



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 03:06 PM UTC

PDB ID : 3L7O / pdb_00003l7o
Title : Crystal structure of Ribose-5-phosphate isomerase A from streptococcus mutants UA159
Authors : Fan, X.X.; Wang, K.T.; Su, X.D.
Deposited on : 2009-12-29
Resolution : 1.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

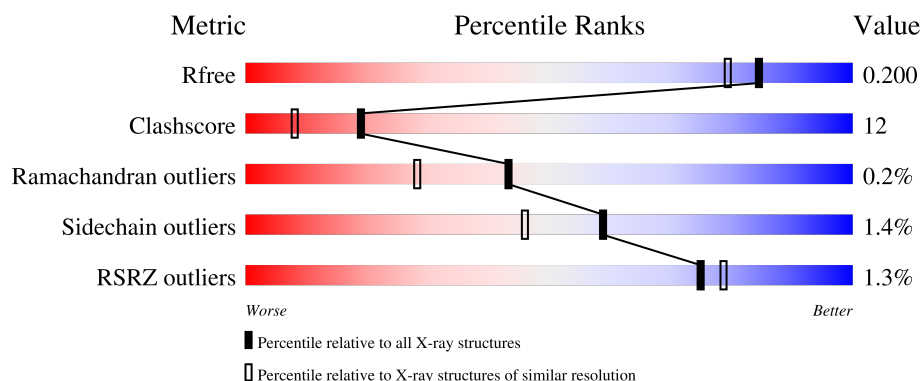
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

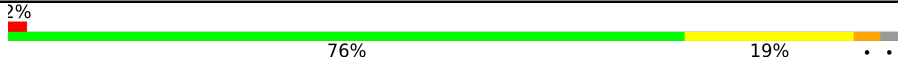
The reported resolution of this entry is 1.70 Å.

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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	225	
1	B	225	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4023 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 Ribose-5-phosphate isomerase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1697	1078	285	328	6			
1	B	224	Total	C	N	O	S	0	0	0
			1715	1087	288	334	6			

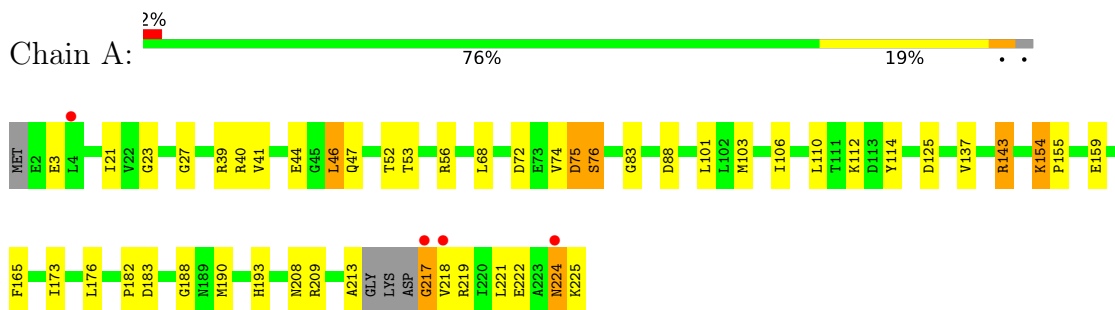
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	322	Total	O	0	0
			322	322		
2	B	289	Total	O	0	0
			289	289		

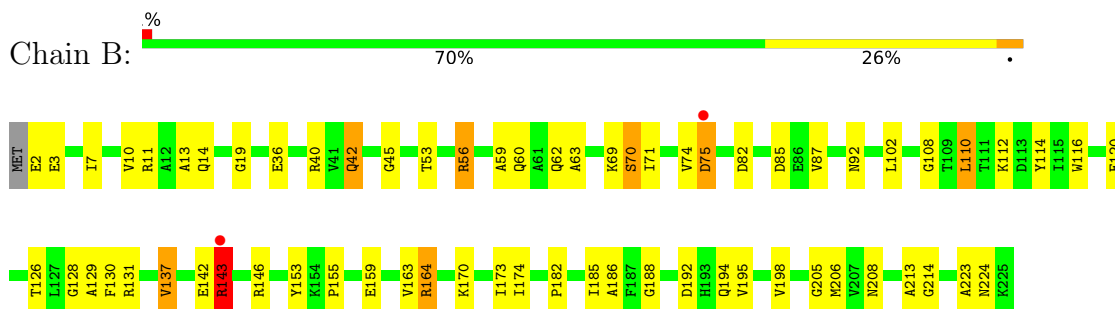
3 Residue-property plots [i](#)

