W O R L D W I D E
wwPDB
P R O T E I N   D A T A   B A N K

Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 07:23 AM GMT

PDB ID	:	3AMC
Title	:	Crystal structures of Thermotoga maritima Cel5A, apo form and dimer/au
Authors	:	Wu, T.H.; Huang, C.H.; Ko, T.P.; Lai, H.L.; Ma, Y.; Cheng, Y.S.; Liu, J.R.; Guo, R.T.
Deposited on	:	2010-08-19
Resolution	:	1.40 Å(reported)

The following versions of software and data (see references (i)) 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 1.40 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.
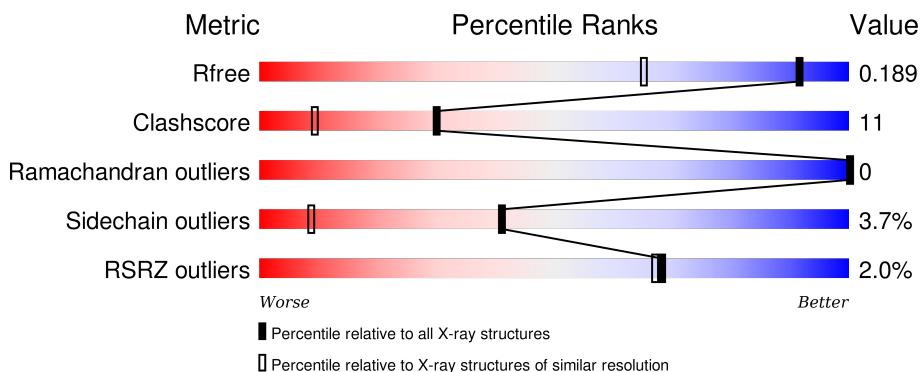
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1199 (1.40-1.40)
Clashscore	102246	1295 (1.40-1.40)
Ramachandran outliers	100387	1259 (1.40-1.40)
Sidechain outliers	100360	1258 (1.40-1.40)
RSRZ outliers	91569	1198 (1.40-1.40)

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	317	2% red bar; 78% green, 16% yellow, · ·
1	B	317	2% red bar; 79% green, 14% yellow, · ·
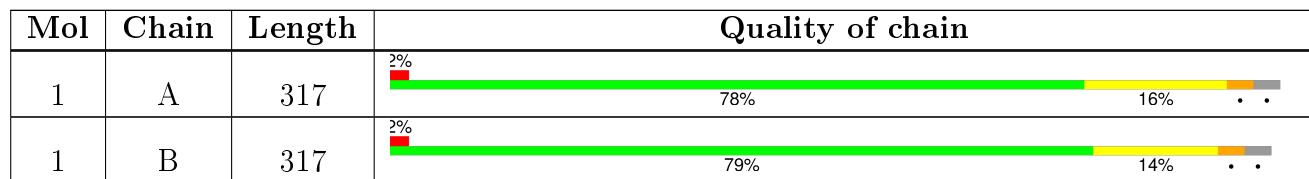
2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5934 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 Endoglucanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	308	Total 2595	C 1689	N 442	O 460	S 4	0	0	0
1	B	306	Total 2579	C 1679	N 440	O 456	S 4	0	0	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	372	Total 372	O 372	0	0
2	B	388	Total 388	O 388	0	0

