



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:40 AM GMT

PDB ID : 2BDQ
Title : Crystal Structure of the Putative Copper Homeostasis Protein CutC from Streptococcus agalactiae, Northeast Structural Genomics Target SaR15.
Authors : Forouhar, F.; Abashidze, M.; Jayaraman, S.; Ho, C.K.; Cooper, B.; Acton, T.B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2005-10-20
Resolution : 2.30 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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

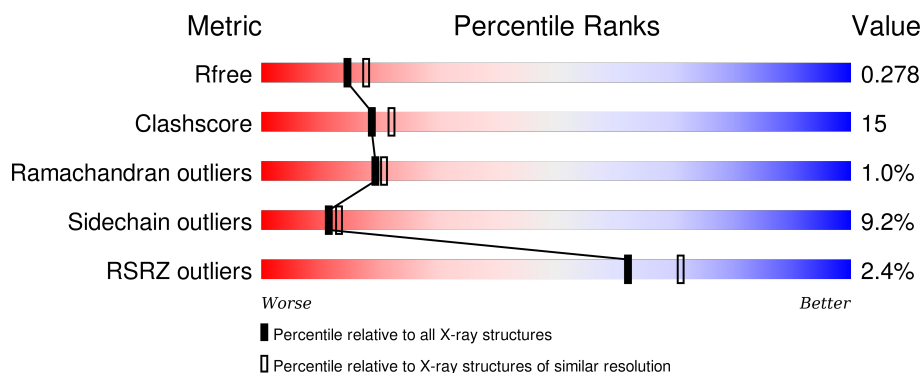
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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 $\leq 5\%$. 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	224	<div> <div>3%</div> <div>67% 21% 6% 6%</div> </div>
1	B	224	<div> <div>%</div> <div>68% 19% 7% 6%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3476 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 copper homeostasis protein CutC.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	210	Total	C	N	O	S	Se	0	0	0
			1630	1028	281	313	3	5			
1	B	210	Total	C	N	O	S	Se	0	0	0
			1630	1028	281	313	3	5			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q8DYB9
A	60	MSE	MET	MODIFIED RESIDUE	UNP Q8DYB9
A	78	MSE	MET	MODIFIED RESIDUE	UNP Q8DYB9
A	125	MSE	MET	MODIFIED RESIDUE	UNP Q8DYB9
A	182	MSE	MET	MODIFIED RESIDUE	UNP Q8DYB9
A	212	MSE	-	SEE REMARK 999	UNP Q8DYB9
A	213	ALA	-	CLONING ARTIFACT	UNP Q8DYB9
A	214	GLY	-	CLONING ARTIFACT	UNP Q8DYB9
A	215	ASP	-	CLONING ARTIFACT	UNP Q8DYB9
A	216	PRO	-	CLONING ARTIFACT	UNP Q8DYB9
A	217	LEU	-	CLONING ARTIFACT	UNP Q8DYB9
A	218	GLU	-	CLONING ARTIFACT	UNP Q8DYB9
A	219	HIS	-	EXPRESSION TAG	UNP Q8DYB9
A	220	HIS	-	EXPRESSION TAG	UNP Q8DYB9
A	221	HIS	-	EXPRESSION TAG	UNP Q8DYB9
A	222	HIS	-	EXPRESSION TAG	UNP Q8DYB9
A	223	HIS	-	EXPRESSION TAG	UNP Q8DYB9
A	224	HIS	-	EXPRESSION TAG	UNP Q8DYB9
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q8DYB9
B	60	MSE	MET	MODIFIED RESIDUE	UNP Q8DYB9
B	78	MSE	MET	MODIFIED RESIDUE	UNP Q8DYB9
B	125	MSE	MET	MODIFIED RESIDUE	UNP Q8DYB9
B	182	MSE	MET	MODIFIED RESIDUE	UNP Q8DYB9
B	212	MSE	-	SEE REMARK 999	UNP Q8DYB9
B	213	ALA	-	CLONING ARTIFACT	UNP Q8DYB9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	214	GLY	-	CLONING ARTIFACT	UNP Q8DYB9
B	215	ASP	-	CLONING ARTIFACT	UNP Q8DYB9
B	216	PRO	-	CLONING ARTIFACT	UNP Q8DYB9
B	217	LEU	-	CLONING ARTIFACT	UNP Q8DYB9
B	218	GLU	-	CLONING ARTIFACT	UNP Q8DYB9
B	219	HIS	-	EXPRESSION TAG	UNP Q8DYB9
B	220	HIS	-	EXPRESSION TAG	UNP Q8DYB9
B	221	HIS	-	EXPRESSION TAG	UNP Q8DYB9
B	222	HIS	-	EXPRESSION TAG	UNP Q8DYB9
B	223	HIS	-	EXPRESSION TAG	UNP Q8DYB9
B	224	HIS	-	EXPRESSION TAG	UNP Q8DYB9

