Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 12:43 PM GMT

PDB ID	:	3RRS
Title	:	Crystal structure analysis of cellobiose phosphorylase from Cellulomonas uda
Authors	:	Van Hoorebeke, A.; Stout, J.; Soetaert, W.; Van Beeumen, J.; Desmet, T.; Savvides, S.
Deposited on	:	2011-04-30
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:

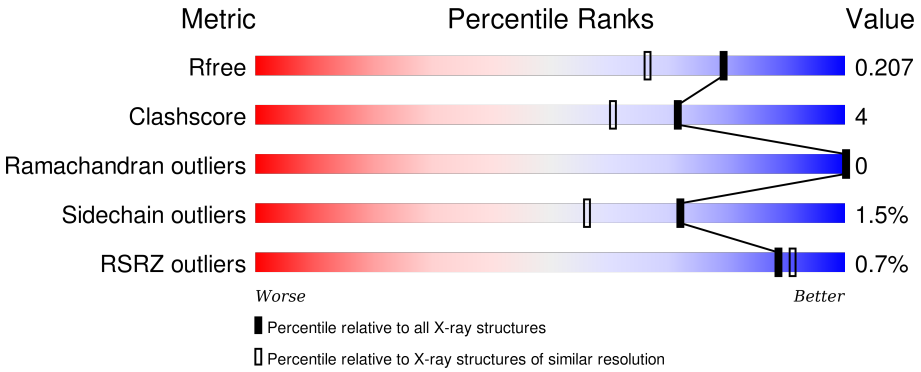
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

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	822	% 91% 9%
1	B	822	% 91% 9%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 27811 atoms, of which 12439 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 Cellobiose phosphorylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	822	12735	4144	6184	1123	1267	17	0	19	0
			Total	C	H	N	O	S			
1	B	822	12850	4186	6255	1125	1267	17	0	27	0

- Molecule 2 is water.

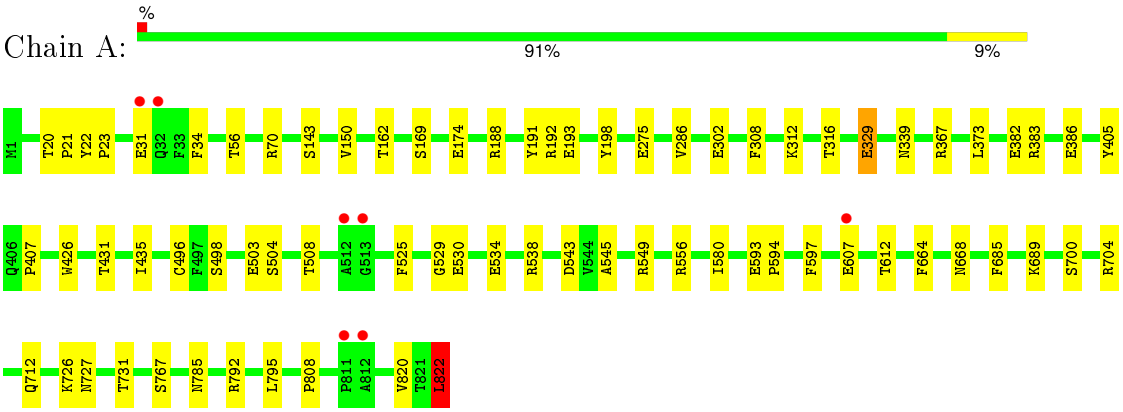
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	1100	1100	1100	0	0
			Total	O		
2	B	1126	1126	1126	0	0

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 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.

