Full wwPDB X-ray Structure Validation Report

Feb 1, 2016 – 10:11 AM GMT

PDB ID	:	3L7O
Title	:	Crystal structure of Ribose-5-phosphate isomerase A from streptococcus mutans 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

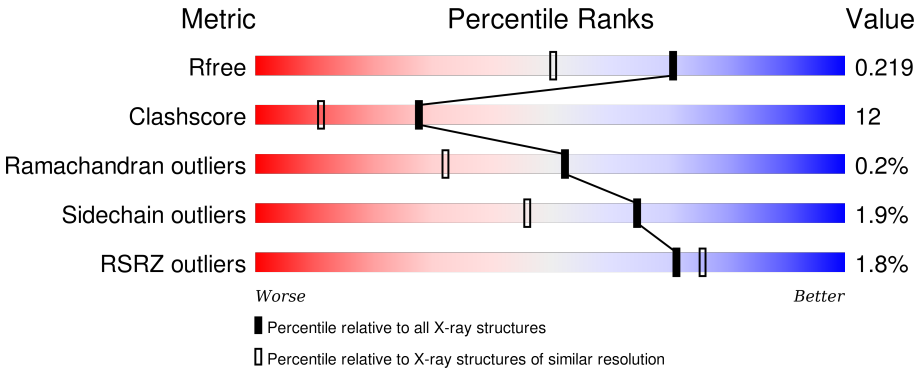
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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}	91344	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (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 on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=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	2% 78% 18%
1	B	225	6% 74% 20%

2 Entry composition (i)

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.

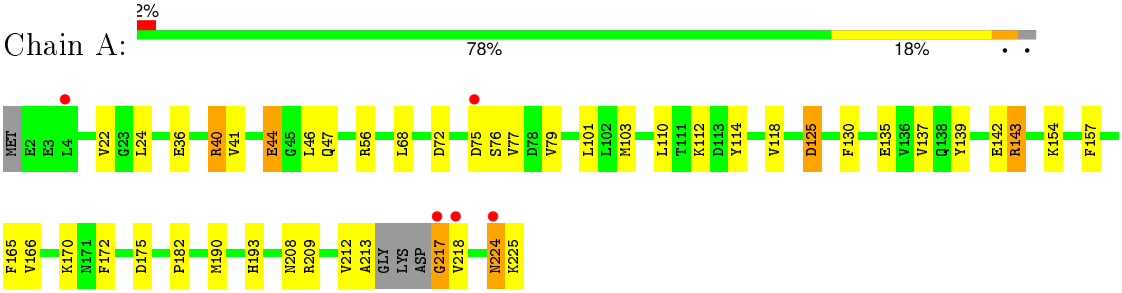
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	322	Total	O	0	0
			322	322		
2	B	289	Total	O	0	0
			289	289		

3 Residue-property plots (i)

These plots are drawn for all protein, RNA 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.

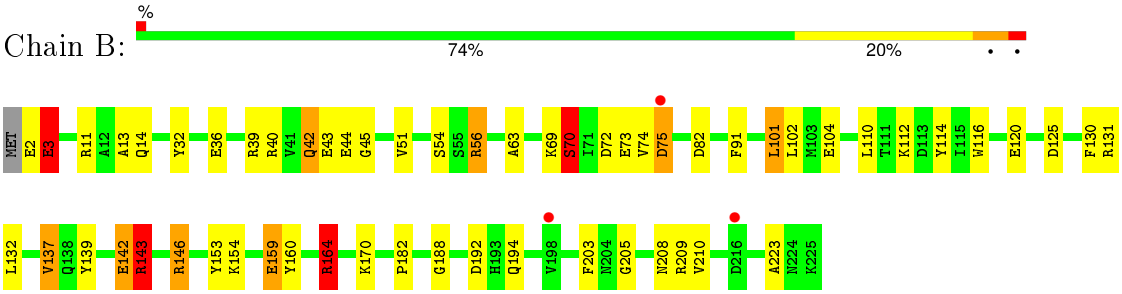
- Molecule 1: Ribose-5-phosphate isomerase A

Chain A:



- Molecule 1: Ribose-5-phosphate isomerase A

Chain B:



4 Data and refinement statistics (i)

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 31.16 – 1.70	Depositor EDS
% Data completeness (in resolution range)	96.1 (32.80-1.70) 97.8 (31.16-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.89 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.197 , 0.214 0.202 , 0.219	Depositor DCC
R_{free} test set	2589 reflections (5.10%)	DCC
Wilson B-factor (Å^2)	16.8	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 40.4	EDS
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.52$, $< L^2 > = 0.36$	Xtriage
Outliers	1 of 100910 reflections (0.001%)	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.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	1.73	20/1722 (1.2%)	1.29	6/2325 (0.3%)
1	B	1.84	41/1741 (2.4%)	1.41	19/2352 (0.8%)
All	All	1.79	61/3463 (1.8%)	1.35	25/4677 (0.5%)

