2.39	0.44
1:E:154:THR:HA	1:E:157:LYS:HE3	2.00	0.44
1:I:65:ALA:HB1	1:I:89:MSE:HE3	2.00	0.44
1:L:157:LYS:HB3	1:L:157:LYS:HE3	1.89	0.44
1:F:111:PRO:HB3	1:K:124:ASN:HB3	2.00	0.44
1:J:13:ASP:O	1:J:16:LEU:HB2	2.17	0.44
1:B:17:SER:HB2	2:B:474:HOH:O	2.17	0.43
1:B:152:ARG:NE	2:B:271:HOH:O	2.51	0.43
1:F:136:GLN:NE2	1:K:136:GLN:HB3	2.33	0.43
1:K:16:LEU:O	1:K:17:SER:C	2.58	0.43
1:F:46:GLU:H	1:F:46:GLU:CD	2.27	0.43
1:I:136:GLN:NE2	2:I:617:HOH:O	2.44	0.43
1:A:154:THR:O	1:A:157:LYS:HB3	2.17	0.43
1:E:142:GLN:OE1	2:E:514:HOH:O	2.21	0.43
1:G:152:ARG:CG	1:L:43:THR:HG22	2.40	0.43
1:N:29:LEU:HD13	1:N:130:ALA:CB	2.49	0.43
1:O:96:GLY:HA3	1:O:117:ILE:HD12	2.01	0.43
1:L:111:PRO:HB3	1:O:124:ASN:CB	2.48	0.43
1:N:25:GLU:HG3	1:N:123:LYS:HG3	2.01	0.43
1:A:76:ILE:O	1:A:80:THR:HG22	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:45:LYS:HE3	1:N:79:HIS:HD2	1.82	0.43
1:N:30:GLY:HA2	2:N:490:HOH:O	2.18	0.43
1:B:16:LEU:O	1:B:20:GLU:HG2	2.19	0.43
1:C:13:ASP:O	1:C:16:LEU:HB2	2.19	0.43
1:F:113:ALA:HB1	1:K:113:ALA:HB1	2.00	0.43
1:I:4:ILE:HG23	1:I:60:ALA:HB2	2.01	0.43
1:N:135:LEU:HD21	1:P:82:LYS:HD2	2.01	0.43
1:C:106:MSE:HE1	1:I:41:HIS:CD2	2.54	0.43
1:M:25:GLU:HG2	2:M:437:HOH:O	2.19	0.43
1:O:45:LYS:O	1:O:49:GLU:HG3	2.19	0.43
1:G:127:ILE:HD13	1:J:109:GLY:HA2	2.00	0.42
1:G:159:ASP:O	1:G:163:GLN:HG2	2.19	0.42
1:H:162:LEU:O	1:H:162:LEU:HD12	2.19	0.42
1:K:116:ALA:O	1:K:121:GLY:HA3	2.19	0.42
1:F:136:GLN:CD	1:K:136:GLN:HB3	2.43	0.42
1:E:113:ALA:HB1	1:I:113:ALA:HB1	2.00	0.42
1:G:87:VAL:HG12	1:G:89:MSE:HE3	2.02	0.42
1:A:42:ARG:HD2	1:B:155:LEU:HD11	2.00	0.42
1:E:16:LEU:O	1:E:17:SER:C	2.62	0.42
1:F:156:LYS:O	1:F:160:GLU:HG3	2.19	0.42
1:J:19:MSE:SE	1:J:65:ALA:HB2	2.69	0.42
1:P:16:LEU:O	1:P:20:GLU:HG2	2.18	0.42
1:B:25:GLU:OE2	1:B:123:LYS:HE2	2.20	0.42
1:I:156:LYS:HB2	2:I:693:HOH:O	2.18	0.42
1:F:150:ALA:O	1:F:153:GLU:HB2	2.20	0.42
1:P:67:GLY:O	1:P:68:LEU:C	2.61	0.42
1:A:9:LEU:HD22	1:A:76:ILE:HD11	2.01	0.42
1:I:64:ALA:HB1	1:I:72:LEU:HG	2.02	0.42
1:O:144:LEU:HD23	1:O:144:LEU:HA	1.83	0.42
1:E:25:GLU:HG3	1:E:123:LYS:HG3	2.01	0.42
1:J:9:LEU:HD11	1:J:50:PHE:HE2	1.85	0.42
1:H:127:ILE:O	1:H:131:GLN:HG3	2.20	0.41
1:P:10:MSE:HE1	1:P:23:PHE:CZ	2.54	0.41
1:D:10:MSE:SE	1:D:16:LEU:HG	2.70	0.41
1:H:106:MSE:HA	1:H:107:PRO:HD3	1.92	0.41
1:H:113:ALA:HB1	1:M:113:ALA:HB1	2.02	0.41
1:N:44:PRO:HD2	2:N:434:HOH:O	2.20	0.41
1:B:42:ARG:NH2	1:F:108:GLY:O	2.53	0.41
1:E:148:ARG:O	1:E:152:ARG:HG3	2.20	0.41
1:F:53:ASN:ND2	2:F:170:HOH:O	2.53	0.41
1:K:98:ASP:OD1	1:K:99:ALA:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:134:ALA:C	1:N:136:GLN:N	2.77	0.41
1:E:159:ASP:O	1:E:163:GLN:HG2	2.20	0.41
1:G:79:HIS:HE1	2:G:467:HOH:O	2.04	0.41
1:G:157:LYS:HE3	1:G:157:LYS:HB3	1.82	0.41