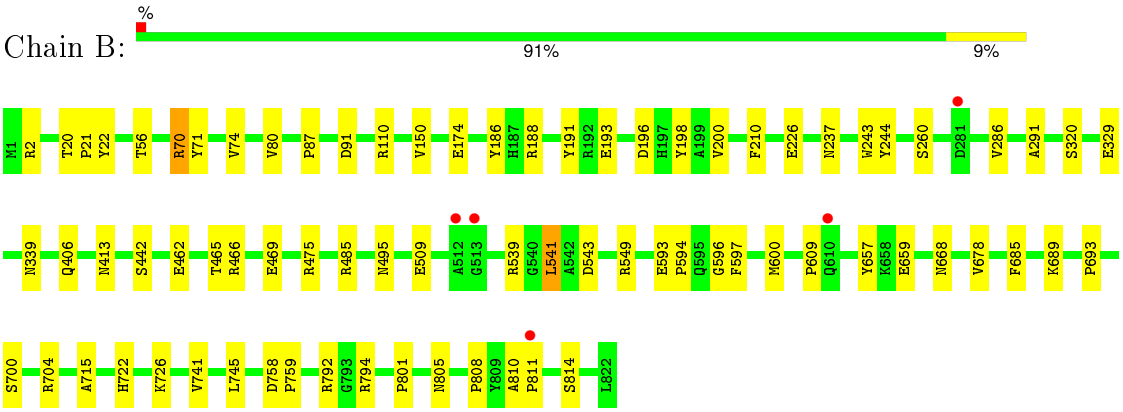
- Molecule 1: Cellobiose phosphorylase

Chain A:



- Molecule 1: Cellobiose phosphorylase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	85.69Å 103.80Å 98.50Å 90.00° 96.51° 90.00°	Depositor
Resolution (Å)	19.75 – 1.70 19.75 – 1.70	Depositor EDS
% Data completeness (in resolution range)	96.6 (19.75-1.70) 96.6 (19.75-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.38 (at 1.70Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.185 , 0.209 0.182 , 0.207	Depositor DCC
R_{free} test set	1812 reflections (1.00%)	DCC
Wilson B-factor (Å^2)	13.0	Xtriage
Anisotropy	0.500	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.44 , 63.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 181275 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	27811	wwPDB-VP
Average B, all atoms (Å^2)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.94% 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	0.79	4/6776 (0.1%)	0.83	5/9231 (0.1%)
1	B	0.81	3/6835 (0.0%)	0.84	11/9312 (0.1%)
All	All	0.80	7/13611 (0.1%)	0.84	16/18543 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	664	PHE	CB-CG	-7.63	1.38	1.51
1	B	22	TYR	CD1-CE1	-6.45	1.29	1.39
1	A	192	ARG	CB-CG	-5.97	1.36	1.52
1	B	659	GLU	CD-OE1	-5.63	1.19	1.25
1	A	556	ARG	CB-CG	-5.62	1.37	1.52
1	A	191	TYR	CE1-CZ	-5.47	1.31	1.38
1	B	244	TYR	CD2-CE2	-5.07	1.31	1.39

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	792	ARG	NE-CZ-NH1	10.01	125.30	120.30
1	A	792	ARG	NE-CZ-NH2	-9.70	115.45	120.30
1	B	543	ASP	CB-CG-OD1	9.43	126.78	118.30
1	A	543	ASP	CB-CG-OD1	9.25	126.62	118.30
1	B	792	ARG	NE-CZ-NH2	-7.18	116.71	120.30
1	B	485	ARG	NE-CZ-NH2	-6.96	116.82	120.30
1	B	543	ASP	CB-CG-OD2	-6.58	112.38	118.30
1	B	792	ARG	NE-CZ-NH1	6.18	123.39	120.30
1	A	822	LEU	CA-CB-CG	5.66	128.32	115.30
1	B	196	ASP	CB-CG-OD1	5.66	123.39	118.30
1	B	549	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	B	475	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	B	2	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	B	70	ARG	NE-CZ-NH1	5.21	122.91	120.30
1	A	556	ARG	NE-CZ-NH1	5.19	122.89	120.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	2	ARG	NE-CZ-NH2	-5.01	117.79	120.30

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	6551	6184	6202	56	0
1	B	6595	6255	6270	56	0
2	A	1100	0	0	8	0
2	B	1126	0	0	14	0
All	All	15372	12439	12472	109	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 4.