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: Ribose-5-phosphate isomerase A



- Molecule 1: Ribose-5-phosphate isomerase A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	48.97Å 101.09Å 98.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.80 – 1.70 32.80 – 1.70	Depositor EDS
% Data completeness (in resolution range)	96.1 (32.80-1.70) 97.9 (32.80-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.197 , 0.214 (Not available) , 0.200	Depositor DCC
R_{free} test set	5029 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 34.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4023	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.72% 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 ⓘ

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	2.00	21/1722 (1.2%)	1.53	9/2325 (0.4%)
1	B	2.13	55/1741 (3.2%)	1.51	7/2352 (0.3%)
All	All	2.06	76/3463 (2.2%)	1.52	16/4677 (0.3%)

All (76) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	137	VAL	CA-CB	12.57	1.67	1.54
1	B	13	ALA	CA-CB	-8.30	1.38	1.53
1	B	71	ILE	CA-C	7.81	1.63	1.52
1	B	2	GLU	N-CA	7.29	1.60	1.46
1	A	165	PHE	CA-C	7.19	1.62	1.52
1	B	206	MET	C-O	7.03	1.32	1.23
1	B	223	ALA	C-O	6.99	1.33	1.24
1	B	195	VAL	CA-CB	6.94	1.62	1.54
1	A	173	ILE	C-O	6.80	1.30	1.24
1	B	188	GLY	N-CA	6.80	1.54	1.45
1	B	182	PRO	CA-C	6.79	1.63	1.52
1	B	170	LYS	N-CA	6.78	1.56	1.46
1	A	155	PRO	C-O	6.72	1.31	1.23
1	B	63	ALA	CA-CB	6.65	1.63	1.53
1	B	10	VAL	CA-CB	6.58	1.62	1.54
1	B	137	VAL	C-O	6.54	1.30	1.24
1	B	143	ARG	CB-CG	6.45	1.71	1.52
1	B	74	VAL	CA-CB	6.36	1.65	1.55
1	B	10	VAL	N-CA	6.29	1.53	1.46
1	B	214	GLY	N-CA	6.24	1.50	1.45
1	A	222	GLU	C-O	6.22	1.31	1.24
1	B	194	GLN	CD-OE1	6.22	1.35	1.23
1	B	102	LEU	C-O	6.13	1.31	1.24
1	A	154	LYS	C-O	6.13	1.26	1.23
1	B	194	GLN	CG-CD	6.12	1.67	1.52
1	B	45	GLY	C-O	6.12	1.32	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	63	ALA	N-CA	6.10	1.53	1.46
1	A	217	GLY	C-O	6.08	1.32	1.23
1	B	53	THR	CA-CB	6.03	1.62	1.53
1	A	183	ASP	N-CA	6.03	1.54	1.46
1	B	59	ALA	N-CA	5.99	1.53	1.46
1	A	23	GLY	C-O	5.96	1.30	1.23
1	A	106	ILE	CA-C	5.94	1.60	1.52
1	B	185	ILE	CA-C	5.88	1.60	1.52
1	B	74	VAL	N-CA	5.87	1.52	1.46
1	A	143	ARG	CB-CG	-5.83	1.34	1.52
1	B	62	GLN	N-CA	5.83	1.53	1.46
1	A	83	GLY	C-O	5.78	1.30	1.23
1	B	195	VAL	C-O	5.77	1.30	1.24
1	A	221	LEU	CA-C	5.71	1.59	1.52
1	A	52	THR	N-CA	5.70	1.52	1.45
1	B	7	ILE	CA-C	5.63	1.59	1.52
1	A	53	THR	C-O	5.58	1.31	1.23
1	B	186	ALA	CA-CB	5.55	1.61	1.53
1	A	154	LYS	CA-CB	5.52	1.57	1.52
1	B	92	ASN	C-O	5.51	1.30	1.23
1	A	46	LEU	C-O	5.47	1.30	1.23
1	B	126	THR	CA-CB	5.41	1.63	1.53
1	B	155	PRO	C-O	5.40	1.30	1.23
1	A	159	GLU	C-O	5.40	1.30	1.23
1	B	223	ALA	CA-CB	-5.39	1.44	1.53
1	B	60	GLN	N-CA	5.37	1.52	1.46
1	B	198	VAL	CA-C	5.36	1.59	1.52
1	B	19	GLY	C-O	-5.34	1.16	1.24
1	B	173	ILE	CA-CB	5.34	1.61	1.54
1	B	128	GLY	N-CA	5.31	1.53	1.45
1	B	75	ASP	CB-CG	5.30	1.65	1.52
1	B	153	TYR	CD1-CE1	5.30	1.54	1.38
1	A	21	ILE	C-O	5.27	1.30	1.24
1	A	188	GLY	C-O	-5.26	1.17	1.23
1	B	174	ILE	CA-C	5.25	1.59	1.52
1	B	126	THR	C-O	5.22	1.29	1.23
1	A	137	VAL	CA-C	5.20	1.58	1.52
1	B	130	PHE	CA-C	5.14	1.59	1.52
1	B	85	ASP	C-O	5.14	1.30	1.24
1	B	42	GLN	CB-CG	5.13	1.67	1.52
1	B	108	GLY	C-O	-5.11	1.17	1.23
1	A	176	LEU	C-O	5.09	1.30	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	129	ALA	CA-CB	5.09	1.61	1.53
1	B	224	ASN	N-CA	5.09	1.52	1.46
1	B	163	VAL	CA-CB	5.07	1.60	1.54
1	B	120	GLU	CA-CB	5.05	1.61	1.53
1	B	11	ARG	CG-CD	5.04	1.67	1.52
1	B	87	VAL	N-CA	5.04	1.52	1.46
1	B	213	ALA	CA-CB	5.04	1.60	1.53
1	B	56	ARG	CZ-NH2	5.03	1.40	1.33

