Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 01:52 am BST

PDB ID	:	3PR7
Title	:	Multi-functional and mechanosensitive receptor binding activity of the Moraxella catarrhalis adhesin UspA1
Authors	:	Agnew, C.R.J.; Zaccai, N.R.; Conners, R.; Brady, R.L.
Deposited on	:	2010-11-29
Resolution	:	2.94 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 following versions of software and data (see references ①) were used in the production of this report:

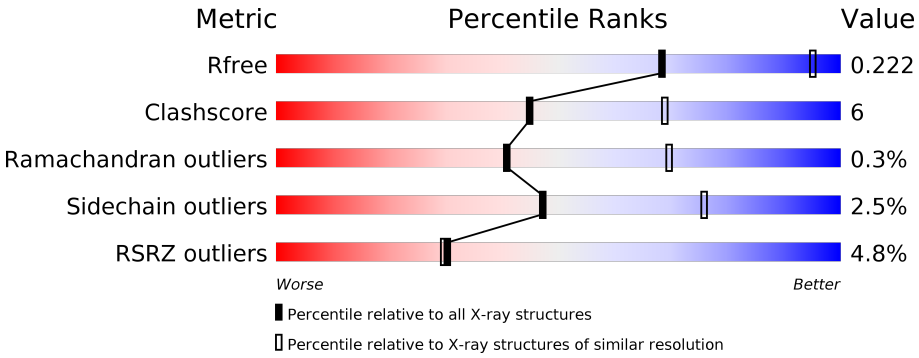
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.94 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2969 (2.98-2.90)
Clashscore	141614	3218 (2.98-2.90)
Ramachandran outliers	138981	3122 (2.98-2.90)
Sidechain outliers	138945	3124 (2.98-2.90)
RSRZ outliers	127900	2902 (2.98-2.90)

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 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 <=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	311	5% 81% 12% • 6%
1	B	311	6% 81% 12% • 6%
1	C	311	3% 80% 13% • 6%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 6177 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 UspA1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	292	Total 2044	C 1225	N 377	O 442	0	1	0
1	B	292	Total 2044	C 1225	N 377	O 442	0	1	0
1	C	292	Total 2044	C 1225	N 377	O 442	0	1	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	346	LYS	-	EXPRESSION TAG	UNP Q9XD56
A	347	HIS	-	EXPRESSION TAG	UNP Q9XD56
A	348	HIS	-	EXPRESSION TAG	UNP Q9XD56
A	349	HIS	-	EXPRESSION TAG	UNP Q9XD56
A	350	HIS	-	EXPRESSION TAG	UNP Q9XD56
A	351	HIS	-	EXPRESSION TAG	UNP Q9XD56
A	352	HIS	-	EXPRESSION TAG	UNP Q9XD56
B	346	LYS	-	EXPRESSION TAG	UNP Q9XD56
B	347	HIS	-	EXPRESSION TAG	UNP Q9XD56
B	348	HIS	-	EXPRESSION TAG	UNP Q9XD56
B	349	HIS	-	EXPRESSION TAG	UNP Q9XD56
B	350	HIS	-	EXPRESSION TAG	UNP Q9XD56
B	351	HIS	-	EXPRESSION TAG	UNP Q9XD56
B	352	HIS	-	EXPRESSION TAG	UNP Q9XD56
C	346	LYS	-	EXPRESSION TAG	UNP Q9XD56
C	347	HIS	-	EXPRESSION TAG	UNP Q9XD56
C	348	HIS	-	EXPRESSION TAG	UNP Q9XD56
C	349	HIS	-	EXPRESSION TAG	UNP Q9XD56
C	350	HIS	-	EXPRESSION TAG	UNP Q9XD56
C	351	HIS	-	EXPRESSION TAG	UNP Q9XD56
C	352	HIS	-	EXPRESSION TAG	UNP Q9XD56

- Molecule 2 is water.