All (109) 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:31:GLU:O	1:A:34:PHE:CZ	2.06	1.09
1:B:539:ARG:HB3	1:B:541:LEU:HD13	1.44	0.96
1:A:383[B]:ARG:H	1:A:383[B]:ARG:HE	1.09	0.94
1:A:504:SER:O	1:A:508[B]:THR:HG22	1.73	0.88
1:A:593:GLU:HB2	1:A:594[A]:PRO:HD3	1.55	0.87
1:A:316:THR:HG21	2:A:2156:HOH:O	1.76	0.85
1:A:503:GLU:HB3	1:A:508[A]:THR:HG21	1.59	0.84
1:A:31:GLU:O	1:A:34:PHE:HZ	1.66	0.78
1:A:275[A]:GLU:HA	1:A:275[A]:GLU:OE1	1.82	0.78
1:A:382:GLU:H	1:A:383[B]:ARG:HH21	1.27	0.78
1:B:186:TYR:CD2	1:B:200[B]:VAL:HG12	2.24	0.72
1:B:593:GLU:HB2	1:B:594[A]:PRO:HD3	1.72	0.71
1:A:383[B]:ARG:H	1:A:383[B]:ARG:NE	1.87	0.71
1:A:545:ALA:O	1:A:549[B]:ARG:HG3	1.91	0.69
1:B:810:ALA:HB1	1:B:811[B]:PRO:HD2	1.72	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:580:ILE:HG12	2:A:1956:HOH:O	1.92	0.69
1:A:704[B]:ARG:HD3	2:A:900:HOH:O	1.92	0.68
1:A:20:THR:HB	1:A:21[B]:PRO:HD2	1.77	0.66
1:B:541:LEU:N	1:B:541:LEU:HD12	2.12	0.65
1:A:593:GLU:N	1:A:594[B]:PRO:HD2	2.11	0.65
1:A:329:GLU:HG2	2:A:2309:HOH:O	1.96	0.65
1:A:496:CYS:SG	1:A:508[B]:THR:HG23	2.37	0.65
1:A:767:SER:HB2	2:A:2115:HOH:O	1.98	0.62
1:A:504:SER:O	1:A:508[B]:THR:CG2	2.44	0.62
1:B:609[B]:PRO:HG3	1:B:678:VAL:O	2.00	0.62
1:B:593:GLU:N	1:B:594[B]:PRO:HD2	2.15	0.62
1:B:722:HIS:HE1	2:B:1014:HOH:O	1.84	0.60
1:A:31:GLU:O	1:A:34:PHE:CE2	2.54	0.59
1:B:200[A]:VAL:HG11	2:B:1975:HOH:O	2.03	0.59
1:A:534:GLU:O	1:A:538:ARG:HG3	2.03	0.59
1:B:811[B]:PRO:HG2	1:B:814:SER:OG	2.02	0.58
1:B:758:ASP:N	1:B:759[A]:PRO:HD3	2.19	0.58
1:B:693:PRO:HA	2:B:1082:HOH:O	2.02	0.57
1:B:200[B]:VAL:HG21	2:B:1975:HOH:O	2.05	0.57
1:A:504:SER:O	1:A:508[A]:THR:HG23	2.03	0.57
1:B:539:ARG:HB3	1:B:541:LEU:CD1	2.29	0.57
1:A:808[B]:PRO:HG2	2:A:1036:HOH:O	2.05	0.56
1:A:726[B]:LYS:NZ	1:B:243:TRP:NE1	2.54	0.56
1:B:320:SER:HB3	2:B:1388:HOH:O	2.04	0.56
1:A:56:THR:HG21	1:A:150[B]:VAL:HG11	1.88	0.56
1:A:496:CYS:SG	1:A:508[B]:THR:CG2	2.95	0.55
1:B:715:ALA:O	1:B:722:HIS:HD2	1.89	0.55
1:B:801:PRO:CG	2:B:2179:HOH:O	2.56	0.53
1:B:596:GLY:O	1:B:600:MET:HG3	2.09	0.53
1:A:198:TYR:CE2	1:A:286[B]:VAL:HG11	2.45	0.52
1:A:726[B]:LYS:NZ	1:B:243:TRP:CE2	2.77	0.52
1:B:329:GLU:CD	1:B:329:GLU:H	2.13	0.52
1:B:21[A]:PRO:HB3	1:B:704[A]:ARG:CZ	2.40	0.51
1:B:541:LEU:CD1	1:B:541:LEU:N	2.73	0.51
1:B:56:THR:HG21	1:B:150[A]:VAL:CG2	2.41	0.51
1:A:382:GLU:O	1:A:386[A]:GLU:HG3	2.10	0.51
1:A:56:THR:HG21	1:A:150[B]:VAL:CG1	2.41	0.51
1:A:22:TYR:CG	1:A:23:PRO:HD2	2.47	0.49
1:A:367:ARG:HB2	1:A:426:TRP:CD1	2.47	0.49
1:B:56:THR:HG21	1:B:150[A]:VAL:HG21	1.95	0.48
1:B:74:VAL:O	1:B:80:VAL:HA	2.13	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:593:GLU:N	1:A:594[B]:PRO:CD	2.74	0.48