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	143	ARG	CG-CD-NE	6.62	126.58	112.00
1	A	76	SER	N-CA-C	-6.38	98.83	108.52
1	A	53	THR	N-CA-C	-6.20	105.36	113.17
1	A	75	ASP	N-CA-C	5.94	119.72	112.23
1	B	131	ARG	NE-CZ-NH2	-5.87	113.92	119.20
1	B	11	ARG	NE-CZ-NH2	5.86	124.48	119.20
1	B	74	VAL	CA-C-N	-5.45	111.50	121.52
1	B	74	VAL	C-N-CA	-5.45	111.50	121.52
1	A	27	GLY	N-CA-C	-5.37	103.06	111.00
1	A	3	GLU	N-CA-C	-5.27	105.53	111.28
1	A	74	VAL	N-CA-CB	-5.25	105.17	112.52
1	A	88	ASP	CB-CA-C	5.21	117.77	109.96
1	B	110	LEU	N-CA-C	-5.14	106.59	112.92
1	A	39	ARG	N-CA-C	5.12	116.86	111.28
1	B	75	ASP	N-CA-CB	5.03	117.87	110.33
1	A	219	ARG	CA-CB-CG	5.02	124.14	114.10

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	1697	0	1693	38	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1715	0	1706	44	0
2	A	322	0	0	25	1
2	B	289	0	0	34	2
All	All	4023	0	3399	80	2

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 12.