3 Residue-property plots ⓘ

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: Endoglucanase
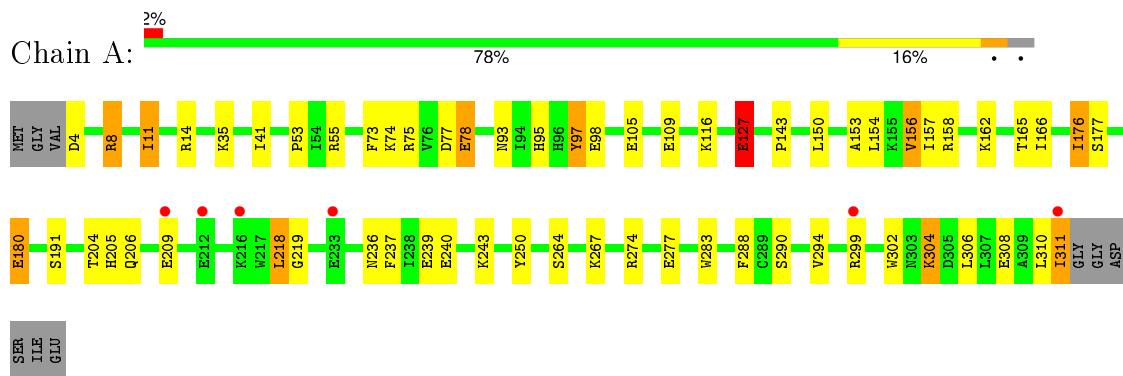
- Molecule 1: Endoglucanase
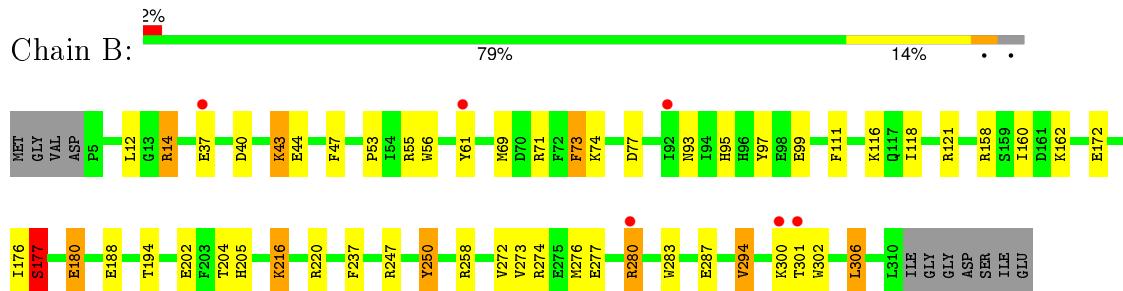
4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	62.92Å 78.34Å 62.98Å 90.00° 97.25° 90.00°	Depositor
Resolution (Å)	25.00 – 1.40 24.42 – 1.40	Depositor EDS
% Data completeness (in resolution range)	97.9 (25.00-1.40) 99.2 (24.42-1.40)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	6.35 (at 1.40Å)	Xtriage
Refinement program	CNS 1.21	Depositor
R, R_{free}	0.184 , 0.209 0.187 , 0.189	Depositor DCC
R_{free} test set	5909 reflections (5.01%)	DCC
Wilson B-factor (Å²)	13.9	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent k_{sol}(e/Å³), B_{sol}(Å²)	0.39 , 42.5	EDS
Estimated twinning fraction	0.017 for l,-k,h	Xtriage
L-test for twinning[2]	$<|L|> = 0.49, <L^2> = 0.32$	Xtriage
Outliers	0 of 118017 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5934	wwPDB-VP
Average B, all atoms (Å²)	15.0	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.*

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[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 (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	1.15	12/2676 (0.4%)	1.13	11/3620 (0.3%)
1	B	1.12	8/2660 (0.3%)	1.12	12/3597 (0.3%)
All	All	1.13	20/5336 (0.4%)	1.13	23/7217 (0.3%)