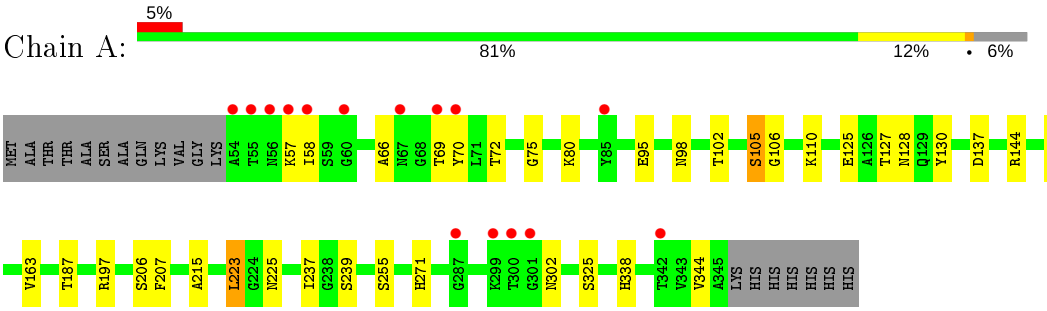
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	14	Total O 14 14	0	0
2	B	7	Total O 7 7	0	0
2	C	24	Total O 24 24	0	0

3 Residue-property plots (i)

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

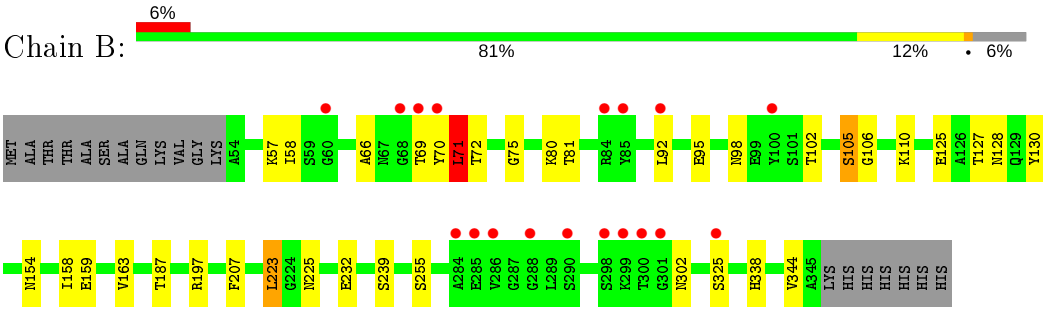
- Molecule 1: UspA1

Chain A:



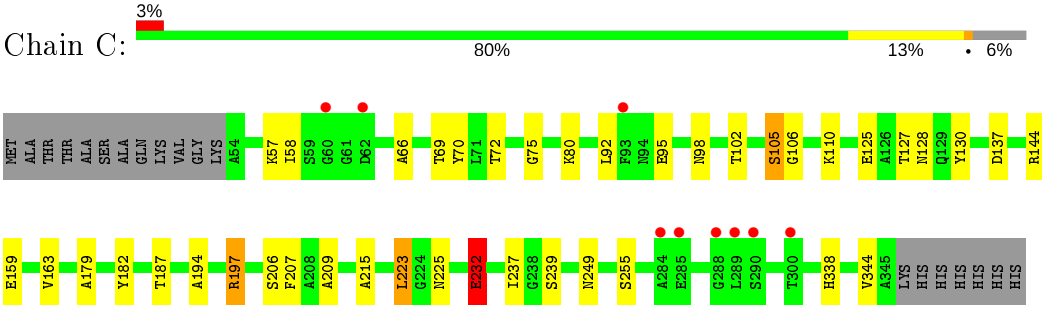
- Molecule 1: UspA1

Chain B:



- Molecule 1: UspA1

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α, β, γ	118.22Å 118.22Å 176.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.60 – 2.94 49.12 – 2.94	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.60-2.94) 99.9 (49.12-2.94)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.55 (at 2.96Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.203 , 0.216 0.209 , 0.222	Depositor DCC
R_{free} test set	1503 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	54.3	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 59.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.043 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6177	wwPDB-VP
Average B, all atoms (Å^2)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.53% 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, 0.333 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	0.33	0/2065	0.60	4/2786 (0.1%)
1	B	0.34	1/2065 (0.0%)	0.61	5/2786 (0.2%)
1	C	0.83	2/2065 (0.1%)	0.77	8/2786 (0.3%)
All	All	0.55	3/6195 (0.0%)	0.67	17/8358 (0.2%)