All (80) 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:137:VAL:HG22	2:B:687:HOH:O	1.29	1.33
1:B:56:ARG:HD2	2:B:603:HOH:O	1.16	1.30
1:A:41:VAL:HA	2:A:664:HOH:O	1.21	1.30
1:B:142:GLU:HG3	2:B:662:HOH:O	1.24	1.28
1:A:46:LEU:HB3	2:A:664:HOH:O	1.30	1.22
1:B:164:ARG:CD	2:B:686:HOH:O	1.91	1.17
2:A:616:HOH:O	1:B:70:SER:HA	1.47	1.12
1:A:190:MET:HG2	2:A:663:HOH:O	1.50	1.10
1:B:164:ARG:CG	2:B:686:HOH:O	1.95	1.10
1:A:44:GLU:HG2	2:A:624:HOH:O	1.52	1.08
1:B:164:ARG:HG2	2:B:686:HOH:O	1.50	1.07
1:A:75:ASP:HB2	2:A:584:HOH:O	1.60	1.02
1:A:110:LEU:HB2	2:B:429:HOH:O	1.60	1.00
1:B:36:GLU:CD	2:B:683:HOH:O	2.05	0.99
1:A:182:PRO:HD2	2:A:629:HOH:O	1.66	0.93
1:A:213:ALA:C	1:A:217:GLY:N	2.28	0.92
1:B:42:GLN:HG2	2:B:585:HOH:O	1.75	0.85
1:B:114:TYR:H	1:B:208:ASN:HD22	1.20	0.85
1:A:68:LEU:HD22	2:A:672:HOH:O	1.79	0.83
1:A:114:TYR:H	1:A:208:ASN:HD22	1.20	0.82
1:B:75:ASP:HB3	2:B:668:HOH:O	1.80	0.82
1:B:110:LEU:HD22	2:B:679:HOH:O	1.81	0.81
1:A:224:ASN:O	1:A:225:LYS:C	2.27	0.76
1:B:14:GLN:HG3	2:B:341:HOH:O	1.85	0.76
1:A:41:VAL:CA	2:A:664:HOH:O	1.99	0.75
1:B:142:GLU:CG	2:B:662:HOH:O	1.98	0.75
1:B:56:ARG:CD	2:B:603:HOH:O	1.94	0.73
1:B:143:ARG:HD2	1:B:146:ARG:HE	1.54	0.73
1:A:72:ASP:OD2	2:A:570:HOH:O	2.06	0.72
1:B:36:GLU:OE2	2:B:683:HOH:O	2.01	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:HIS:HD2	1:B:192:ASP:OD2	1.75	0.69
1:A:218:VAL:HG22	2:A:281:HOH:O	1.94	0.68
1:B:75:ASP:O	2:B:679:HOH:O	2.12	0.68
1:B:164:ARG:HD2	2:B:686:HOH:O	1.73	0.66
1:A:103:MET:HE1	2:A:626:HOH:O	1.96	0.66
1:B:3:GLU:HG2	2:B:317:HOH:O	1.96	0.65
1:B:143:ARG:CD	1:B:146:ARG:HE	2.09	0.65
1:B:205:GLY:HA2	2:B:645:HOH:O	1.98	0.64
1:A:154:LYS:HD3	2:A:550:HOH:O	1.99	0.63
1:A:110:LEU:HD22	1:B:143:ARG:HG3	1.82	0.61
1:B:69:LYS:HE3	2:B:704:HOH:O	1.99	0.61
1:A:72:ASP:CG	2:A:570:HOH:O	2.43	0.61
1:A:76:SER:CB	1:A:112:LYS:NZ	2.64	0.61
1:B:142:GLU:CD	2:B:662:HOH:O	2.33	0.60
1:A:47:GLN:HG3	2:A:638:HOH:O	2.01	0.59
1:B:40:ARG:HG3	2:B:683:HOH:O	2.04	0.57
1:A:114:TYR:H	1:A:208:ASN:ND2	1.98	0.56
1:B:143:ARG:CD	1:B:146:ARG:HH21	2.18	0.56
1:B:164:ARG:NH1	2:B:686:HOH:O	2.39	0.55
1:B:69:LYS:CE	2:B:704:HOH:O	2.52	0.55
1:B:143:ARG:HD2	1:B:146:ARG:NE	2.23	0.54
1:A:76:SER:CB	1:A:112:LYS:HZ1	2.21	0.54
1:A:41:VAL:N	2:A:664:HOH:O	2.34	0.53
1:B:36:GLU:CG	2:B:683:HOH:O	2.49	0.53
1:B:143:ARG:HD3	1:B:146:ARG:HH21	1.74	0.53
1:B:112:LYS:CE	2:B:614:HOH:O	2.57	0.52
1:B:114:TYR:H	1:B:208:ASN:ND2	2.01	0.51
1:A:143:ARG:HD3	2:B:577:HOH:O	2.11	0.51
1:A:209:ARG:CZ	2:A:677:HOH:O	2.58	0.50
1:B:164:ARG:CZ	2:B:686:HOH:O	2.59	0.49
1:B:36:GLU:HG3	2:B:683:HOH:O	2.11	0.49
1:B:112:LYS:HD2	2:B:614:HOH:O	2.13	0.48
1:B:112:LYS:CD	2:B:614:HOH:O	2.62	0.48
1:A:209:ARG:CZ	1:A:209:ARG:HB2	2.44	0.48
1:A:46:LEU:N	2:A:664:HOH:O	2.47	0.47
1:A:75:ASP:CB	2:A:584:HOH:O	2.39	0.47
1:A:40:ARG:C	2:A:664:HOH:O	2.57	0.45
1:B:82:ASP:O	1:B:116:TRP:HA	2.16	0.45
1:A:101:LEU:H	1:A:101:LEU:HD23	1.81	0.45
1:A:209:ARG:NH2	2:A:677:HOH:O	2.50	0.45
1:A:110:LEU:HG	1:A:110:LEU:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:ARG:HD2	1:B:143:ARG:HA	1.68	0.43
1:B:40:ARG:CG	2:B:683:HOH:O	2.65	0.43
1:A:46:LEU:CB	2:A:664:HOH:O	2.15	0.43
1:B:164:ARG:NE	2:B:686:HOH:O	2.31	0.43
1:A:225:LYS:HE3	1:A:225:LYS:HB3	1.82	0.42
1:A:46:LEU:CA	2:A:664:HOH:O	2.59	0.41
1:A:209:ARG:NH1	2:A:677:HOH:O	2.54	0.41
1:B:112:LYS:HE3	2:B:614:HOH:O	2.20	0.41
1:A:125:ASP:HB2	2:A:580:HOH:O	2.20	0.40