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	A	0	1
1	B	0	1
All	All	0	2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	61	TYR	CD2-CE2	9.20	1.53	1.39
1	A	78	GLU	CD-OE2	8.49	1.34	1.25
1	A	290	SER	CB-OG	-7.05	1.33	1.42
1	A	98	GLU	CG-CD	7.02	1.62	1.51
1	A	78	GLU	CD-OE1	6.79	1.33	1.25
1	B	177	SER	CA-CB	6.79	1.63	1.52
1	A	274	ARG	CG-CD	6.43	1.68	1.51
1	B	111	PHE	CD2-CE2	6.16	1.51	1.39
1	A	105	GLU	CB-CG	-6.07	1.40	1.52
1	B	202	GLU	CG-CD	5.96	1.60	1.51
1	A	127	GLU	CG-CD	5.95	1.60	1.51
1	A	109	GLU	CG-CD	5.79	1.60	1.51
1	B	61	TYR	CZ-OH	5.58	1.47	1.37
1	B	61	TYR	CE2-CZ	5.51	1.45	1.38
1	B	258	ARG	CZ-NH2	-5.43	1.25	1.33
1	B	177	SER	CB-OG	-5.42	1.35	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	78	GLU	CG-CD	5.36	1.59	1.51
1	A	264	SER	CB-OG	5.19	1.49	1.42
1	A	156	VAL	CB-CG1	-5.17	1.42	1.52
1	A	98	GLU	CB-CG	5.05	1.61	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	274	ARG	NE-CZ-NH1	-11.20	114.70	120.30
1	B	61	TYR	CA-CB-CG	-10.57	93.31	113.40
1	A	77	ASP	CB-CG-OD1	10.00	127.30	118.30
1	B	77	ASP	CB-CG-OD1	8.74	126.17	118.30
1	B	14	ARG	NE-CZ-NH1	8.45	124.53	120.30
1	B	71	ARG	NE-CZ-NH1	8.14	124.37	120.30
1	B	121	ARG	NE-CZ-NH1	-7.99	116.31	120.30
1	A	55	ARG	NE-CZ-NH2	-7.79	116.40	120.30
1	A	310	LEU	C-N-CA	7.35	140.08	121.70
1	A	274	ARG	CA-CB-CG	7.35	129.57	113.40
1	B	177	SER	CA-CB-OG	-7.11	92.01	111.20
1	A	274	ARG	NE-CZ-NH1	6.57	123.58	120.30
1	A	274	ARG	NE-CZ-NH2	-5.99	117.30	120.30
1	B	274	ARG	NE-CZ-NH2	5.96	123.28	120.30
1	B	247	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	B	294	VAL	CG1-CB-CG2	-5.77	101.67	110.90
1	A	288	PHE	CB-CG-CD2	-5.55	116.91	120.80
1	B	306	LEU	CB-CG-CD1	-5.52	101.61	111.00
1	B	55	ARG	NE-CZ-NH2	-5.51	117.55	120.30
1	A	218	LEU	CA-CB-CG	5.39	127.69	115.30
1	A	97	TYR	CB-CG-CD1	-5.10	117.94	121.00
1	A	75	ARG	NE-CZ-NH2	-5.08	117.76	120.30
1	A	267	LYS	CD-CE-NZ	-5.01	100.19	111.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	250	TYR	Sidechain
1	B	250	TYR	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	2595	0	2528	63	0
1	B	2579	0	2514	50	0
2	A	372	0	0	9	0
2	B	388	0	0	11	0
All	All	5934	0	5042	113	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 11.