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	C	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	232	GLU	CD-OE2	34.31	1.63	1.25
1	B	232	GLU	CD-OE2	6.98	1.33	1.25
1	C	232	GLU	CB-CG	-6.58	1.39	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	197	ARG	NE-CZ-NH1	-15.61	112.50	120.30
1	C	197	ARG	NE-CZ-NH2	14.55	127.58	120.30
1	C	144	ARG	NE-CZ-NH2	-11.90	114.35	120.30
1	C	144	ARG	NE-CZ-NH1	10.48	125.54	120.30
1	A	197	ARG	NE-CZ-NH1	7.89	124.25	120.30
1	C	197	ARG	CD-NE-CZ	7.67	134.34	123.60
1	B	144	ARG	NE-CZ-NH1	-7.66	116.47	120.30
1	B	197	ARG	NE-CZ-NH1	7.45	124.03	120.30
1	A	197	ARG	NE-CZ-NH2	-7.40	116.60	120.30
1	B	197	ARG	NE-CZ-NH2	-7.30	116.65	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	144	ARG	NE-CZ-NH1	-7.29	116.66	120.30
1	B	144	ARG	NE-CZ-NH2	7.25	123.92	120.30
1	A	144	ARG	NE-CZ-NH2	7.01	123.81	120.30
1	C	144	ARG	CD-NE-CZ	5.90	131.87	123.60
1	B	71	LEU	CB-CG-CD1	5.59	120.50	111.00
1	C	197	ARG	CG-CD-NE	-5.53	100.18	111.80
1	C	232	GLU	OE1-CD-OE2	-5.47	116.73	123.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	232	GLU	Sidechain

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	2044	0	1963	32	0
1	B	2044	0	1963	29	0
1	C	2044	0	1963	33	0
2	A	14	0	0	0	0
2	B	7	0	0	0	0
2	C	24	0	0	0	0
All	All	6177	0	5889	78	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 6.