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	99	Total O 99 99	0	0
2	B	117	Total O 117 117	0	0

- Molecule 1: copper homeostasis protein CutC



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	48.46 Å 71.54 Å 128.19 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.25 – 2.30 47.74 – 2.29	Depositor EDS
% Data completeness (in resolution range)	94.1 (29.25-2.30) 97.9 (47.74-2.29)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.60 (at 2.29 Å)	Xtriage
Refinement program	XTALVIEW, CNS 1.1	Depositor
R, R_{free}	0.221 , 0.271 0.232 , 0.278	Depositor DCC
R_{free} test set	1998 reflections (11.12%)	DCC
Wilson B-factor (Å ²)	21.3	Xtriage
Anisotropy	0.393	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 38496 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3476	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.31% 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, 0.375 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	0.37	1/1648 (0.1%)	0.59	0/2224
1	B	0.37	1/1648 (0.1%)	0.59	0/2224
All	All	0.37	2/3296 (0.1%)	0.59	0/4448

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	159	GLU	CD-OE2	6.90	1.33	1.25
1	B	159	GLU	CD-OE2	6.65	1.32	1.25

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1630	0	1658	50	0
1	B	1630	0	1658	53	0
2	A	99	0	0	3	0
2	B	117	0	0	8	0
All	All	3476	0	3316	101	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 15.