All (2) 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	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:LYS:CA	2:B:341:HOH:O[3_645]	1.94	0.26
2:A:255:HOH:O	2:B:556:HOH:O[4_555]	2.14	0.06

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	217/225 (96%)	214 (99%)	2 (1%)	1 (0%)	24	12
1	B	222/225 (99%)	219 (99%)	3 (1%)	0	100	100
All	All	439/450 (98%)	433 (99%)	5 (1%)	1 (0%)	43	28

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	224	ASN

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	170/184 (92%)	169 (99%)	1 (1%)	78	72
1	B	182/184 (99%)	178 (98%)	4 (2%)	45	29
All	All	352/368 (96%)	347 (99%)	5 (1%)	59	46

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ARG
1	B	70	SER
1	B	143	ARG
1	B	159	GLU
1	B	164	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	GLN
1	A	193	HIS
1	A	194	GLN
1	A	208	ASN
1	B	90	ASN
1	B	204	ASN
1	B	208	ASN
1	B	224	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	221/225 (98%)	-0.36	4 (1%) 67 72	9, 14, 27, 41	0
1	B	224/225 (99%)	-0.28	2 (0%) 81 83	10, 15, 31, 46	0
All	All	445/450 (98%)	-0.32	6 (1%) 75 79	9, 14, 30, 46	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	217	GLY	3.6
1	B	75	ASP	3.1
1	A	224	ASN	2.5
1	A	218	VAL	2.5
1	B	143	ARG	2.3
1	A	4	LEU	2.1

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.