All (78) 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:A:72:THR:HG21	1:B:75:GLY:HA3	1.45	0.96
1:B:72:THR:HG21	1:C:75:GLY:HA3	1.47	0.93
1:A:75:GLY:HA3	1:C:72:THR:HG21	1.49	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:ILE:HG22	1:A:159:GLU:N	2.10	0.67
1:B:158:ILE:HG22	1:B:159:GLU:N	2.09	0.66
1:C:158:ILE:HG22	1:C:159:GLU:N	2.10	0.66
1:B:72:THR:CG2	1:C:75:GLY:HA3	2.27	0.58
1:C:69:THR:HG22	1:C:70:TYR:CD2	2.41	0.56
1:A:72:THR:CG2	1:B:75:GLY:HA3	2.28	0.55
1:A:75:GLY:HA3	1:C:72:THR:CG2	2.30	0.55
1:B:158:ILE:CG2	1:B:159:GLU:N	2.70	0.55
1:A:69:THR:HG22	1:A:70:TYR:CD2	2.43	0.54
1:B:69:THR:HG22	1:B:70:TYR:CD2	2.42	0.54
1:B:158:ILE:CG2	1:B:159:GLU:H	2.21	0.53
1:B:71:LEU:HG	1:B:81:THR:HB	1.91	0.53
1:A:207:PHE:HE2	1:A:223[B]:LEU:HD21	1.74	0.53
1:A:158:ILE:CG2	1:A:159:GLU:N	2.70	0.53
1:B:158:ILE:HG22	1:B:159:GLU:H	1.72	0.53
1:C:158:ILE:CG2	1:C:159:GLU:N	2.72	0.53
1:A:158:ILE:CG2	1:A:159:GLU:H	2.22	0.52
1:C:158:ILE:HG22	1:C:159:GLU:H	1.74	0.52
1:C:239:SER:O	1:C:255:SER:HA	2.10	0.52
1:A:239:SER:O	1:A:255:SER:HA	2.09	0.52
1:B:207:PHE:HE2	1:B:223[B]:LEU:HD21	1.74	0.52
1:A:158:ILE:HG22	1:A:159:GLU:H	1.73	0.52
1:C:110:LYS:HB2	1:C:125:GLU:HG2	1.92	0.51
1:C:207:PHE:HE2	1:C:223[B]:LEU:HD21	1.74	0.51
1:A:154:ASN:HB3	1:A:163:VAL:HG11	1.92	0.51
1:B:225:ASN:O	1:B:239:SER:HA	2.11	0.51
1:B:239:SER:O	1:B:255:SER:HA	2.11	0.51
1:B:110:LYS:HB2	1:B:125:GLU:HG2	1.93	0.50
1:A:110:LYS:HB2	1:A:125:GLU:HG2	1.93	0.50
1:C:225:ASN:O	1:C:239:SER:HA	2.10	0.50
1:C:57:LYS:HB2	1:C:72:THR:HG22	1.93	0.50
1:C:182:TYR:HB3	1:C:197:ARG:HG3	1.92	0.50
1:A:225:ASN:O	1:A:239:SER:HA	2.11	0.50
1:C:158:ILE:CG2	1:C:159:GLU:H	2.25	0.50
1:A:57:LYS:HB2	1:A:72:THR:HG22	1.93	0.49
1:B:207:PHE:CE2	1:B:223[B]:LEU:HD21	2.47	0.49
1:B:57:LYS:HB2	1:B:72:THR:HG22	1.94	0.49
1:C:154:ASN:HB3	1:C:163:VAL:HG11	1.94	0.49
1:A:207:PHE:CE2	1:A:223[B]:LEU:HD21	2.47	0.49
1:A:102:THR:HG23	1:B:105:SER:HB3	1.93	0.48
1:B:344:VAL:O	1:B:344:VAL:HG12	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:207:PHE:CE2	1:C:223[B]:LEU:HD21	2.48	0.48
1:A:58:ILE:HG21	1:A:66:ALA:HB2	1.96	0.48
1:B:58:ILE:HG21	1:B:66:ALA:HB2	1.96	0.48
1:A:137:ASP:HB2	1:C:130:TYR:CD1	2.49	0.47
1:A:344:VAL:O	1:A:344:VAL:HG12	2.14	0.47
1:B:105:SER:OG	1:B:106:GLY:N	2.45	0.47
1:C:344:VAL:HG12	1:C:344:VAL:O	2.13	0.47
1:C:338:HIS:CD2	1:C:338:HIS:C	2.88	0.47
1:A:105:SER:HB3	1:C:102:THR:HG23	1.97	0.47
1:A:338:HIS:ND1	1:B:325:SER:HA	2.29	0.47
1:A:325:SER:HA	1:C:338:HIS:ND1	2.30	0.46
1:C:58:ILE:HG21	1:C:66:ALA:HB2	1.96	0.46
1:B:154:ASN:HB3	1:B:163:VAL:HG11	1.97	0.46
1:B:338:HIS:C	1:B:338:HIS:CD2	2.89	0.46
1:A:338:HIS:CD2	1:A:338:HIS:C	2.89	0.45
1:B:102:THR:HG23	1:C:105:SER:HB3	1.99	0.45
1:B:80:LYS:HB2	1:B:95:GLU:HG2	1.99	0.44
1:C:80:LYS:HB2	1:C:95:GLU:HG2	1.99	0.44
1:A:80:LYS:HB2	1:A:95:GLU:HG2	1.99	0.44
1:C:105:SER:OG	1:C:106:GLY:N	2.47	0.43
1:B:130:TYR:CD1	1:C:137:ASP:HB2	2.53	0.43
1:B:127:THR:O	1:B:128:ASN:CB	2.67	0.43
1:C:127:THR:O	1:C:128:ASN:CB	2.66	0.43
1:A:127:THR:O	1:A:128:ASN:CB	2.65	0.42
1:A:105:SER:OG	1:A:106:GLY:N	2.42	0.41
1:A:130:TYR:CD1	1:B:137:ASP:HB2	2.54	0.41
1:A:102:THR:CG2	1:B:105:SER:HB3	2.51	0.41
1:A:206:SER:HB2	1:A:215:ALA:HB1	2.03	0.41
1:C:194:ALA:HB3	1:C:209:ALA:HB2	2.03	0.40
1:C:223[B]:LEU:HB2	1:C:237:ILE:HG12	2.02	0.40
1:A:223[B]:LEU:HB2	1:A:237:ILE:HG12	2.03	0.40
1:C:179:ALA:HB3	1:C:194:ALA:O	2.22	0.40
1:C:206:SER:HB2	1:C:215:ALA:HB1	2.03	0.40
1:A:271:HIS:HB2	1:C:249:ASN:HB3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	291/311 (94%)	278 (96%)	12 (4%)	1 (0%)	41	69
1	B	291/311 (94%)	276 (95%)	14 (5%)	1 (0%)	41	69
1	C	291/311 (94%)	278 (96%)	12 (4%)	1 (0%)	41	69
All	All	873/933 (94%)	832 (95%)	38 (4%)	3 (0%)	41	69