1:B:200[B]:VAL:HG11	1:B:291:ALA:HB1	1.96	0.48
1:B:291:ALA:HA	2:B:1975:HOH:O	2.13	0.48
1:B:801:PRO:HG3	2:B:2179:HOH:O	2.13	0.48
1:B:191:TYR:CE2	1:B:286[B]:VAL:HG22	2.49	0.47
1:A:785:ASN:HA	1:A:822:LEU:HD13	1.96	0.47
1:B:21[B]:PRO:O	1:B:704[B]:ARG:HD2	2.16	0.46
1:A:275[A]:GLU:OE1	1:A:275[A]:GLU:CA	2.59	0.46
1:A:162[A]:THR:HG22	2:B:878:HOH:O	2.14	0.46
1:A:308:PHE:CZ	1:A:312:LYS:HE3	2.50	0.46
1:A:685:PHE:O	1:A:689:LYS:HG3	2.16	0.45
1:A:169:SER:HB3	1:B:657:TYR:HB3	1.98	0.45
1:B:685:PHE:O	1:B:689:LYS:HG3	2.17	0.45
1:B:191:TYR:CD2	1:B:286[B]:VAL:HG22	2.51	0.45
1:B:593:GLU:N	1:B:594[B]:PRO:CD	2.80	0.44
1:B:700:SER:HA	2:B:1024:HOH:O	2.17	0.44
1:A:198:TYR:CE2	1:A:286[B]:VAL:CG1	3.01	0.44
1:B:91:ASP:O	1:B:110:ARG:HD2	2.18	0.44
1:A:795:LEU:CD2	1:A:820:VAL:HG22	2.47	0.44
1:B:495:ASN:O	1:B:509:GLU:HA	2.17	0.44
1:B:805:ASN:HB3	2:B:2246:HOH:O	2.17	0.44
1:A:174:GLU:HG3	1:A:188:ARG:HD3	1.99	0.43
1:B:191:TYR:CD2	1:B:286[A]:VAL:HG12	2.53	0.43
1:B:726[B]:LYS:HE2	2:B:886:HOH:O	2.18	0.43
1:B:465:THR:O	1:B:469:GLU:HG2	2.18	0.43
1:A:431:THR:O	1:A:435:ILE:HG13	2.18	0.43
1:A:700:SER:HA	2:A:1485:HOH:O	2.19	0.43
1:A:22:TYR:CD2	1:A:23:PRO:HD2	2.54	0.42
1:A:607:GLU:HA	2:A:1921:HOH:O	2.19	0.42
1:B:539:ARG:HG2	1:B:541:LEU:HD11	2.01	0.42
1:B:210:PHE:O	1:B:237:ASN:HA	2.19	0.42
1:A:302[B]:GLU:H	1:A:302[B]:GLU:CD	2.23	0.42
1:B:198:TYR:CE2	1:B:286[B]:VAL:HG11	2.55	0.42
1:B:704[B]:ARG:NH2	2:B:1479:HOH:O	2.53	0.41
1:A:712:GLN:HB3	1:A:727:ASN:HB2	2.02	0.41
1:A:496:CYS:SG	1:A:498[B]:SER:OG	2.68	0.41
1:B:174:GLU:HG3	1:B:188:ARG:HD3	2.02	0.41
1:B:20:THR:HB	1:B:21[B]:PRO:HD2	2.02	0.41
1:B:186:TYR:CE2	1:B:200[B]:VAL:HG12	2.55	0.41
1:A:373:LEU:O	1:A:373:LEU:HD23	2.19	0.41
1:B:741:VAL:HA	1:B:745:LEU:HD12	2.01	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:593:GLU:HB2	1:A:594[A]:PRO:CD	2.39	0.41
1:A:405:TYR:O	1:A:407:PRO:HD3	2.21	0.41
1:A:530:GLU:CD	1:A:549[B]:ARG:HH11	2.24	0.41
1:B:811[B]:PRO:HG2	1:B:814:SER:HG	1.85	0.40
1:B:71:TYR:CZ	1:B:87:PRO:HD3	2.56	0.40
1:B:808:PRO:HG3	2:B:1979:HOH:O	2.21	0.40
1:A:525:PHE:O	1:A:529:GLY:HA3	2.21	0.40
1:B:406:GLN:OE1	1:B:413:ASN:HB2	2.22	0.40
1:B:462:GLU:O	1:B:466:ARG:HG3	2.22	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	839/822 (102%)	801 (96%)	38 (4%)	0	100	100
1	B	847/822 (103%)	811 (96%)	36 (4%)	0	100	100
All	All	1686/1644 (103%)	1612 (96%)	74 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	688/675 (102%)	678 (98%)	10 (2%)	72	56
1	B	693/675 (103%)	683 (99%)	10 (1%)	74	59
All	All	1381/1350 (102%)	1361 (99%)	20 (1%)	72	59