All (113) 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:14:ARG:HD3	1:A:283:TRP:CZ2	1.74	1.20
1:A:35:LYS:HE2	2:A:679:HOH:O	1.41	1.16
1:B:40:ASP:O	1:B:44:GLU:HG2	1.49	1.12
1:B:43:LYS:HD3	2:B:1160:HOH:O	1.49	1.09
1:A:150:LEU:O	1:A:154:LEU:HD13	1.50	1.09
1:A:14:ARG:HD3	1:A:283:TRP:HZ2	1.08	1.07
1:B:99:GLU:OE2	2:B:751:HOH:O	1.84	0.95
1:B:180:GLU:H	1:B:180:GLU:CD	1.74	0.90
1:B:14:ARG:HD2	2:B:1161:HOH:O	1.72	0.90
1:A:14:ARG:CD	1:A:283:TRP:CZ2	2.56	0.89
1:B:93:ASN:HD21	1:B:95:HIS:HD2	1.20	0.88
1:A:93:ASN:HD21	1:A:95:HIS:HD2	1.21	0.86
1:A:236:ASN:O	1:A:240:GLU:HG2	1.76	0.85
1:A:176:ILE:H	1:A:176:ILE:CD1	1.92	0.82
1:A:150:LEU:O	1:A:154:LEU:CD1	2.28	0.81
1:A:299:ARG:HG2	1:A:299:ARG:HH21	1.45	0.80
1:B:176:ILE:HD11	1:B:237:PHE:CD2	2.19	0.78
1:A:14:ARG:CD	1:A:283:TRP:HZ2	1.94	0.77
1:A:176:ILE:HD12	1:A:176:ILE:H	1.49	0.77
1:A:180:GLU:CD	1:A:180:GLU:H	1.87	0.77
1:A:308:GLU:O	1:A:311:ILE:HG22	1.86	0.76
1:A:127:GLU:CD	1:A:127:GLU:H	1.90	0.75
1:A:8:ARG:HD3	1:A:165:THR:OG1	1.91	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:ARG:NH1	1:A:311:ILE:HD13	2.06	0.70
1:A:41:ILE:HD13	2:A:875:HOH:O	1.92	0.69
1:B:176:ILE:HD11	1:B:237:PHE:HD2	1.53	0.69
1:A:176:ILE:N	1:A:176:ILE:CD1	2.57	0.68
1:B:56:TRP:CZ2	1:B:118:ILE:HD12	2.32	0.65
1:B:180:GLU:HG3	1:B:237:PHE:HZ	1.62	0.65
1:A:299:ARG:HG2	1:A:299:ARG:NH2	2.11	0.64
1:A:153:ALA:O	1:A:156:VAL:HG13	1.98	0.64
1:B:43:LYS:HA	1:B:43:LYS:HE3	1.80	0.64
1:B:43:LYS:HA	1:B:43:LYS:CE	2.27	0.63
1:A:158:ARG:HB3	1:A:162:LYS:HE3	1.81	0.63
1:A:14:ARG:HH12	1:A:311:ILE:HD13	1.62	0.63
1:A:14:ARG:HB3	1:A:283:TRP:CH2	2.33	0.63
1:B:14:ARG:CZ	1:B:277:GLU:OE2	2.48	0.62
1:B:280:ARG:HH11	1:B:280:ARG:HG3	1.65	0.62
1:A:14:ARG:NH2	1:A:277:GLU:HG3	2.13	0.62
1:A:116:LYS:HE2	2:A:1135:HOH:O	2.00	0.61
1:B:43:LYS:CE	1:B:47:PHE:O	2.49	0.61
1:B:43:LYS:HE3	1:B:47:PHE:O	1.99	0.61
1:B:280:ARG:HH11	1:B:280:ARG:CG	2.13	0.61
1:B:43:LYS:NZ	1:B:47:PHE:O	2.34	0.60
1:A:176:ILE:HD13	1:A:177:SER:H	1.65	0.60
1:A:176:ILE:HD13	1:A:177:SER:N	2.17	0.59
1:B:180:GLU:CD	1:B:180:GLU:N	2.53	0.58
1:A:153:ALA:O	1:A:156:VAL:CG1	2.51	0.58
1:B:69:MET:CE	2:B:752:HOH:O	2.51	0.58
1:A:14:ARG:HB3	1:A:283:TRP:CZ2	2.40	0.56
1:B:280:ARG:CG	1:B:280:ARG:NH1	2.67	0.56
1:A:180:GLU:HG3	1:A:237:PHE:HZ	1.71	0.56
1:B:294:VAL:HG12	1:B:302:TRP:HE3	1.71	0.55
1:A:302:TRP:O	1:A:304:LYS:HD2	2.06	0.55
1:A:209:GLU:N	1:A:209:GLU:OE1	2.27	0.54
1:B:172:GLU:OE1	1:B:177:SER:HB3	2.06	0.54
1:B:56:TRP:CH2	1:B:118:ILE:CD1	2.91	0.53
1:A:95:HIS:HE1	2:A:519:HOH:O	1.91	0.53
1:A:74:LYS:O	1:A:78:GLU:HG3	2.07	0.53
1:A:8:ARG:NH2	1:A:11:ILE:HG21	2.24	0.53
1:A:127:GLU:N	1:A:127:GLU:CD	2.63	0.52
1:B:158:ARG:HD2	1:B:188:GLU:OE1	2.11	0.51
1:A:14:ARG:HH21	1:A:277:GLU:CD	2.13	0.51
1:A:206:GLN:NE2	1:A:219:GLY:H	2.08	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:THR:OG1	1:B:205:HIS:HD2	1.94	0.50
1:A:176:ILE:N	1:A:176:ILE:HD13	2.27	0.50
1:B:300:LYS:O	1:B:301:THR:CG2	2.59	0.50
1:B:56:TRP:CZ3	1:B:118:ILE:HD11	2.47	0.50
1:B:56:TRP:CE3	1:B:118:ILE:HD11	2.46	0.50
1:B:158:ARG:NE	2:B:1086:HOH:O	2.46	0.49
1:A:116:LYS:CE	2:A:1135:HOH:O	2.59	0.49
1:A:14:ARG:CG	1:A:283:TRP:CZ2	2.96	0.48
1:B:95:HIS:HE1	2:B:728:HOH:O	1.97	0.48
1:A:53:PRO:HB3	1:A:95:HIS:CD2	2.49	0.47
1:B:56:TRP:CZ2	1:B:118:ILE:CD1	2.97	0.47
1:A:299:ARG:CG	1:A:299:ARG:NH2	2.76	0.47
1:A:204:THR:OG1	1:A:205:HIS:HD2	1.96	0.47
1:B:69:MET:HE1	2:B:752:HOH:O	2.14	0.47
1:B:294:VAL:HG22	1:B:306:LEU:HB3	1.97	0.47
1:B:14:ARG:HD3	1:B:283:TRP:CZ2	2.50	0.46
1:A:14:ARG:HH21	1:A:277:GLU:CG	2.30	0.45
1:B:74:LYS:HB2	1:B:74:LYS:NZ	2.31	0.45
1:A:14:ARG:HH21	1:A:277:GLU:HG3	1.81	0.45
1:B:73:PHE:HE2	1:B:118:ILE:HD13	1.82	0.45
1:B:300:LYS:C	1:B:301:THR:HG23	2.37	0.44
1:B:272:VAL:O	1:B:276:MET:HG3	2.18	0.44
1:A:294:VAL:HG12	1:A:294:VAL:O	2.17	0.44
1:A:93:ASN:HD21	1:A:95:HIS:CD2	2.13	0.44
1:A:153:ALA:HA	1:A:156:VAL:CG1	2.48	0.43
1:B:220:ARG:NH2	2:B:691:HOH:O	2.46	0.43
1:B:53:PRO:HB3	1:B:95:HIS:CD2	2.53	0.43
1:A:116:LYS:NZ	2:A:1135:HOH:O	2.42	0.43
1:B:93:ASN:HD21	1:B:95:HIS:CD2	2.13	0.43
1:A:294:VAL:HG12	1:A:302:TRP:HE3	1.83	0.43
1:B:158:ARG:NH2	1:B:162:LYS:HE3	2.34	0.43
1:A:180:GLU:N	1:A:180:GLU:CD	2.64	0.43
1:A:35:LYS:CE	2:A:679:HOH:O	2.25	0.43
1:B:194:THR:HA	1:B:250:TYR:O	2.19	0.42
1:A:156:VAL:HG13	1:A:157:ILE:N	2.34	0.42
1:A:4:ASP:N	2:A:1005:HOH:O	2.52	0.42
1:B:116:LYS:HG3	1:B:160:ILE:HD11	2.01	0.42
1:A:306:LEU:HD23	1:A:306:LEU:HA	1.94	0.42
1:A:93:ASN:ND2	1:A:95:HIS:HD2	2.03	0.41
1:B:220:ARG:NE	2:B:691:HOH:O	2.48	0.41
1:A:156:VAL:HG13	1:A:157:ILE:H	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:ILE:O	1:A:191:SER:HA	2.21	0.41
1:B:37:GLU:HG3	2:B:618:HOH:O	2.20	0.41
1:A:11:ILE:HD11	2:A:1049:HOH:O	2.21	0.41
1:B:180:GLU:HB2	2:B:689:HOH:O	2.19	0.41
1:B:216:LYS:HB2	1:B:216:LYS:HE2	1.45	0.41
1:A:239:GLU:OE2	1:A:243:LYS:HE2	2.20	0.40
1:A:153:ALA:HA	1:A:156:VAL:HG12	2.04	0.40
1:B:273:VAL:HA	1:B:276:MET:HE2	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	306/317 (96%)	299 (98%)	7 (2%)	0	100	100
1	B	304/317 (96%)	296 (97%)	8 (3%)	0	100	100
All	All	610/634 (96%)	595 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	271/277 (98%)	260 (96%)	11 (4%)	37	6