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	105	SER
1	B	105	SER
1	C	105	SER

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	204/218 (94%)	199 (98%)	5 (2%)	47	76
1	B	204/218 (94%)	197 (97%)	7 (3%)	37	68
1	C	204/218 (94%)	198 (97%)	6 (3%)	42	73
All	All	612/654 (94%)	594 (97%)	18 (3%)	47	73

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	A	187	THR
1	A	223[A]	LEU
1	A	223[B]	LEU
1	A	302	ASN
1	B	71	LEU
1	B	92	LEU
1	B	98	ASN
1	B	187	THR
1	B	223[A]	LEU
1	B	223[B]	LEU
1	B	302	ASN
1	C	92	LEU
1	C	98	ASN
1	C	187	THR
1	C	223[A]	LEU
1	C	223[B]	LEU
1	C	232	GLU

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

Mol	Chain	Res	Type
1	C	302	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 (i)

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	292/311 (93%)	0.12	15 (5%)	28	27	34, 58, 112, 159	0
1	B	292/311 (93%)	0.26	18 (6%)	20	18	39, 63, 114, 144	0
1	C	292/311 (93%)	0.09	9 (3%)	49	48	29, 52, 121, 151	0
All	All	876/933 (93%)	0.16	42 (4%)	30	30	29, 58, 115, 159	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	300	THR	4.8
1	C	60	GLY	4.7
1	A	287	GLY	4.4
1	B	299	LYS	4.2
1	C	300	THR	3.8
1	A	57	LYS	3.6
1	A	299	LYS	3.6
1	A	300	THR	3.5
1	A	55	THR	3.5
1	B	288	GLY	3.5
1	B	60	GLY	3.4
1	C	288	GLY	3.4
1	B	84	ARG	3.3
1	C	285	GLU	3.2
1	A	69	THR	3.1
1	A	67	ASN	3.0
1	A	85	TYR	2.9
1	B	286	VAL	2.7
1	A	58	ILE	2.6
1	B	85	TYR	2.5
1	A	301	GLY	2.5
1	B	298	SER	2.5
1	A	60	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	70	TYR	2.4
1	B	285	GLU	2.4
1	C	290	SER	2.4
1	B	70	TYR	2.4
1	B	69	THR	2.3
1	B	68	GLY	2.3
1	B	301	GLY	2.3
1	B	100	TYR	2.2
1	A	54	ALA	2.1
1	C	284	ALA	2.1
1	A	56	ASN	2.1
1	C	289	LEU	2.1
1	C	62	ASP	2.1
1	B	290	SER	2.1
1	C	93	PHE	2.1
1	B	284	ALA	2.1
1	B	325	SER	2.0
1	B	92	LEU	2.0
1	A	342	THR	2.0

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

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

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.