1:H:108:GLY:HA3	1:M:123:LYS:HD2	2.02	0.41
1:I:42:ARG:CD	2:I:483:HOH:O	2.66	0.41
1:F:25:GLU:OE2	1:F:123:LYS:HE2	2.21	0.41
1:N:41:HIS:NE2	1:N:71:HIS:CD2	2.89	0.41
1:P:81:LEU:HD23	1:P:81:LEU:HA	1.94	0.41
1:N:43:THR:HB	1:N:46:GLU:OE1	2.21	0.41
1:B:98:ASP:OD1	1:B:98:ASP:N	2.53	0.41
1:C:16:LEU:O	1:C:20:GLU:HG2	2.20	0.41
1:H:155:LEU:HD11	1:N:42:ARG:CD	2.48	0.41
1:N:141:ALA:O	1:N:145:VAL:HG23	2.20	0.41
1:L:4:ILE:HG23	1:L:60:ALA:HB2	2.03	0.40
1:B:4:ILE:HG23	1:B:60:ALA:HB2	2.03	0.40
1:E:116:ALA:O	1:E:121:GLY:HA3	2.21	0.40
1:N:31:ILE:HA	1:N:32:PRO:HD3	1.82	0.40
1:B:53:ASN:O	1:B:57:ARG:HG3	2.21	0.40
1:D:146:GLN:HB3	2:D:627:HOH:O	2.21	0.40
1:M:81:LEU:HD23	1:M:81:LEU:HA	1.96	0.40
1:O:82:LYS:HD3	1:O:82:LYS:HA	1.86	0.40
1:C:87:VAL:HA	1:C:88:PRO:HD3	1.90	0.40
1:K:90:ALA:HB2	1:K:117:ILE:CG2	2.50	0.40
1:A:124:ASN:HB3	1:C:111:PRO:HB3	2.04	0.40
1:M:98:ASP:OD1	1:M:98:ASP:N	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	162/169 (96%)	159 (98%)	3 (2%)	0	100	100
1	B	162/169 (96%)	162 (100%)	0	0	100	100
1	C	161/169 (95%)	157 (98%)	4 (2%)	0	100	100
1	D	162/169 (96%)	159 (98%)	3 (2%)	0	100	100
1	E	162/169 (96%)	159 (98%)	3 (2%)	0	100	100
1	F	161/169 (95%)	161 (100%)	0	0	100	100
1	G	164/169 (97%)	161 (98%)	3 (2%)	0	100	100
1	H	161/169 (95%)	158 (98%)	3 (2%)	0	100	100
1	I	163/169 (96%)	158 (97%)	5 (3%)	0	100	100
1	J	162/169 (96%)	160 (99%)	2 (1%)	0	100	100
1	K	162/169 (96%)	159 (98%)	3 (2%)	0	100	100
1	L	161/169 (95%)	158 (98%)	3 (2%)	0	100	100
1	M	162/169 (96%)	158 (98%)	4 (2%)	0	100	100
1	N	162/169 (96%)	155 (96%)	7 (4%)	0	100	100
1	O	162/169 (96%)	157 (97%)	5 (3%)	0	100	100
1	P	162/169 (96%)	155 (96%)	6 (4%)	1 (1%)	21	23
All	All	2591/2704 (96%)	2536 (98%)	54 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	P	68	LEU

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	122/121 (101%)	122 (100%)	0	100	100
1	B	122/121 (101%)	121 (99%)	1 (1%)	73	85
1	C	121/121 (100%)	119 (98%)	2 (2%)	53	69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	122/121 (101%)	119 (98%)	3 (2%)	42	56
1	E	122/121 (101%)	122 (100%)	0	100	100
1	F	121/121 (100%)	121 (100%)	0	100	100
1	G	124/121 (102%)	123 (99%)	1 (1%)	73	85
1	H	121/121 (100%)	120 (99%)	1 (1%)	73	85
1	I	123/121 (102%)	123 (100%)	0	100	100
1	J	122/121 (101%)	121 (99%)	1 (1%)	73	85
1	K	122/121 (101%)	121 (99%)	1 (1%)	73	85
1	L	121/121 (100%)	119 (98%)	2 (2%)	53	69
1	M	122/121 (101%)	122 (100%)	0	100	100
1	N	122/121 (101%)	121 (99%)	1 (1%)	73	85
1	O	122/121 (101%)	120 (98%)	2 (2%)	55	71
1	P	122/121 (101%)	122 (100%)	0	100	100
All	All	1951/1936 (101%)	1936 (99%)	15 (1%)	73	85

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

Mol	Chain	Res	Type
1	B	14	SER
1	C	102	SER
1	C	154	THR
1	D	4	ILE
1	D	13	ASP
1	D	139	SER