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	269/277 (97%)	260 (97%)	9 (3%)	45	10
All	All	540/554 (98%)	520 (96%)	20 (4%)	41	9

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

Mol	Chain	Res	Type
1	A	8	ARG
1	A	11	ILE
1	A	73	PHE
1	A	97	TYR
1	A	127	GLU
1	A	143	PRO
1	A	176	ILE
1	A	180	GLU
1	A	218	LEU
1	A	304	LYS
1	A	311	ILE
1	B	12	LEU
1	B	43	LYS
1	B	73	PHE
1	B	97	TYR
1	B	177	SER
1	B	180	GLU
1	B	216	LYS
1	B	280	ARG
1	B	287	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	95	HIS
1	A	205	HIS
1	A	206	GLN
1	A	230	HIS
1	B	59	HIS
1	B	95	HIS
1	B	205	HIS
1	B	206	GLN
1	B	230	HIS
1	B	245	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 (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	308/317 (97%)	0.15	6 (1%)	70	68	8, 14, 22, 28	0
1	B	306/317 (96%)	0.16	6 (1%)	68	67	8, 13, 21, 26	0
All	All	614/634 (96%)	0.15	12 (1%)	68	67	8, 13, 22, 28	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	61	TYR	5.9
1	A	212	GLU	4.0
1	A	311	ILE	3.7
1	B	300	LYS	3.4
1	A	209	GLU	2.6
1	B	280	ARG	2.4
1	A	216	LYS	2.2
1	A	233	GLU	2.2
1	B	37	GLU	2.2
1	B	92	ILE	2.1
1	B	301	THR	2.0
1	A	299	ARG	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.