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	ARG
1	A	143	SER
1	A	193	GLU
1	A	329	GLU
1	A	339	ASN
1	A	597	PHE
1	A	612	THR
1	A	668	ASN
1	A	731	THR
1	A	822	LEU
1	B	70	ARG
1	B	193	GLU
1	B	226	GLU
1	B	260	SER
1	B	339	ASN
1	B	442	SER
1	B	541	LEU
1	B	597	PHE
1	B	668	ASN
1	B	794	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	111	ASN
1	A	126	ASN
1	A	161	GLN
1	A	414	ASN
1	A	511	GLN
1	A	610	GLN
1	B	126	ASN
1	B	130	GLN
1	B	722	HIS

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 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	822/822 (100%)	-0.20	7 (0%) 85 88	8, 13, 24, 38	0
1	B	822/822 (100%)	-0.16	5 (0%) 90 92	7, 13, 22, 39	0
All	All	1644/1644 (100%)	-0.18	12 (0%) 89 91	7, 13, 23, 39	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	512	ALA	4.7
1	A	513	GLY	3.8
1	A	512	ALA	2.8
1	B	811[A]	PRO	2.8
1	B	513	GLY	2.6
1	A	811	PRO	2.5
1	A	812	ALA	2.5
1	B	610	GLN	2.5
1	A	31	GLU	2.4
1	A	32	GLN	2.3
1	A	607	GLU	2.3
1	B	281	ASP	2.2

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.