All (61) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	153	TYR	CD1-CE1	10.13	1.54	1.39
1	B	91	PHE	CE1-CZ	8.27	1.53	1.37
1	B	120	GLU	CG-CD	7.71	1.63	1.51
1	B	73	GLU	CG-CD	7.67	1.63	1.51
1	B	3	GLU	CG-CD	7.31	1.62	1.51
1	A	130	PHE	CE1-CZ	7.21	1.51	1.37
1	B	143	ARG	CB-CG	7.12	1.71	1.52
1	B	194	GLN	CG-CD	7.09	1.67	1.51
1	B	2	GLU	N-CA	6.88	1.60	1.46
1	A	165	PHE	CD1-CE1	6.85	1.52	1.39
1	B	139	TYR	CD2-CE2	6.62	1.49	1.39
1	B	13	ALA	CA-CB	-6.59	1.38	1.52
1	A	143	ARG	CB-CG	-6.51	1.34	1.52
1	B	75	ASP	CB-CG	6.45	1.65	1.51
1	B	160	TYR	CE1-CZ	6.41	1.46	1.38
1	A	44	GLU	CD-OE1	6.29	1.32	1.25
1	B	11	ARG	CG-CD	6.25	1.67	1.51
1	A	137	VAL	CB-CG2	6.23	1.66	1.52
1	A	157	PHE	CG-CD2	6.14	1.48	1.38
1	B	54	SER	CB-OG	6.04	1.50	1.42
1	B	43	GLU	CB-CG	5.94	1.63	1.52
1	B	137	VAL	CA-CB	5.88	1.67	1.54
1	A	172	PHE	CD1-CE1	5.86	1.50	1.39
1	B	44	GLU	CD-OE2	-5.74	1.19	1.25
1	A	125	ASP	CB-CG	5.73	1.63	1.51
1	B	104	GLU	CB-CG	5.72	1.63	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	130	PHE	CE2-CZ	5.70	1.48	1.37
1	A	157	PHE	CE1-CZ	5.69	1.48	1.37
1	B	32	TYR	CE2-CZ	5.67	1.46	1.38
1	A	77	VAL	CB-CG1	5.67	1.64	1.52
1	B	42	GLN	CB-CG	5.66	1.67	1.52
1	A	79	VAL	CB-CG1	5.54	1.64	1.52
1	B	142	GLU	CD-OE1	5.52	1.31	1.25
1	B	188	GLY	N-CA	5.50	1.54	1.46
1	B	203	PHE	CE1-CZ	5.48	1.47	1.37
1	B	51	VAL	CB-CG2	5.44	1.64	1.52
1	B	154	LYS	CB-CG	5.41	1.67	1.52
1	B	63	ALA	CA-CB	5.34	1.63	1.52
1	B	56	ARG	CZ-NH2	5.34	1.40	1.33
1	B	160	TYR	CG-CD2	5.32	1.46	1.39
1	A	22	VAL	CB-CG1	5.32	1.64	1.52
1	A	217	GLY	C-O	5.30	1.32	1.23
1	B	74	VAL	CA-CB	5.27	1.65	1.54
1	B	130	PHE	CD1-CE1	5.26	1.49	1.39
1	B	159	GLU	CG-CD	-5.25	1.44	1.51
1	B	70	SER	CA-CB	5.25	1.60	1.52
1	B	45	GLY	C-O	5.24	1.32	1.23
1	A	118	VAL	CB-CG2	5.24	1.63	1.52
1	B	170	LYS	N-CA	5.21	1.56	1.46
1	A	44	GLU	CG-CD	5.20	1.59	1.51
1	A	139	TYR	CD2-CE2	5.19	1.47	1.39
1	B	194	GLN	CD-OE1	5.19	1.35	1.24
1	A	142	GLU	CB-CG	5.19	1.62	1.52
1	A	135	GLU	CB-CG	5.17	1.61	1.52
1	A	166	VAL	CB-CG1	5.17	1.63	1.52
1	B	223	ALA	C-O	5.12	1.33	1.23
1	B	3	GLU	CD-OE1	5.11	1.31	1.25
1	A	212	VAL	CB-CG2	5.09	1.63	1.52
1	B	182	PRO	CA-C	5.08	1.63	1.52
1	B	210	VAL	CB-CG1	5.05	1.63	1.52
1	B	143	ARG	CG-CD	5.04	1.64	1.51