1	G	4	ILE
1	H	142	GLN
1	J	17	SER
1	K	139	SER
1	L	104	VAL
1	L	154	THR
1	N	112	VAL
1	O	4	ILE
1	O	117	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	71	HIS
1	A	136	GLN
1	A	146	GLN
1	B	71	HIS
1	D	71	HIS
1	E	56	ASN
1	E	71	HIS
1	F	36	HIS
1	F	71	HIS
1	F	136	GLN
1	G	2	ASN
1	G	147	GLN
1	G	165	GLN
1	I	2	ASN
1	I	36	HIS
1	I	71	HIS
1	J	56	ASN
1	K	71	HIS
1	K	147	GLN
1	L	36	HIS
1	L	163	GLN
1	M	71	HIS
1	M	136	GLN
1	M	161	ASN
1	N	71	HIS
1	O	36	HIS
1	O	146	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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	160/169 (94%)	-0.19	0 100 100	16, 29, 45, 71	0
1	B	160/169 (94%)	-0.27	3 (1%) 66 63	17, 28, 46, 66	0
1	C	159/169 (94%)	-0.25	0 100 100	17, 30, 44, 55	0
1	D	160/169 (94%)	-0.33	0 100 100	15, 26, 43, 61	0
1	E	160/169 (94%)	-0.02	3 (1%) 66 63	20, 35, 51, 70	0
1	F	159/169 (94%)	-0.10	4 (2%) 58 55	19, 32, 54, 71	0
1	G	161/169 (95%)	-0.11	0 100 100	20, 33, 49, 57	0
1	H	159/169 (94%)	0.06	6 (3%) 44 41	23, 34, 61, 73	0
1	I	161/169 (95%)	0.06	7 (4%) 40 36	20, 30, 62, 76	0
1	J	160/169 (94%)	-0.05	2 (1%) 75 73	20, 33, 52, 68	0
1	K	160/169 (94%)	0.12	4 (2%) 58 55	24, 37, 55, 76	0
1	L	159/169 (94%)	0.37	8 (5%) 34 31	27, 41, 72, 87	0
1	M	160/169 (94%)	0.20	2 (1%) 75 73	26, 41, 58, 73	0
1	N	160/169 (94%)	0.64	8 (5%) 34 31	32, 49, 69, 80	0
1	O	160/169 (94%)	0.43	4 (2%) 58 55	33, 50, 66, 77	0
1	P	160/169 (94%)	0.96	15 (9%) 14 11	37, 53, 70, 81	0
All	All	2558/2704 (94%)	0.10	66 (2%) 57 54	15, 36, 63, 87	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	164	THR	4.2
1	N	91	GLY	4.0
1	H	91	GLY	3.7
1	F	166	LEU	3.6
1	I	162	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
1	H	153	GLU	3.5
1	P	58	GLY	3.4
1	N	137	ASP	3.4
1	L	157	LYS	3.3
1	L	166	LEU	3.2
1	P	139	SER	3.1
1	P	155	LEU	2.9
1	N	166	LEU	2.9
1	L	165	GLN	2.8
1	P	145	VAL	2.8
1	I	166	LEU	2.8
1	K	157	LYS	2.7
1	H	17	SER	2.7
1	J	165	GLN	2.6
1	L	160	GLU	2.6
1	E	91	GLY	2.6
1	L	109	GLY	2.6
1	H	155	LEU	2.6
1	I	137	ASP	2.5
1	I	161	ASN	2.5
1	P	154	THR	2.5
1	O	17	SER	2.5
1	L	162	LEU	2.5
1	P	31	ILE	2.4
1	K	17	SER	2.4
1	P	143	LYS	2.3
1	P	149	THR	2.3
1	F	157	LYS	2.3
1	J	13	ASP	2.3
1	N	141	ALA	2.3
1	E	3	LYS	2.3
1	P	130	ALA	2.3
1	O	4	ILE	2.3
1	L	154	THR	2.2
1	E	166	LEU	2.2
1	N	155	LEU	2.2
1	L	49	GLU	2.2
1	P	48	VAL	2.2
1	M	91	GLY	2.2
1	I	3	LYS	2.2
1	I	165	GLN	2.2
1	P	134	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	N	93	SER	2.2
1	N	13	ASP	2.1
1	P	3	LYS	2.1
1	P	157	LYS	2.1
1	B	166	LEU	2.1
1	M	166	LEU	2.1
1	N	146	GLN	2.1
1	O	5	PHE	2.1
1	K	49	GLU	2.1
1	B	13	ASP	2.1
1	F	164	THR	2.0
1	H	164	THR	2.0
1	O	12	SER	2.0
1	F	165	GLN	2.0
1	H	4	ILE	2.0
1	K	91	GLY	2.0
1	B	3	LYS	2.0
1	P	56	ASN	2.0
1	P	161	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.