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	131	ARG	NE-CZ-NH2	-12.77	113.92	120.30
1	B	143	ARG	NE-CZ-NH1	10.72	125.66	120.30
1	B	131	ARG	NE-CZ-NH1	10.16	125.38	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	72	ASP	CB-CG-OD2	8.84	126.26	118.30
1	B	11	ARG	NE-CZ-NH2	8.35	124.48	120.30
1	B	143	ARG	CG-CD-NE	7.04	126.58	111.80
1	B	72	ASP	CB-CG-OD1	-6.80	112.18	118.30
1	B	101	LEU	CB-CG-CD2	6.48	122.02	111.00
1	B	164	ARG	NE-CZ-NH1	6.30	123.45	120.30
1	B	146	ARG	NE-CZ-NH1	6.25	123.42	120.30
1	B	75	ASP	CB-CG-OD1	6.15	123.83	118.30
1	B	102	LEU	CB-CG-CD2	-6.03	100.76	111.00
1	A	40	ARG	NE-CZ-NH2	-5.94	117.33	120.30
1	A	24	LEU	CB-CG-CD1	-5.92	100.94	111.00
1	B	209	ARG	NE-CZ-NH1	-5.92	117.34	120.30
1	A	79	VAL	CG1-CB-CG2	-5.78	101.66	110.90
1	B	39	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	A	170	LYS	CD-CE-NZ	-5.60	98.82	111.70
1	A	175	ASP	CB-CG-OD1	5.38	123.14	118.30
1	B	143	ARG	CD-NE-CZ	5.31	131.04	123.60
1	B	125	ASP	CB-CG-OD2	-5.27	113.56	118.30
1	A	36	GLU	CG-CD-OE1	5.20	128.70	118.30
1	B	56	ARG	NE-CZ-NH1	-5.09	117.75	120.30
1	B	132	LEU	CB-CG-CD2	-5.07	102.37	111.00
1	B	164	ARG	CG-CD-NE	5.07	122.45	111.80

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
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	0.94
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	0.87
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:B:36:GLU:OE2	2:B:683:HOH:O	2.01	0.78
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:A:224:ASN:O	1:A:225:LYS:C	2.27	0.71
1:A:193:HIS:HD2	1:B:192:ASP:OD2	1.75	0.69
1:B:75:ASP:O	2:B:679:HOH:O	2.12	0.68
1:A:218:VAL:HG22	2:A:281:HOH:O	1.94	0.68
1:B:164:ARG:HD2	2:B:686:HOH:O	1.73	0.66
1:B:143:ARG:CD	1:B:146:ARG:HE	2.09	0.65
1:B:3:GLU:HG2	2:B:317:HOH:O	1.96	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:76:SER:CB	1:A:112:LYS:NZ	2.64	0.61
1:B:36:GLU:CG	2:B:683:HOH:O	2.49	0.59
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:72:ASP:CG	2:A:570:HOH:O	2.43	0.56
1:B:142:GLU:CD	2:B:662:HOH:O	2.33	0.56
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:41:VAL:N	2:A:664:HOH:O	2.34	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:103:MET:HE1	2:A:626:HOH:O	2.14	0.46
1:A:76:SER:CB	1:A:112:LYS:HZ2	2.29	0.46
1:B:82:ASP:O	1:B:116:TRP:HA	2.16	0.45
1:A:209:ARG:NH2	2:A:677:HOH:O	2.50	0.45
1:A:101:LEU:H	1:A:101:LEU:HD23	1.81	0.45
1:A:110:LEU:HG	1:A:110:LEU:O	2.17	0.44
1:B:143:ARG:HD2	1:B:143:ARG:HA	1.68	0.43
1:A:40:ARG:C	2:A:664:HOH:O	2.57	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:HB3	1:A:225:LYS:HE3	1.82	0.42
1:A:46:LEU:CA	2:A:664:HOH:O	2.59	0.41
1:B:112:LYS:HE3	2:B:614:HOH:O	2.20	0.41
1:A:209:ARG:NH1	2:A:677:HOH:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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 (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	217/225 (96%)	214 (99%)	2 (1%)	1 (0%)	34 15
1	B	222/225 (99%)	219 (99%)	3 (1%)	0	100 100
All	All	439/450 (98%)	433 (99%)	5 (1%)	1 (0%)	52 32

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	224	ASN

5.3.2 Protein sidechains (i)

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	180/184 (98%)	179 (99%)	1 (1%)	90 85
1	B	182/184 (99%)	176 (97%)	6 (3%)	45 22
All	All	362/368 (98%)	355 (98%)	7 (2%)	65 46

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ARG
1	B	3	GLU
1	B	70	SER
1	B	101	LEU
1	B	143	ARG
1	B	159	GLU
1	B	164	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	ASN
1	A	138	GLN
1	A	193	HIS
1	A	208	ASN
1	B	90	ASN
1	B	194	GLN
1	B	204	ASN
1	B	208	ASN
1	B	224	ASN

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

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, 95^{th} 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%)	-0.04	5 (2%)	64 68	9, 14, 27, 41	0
1	B	224/225 (99%)	-0.02	3 (1%)	79 83	10, 15, 31, 46	0
All	All	445/450 (98%)	-0.03	8 (1%)	71 76	9, 14, 30, 46	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	217	GLY	4.9
1	B	75	ASP	4.3
1	A	218	VAL	3.1
1	A	75	ASP	2.4
1	A	224	ASN	2.2
1	B	216	ASP	2.2
1	A	4	LEU	2.1
1	B	198	VAL	2.1

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 (i)

There are no such residues in this entry.