All (101) 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:28:CYS:HB3	1:A:60:MSE:HB2	1.64	0.79
1:B:28:CYS:HB3	1:B:60:MSE:HB2	1.65	0.79
1:A:71:ASN:ND2	1:A:74:GLU:H	1.82	0.76
1:B:71:ASN:ND2	1:B:74:GLU:H	1.85	0.75
1:A:77:ILE:HD11	1:B:11:LEU:HD21	1.70	0.74
1:A:71:ASN:C	1:A:71:ASN:HD22	1.91	0.74
1:B:50:LEU:HB3	1:B:55:ILE:HG23	1.70	0.74
1:B:71:ASN:C	1:B:71:ASN:HD22	1.91	0.73
1:A:50:LEU:HB3	1:A:55:ILE:HG23	1.70	0.73
1:B:191:ASN:HD22	1:B:192:TYR:N	1.90	0.70
1:A:25:VAL:HG11	1:A:50:LEU:HD13	1.74	0.67
1:A:191:ASN:HD22	1:A:192:TYR:N	1.92	0.67
1:B:25:VAL:HG11	1:B:50:LEU:HD13	1.74	0.67
1:B:25:VAL:CG1	1:B:50:LEU:HD13	2.27	0.65
1:A:25:VAL:CG1	1:A:50:LEU:HD13	2.27	0.64
1:A:28:CYS:HA	1:A:38:PRO:HD3	1.80	0.63
1:B:28:CYS:HA	1:B:38:PRO:HD3	1.81	0.61
1:A:83:LEU:HD13	1:A:115:ALA:HB1	1.83	0.59
1:B:83:LEU:HD13	1:B:115:ALA:HB1	1.85	0.59
1:A:105:ASP:OD1	1:A:107:GLU:HG2	2.02	0.59
1:B:209:ILE:HG23	1:B:210:THR:HG23	1.86	0.58
1:A:64:ARG:HD2	1:A:65:GLY:O	2.03	0.58
1:B:105:ASP:OD1	1:B:107:GLU:HG2	2.04	0.58
1:B:33:VAL:CG2	1:B:65:GLY:HA3	2.35	0.57
1:A:209:ILE:HG23	1:A:210:THR:HG23	1.85	0.57
1:B:64:ARG:HD2	1:B:65:GLY:O	2.05	0.57
1:A:28:CYS:HB3	1:A:60:MSE:CB	2.35	0.57
1:A:33:VAL:HG21	2:A:272:HOH:O	2.05	0.56
1:A:11:LEU:HD21	1:B:77:ILE:HD11	1.88	0.56
1:B:28:CYS:HB3	1:B:60:MSE:CB	2.34	0.56
1:A:33:VAL:CG2	1:A:65:GLY:HA3	2.36	0.55
1:B:188:THR:H	1:B:191:ASN:HD21	1.54	0.55
1:B:33:VAL:HG21	1:B:65:GLY:HA3	1.89	0.55
1:B:23:SER:HB3	2:B:309:HOH:O	2.06	0.54
1:B:161:ILE:HD13	1:B:161:ILE:O	2.08	0.54
1:A:161:ILE:HD13	1:A:161:ILE:O	2.08	0.54
1:A:106:THR:HG23	1:A:145:LEU:HD21	1.90	0.54
1:B:25:VAL:HG11	1:B:50:LEU:CD1	2.38	0.54
1:A:188:THR:H	1:A:191:ASN:HD21	1.55	0.53
1:B:47:ASN:ND2	1:B:57:VAL:HG11	2.25	0.52
1:A:33:VAL:HG21	1:A:65:GLY:HA3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:VAL:HG11	1:A:50:LEU:CD1	2.38	0.52
1:A:47:ASN:ND2	1:A:57:VAL:HG11	2.25	0.52
1:B:202:LYS:HD2	2:B:275:HOH:O	2.10	0.51
1:A:71:ASN:HD21	1:A:74:GLU:H	1.54	0.51
1:B:71:ASN:HD21	1:B:74:GLU:H	1.58	0.50
1:B:93:LEU:HD12	1:B:116:THR:HG23	1.92	0.50
1:B:106:THR:HG23	1:B:145:LEU:HD21	1.92	0.50
1:A:116:THR:HG21	1:A:121:LEU:HD21	1.93	0.50
1:A:86:VAL:HG11	1:A:117:GLN:NE2	2.27	0.50
1:B:166:LYS:HD2	1:B:166:LYS:N	2.26	0.50
1:B:33:VAL:HG23	2:B:233:HOH:O	2.11	0.49
1:A:166:LYS:HD2	1:A:166:LYS:N	2.27	0.49
1:A:93:LEU:HD12	1:A:116:THR:HG23	1.94	0.49
1:A:21:ILE:HD11	2:A:283:HOH:O	2.13	0.49
1:B:116:THR:HG21	1:B:121:LEU:HD21	1.94	0.48
1:A:71:ASN:ND2	1:A:71:ASN:C	2.65	0.48
1:B:106:THR:O	1:B:110:GLU:HG3	2.13	0.48
1:B:137:LYS:HE3	2:B:299:HOH:O	2.13	0.48
1:B:23:SER:HA	2:B:256:HOH:O	2.14	0.48
1:B:71:ASN:C	1:B:71:ASN:ND2	2.64	0.48
1:B:86:VAL:HG11	1:B:117:GLN:NE2	2.30	0.47
1:B:73:LEU:O	1:B:73:LEU:HD23	2.14	0.47
1:B:71:ASN:HD22	1:B:74:GLU:H	1.62	0.47
1:A:106:THR:O	1:A:110:GLU:HG3	2.14	0.47
1:A:94:VAL:HA	1:A:122:VAL:HG13	1.98	0.46
1:A:71:ASN:HD22	1:A:74:GLU:H	1.61	0.46
1:B:94:VAL:HA	1:B:122:VAL:HG13	1.98	0.46
1:A:40:TYR:HE1	1:A:84:ARG:HG3	1.81	0.45
1:B:188:THR:H	1:B:191:ASN:ND2	2.13	0.45
1:A:73:LEU:O	1:A:73:LEU:HD23	2.17	0.45
1:A:188:THR:H	1:A:191:ASN:ND2	2.14	0.45
1:B:191:ASN:O	1:B:195:ILE:HD13	2.17	0.45
1:B:16:ARG:HB2	2:B:311:HOH:O	2.16	0.45
1:B:94:VAL:HG22	1:B:122:VAL:CG1	2.47	0.45
1:A:191:ASN:O	1:A:195:ILE:HD13	2.16	0.44
1:A:27:LEU:HD22	1:A:38:PRO:HG3	2.00	0.44
1:A:94:VAL:HG22	1:A:122:VAL:CG1	2.48	0.44
1:A:191:ASN:HD22	1:A:191:ASN:C	2.17	0.44
1:A:157:ASN:HB3	1:A:159:GLU:CD	2.39	0.43
1:B:82:ILE:O	1:B:86:VAL:HG12	2.18	0.43
1:A:6:PHE:HB2	1:A:22:ILE:HD13	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:TYR:HE1	1:B:84:ARG:HG3	1.83	0.43
1:A:162:ILE:HG23	1:A:163:GLU:HG2	2.01	0.43
1:B:125:MSE:HB2	2:B:235:HOH:O	2.19	0.42
1:B:116:THR:HG22	1:B:116:THR:O	2.18	0.42
1:A:173:GLU:HB3	2:A:297:HOH:O	2.18	0.42
1:A:33:VAL:HG22	1:A:65:GLY:HA3	2.01	0.42
1:A:1:MSE:HB3	1:A:2:ILE:H	1.66	0.42
1:B:157:ASN:HB3	1:B:159:GLU:CD	2.39	0.42
1:A:50:LEU:CD1	1:A:57:VAL:HG22	2.49	0.42
1:B:67:ASN:HB3	1:B:69:VAL:H	1.85	0.42
1:B:137:LYS:H	1:B:137:LYS:HZ3	1.68	0.41
1:B:35:GLY:O	1:B:62:ARG:HD3	2.20	0.41
1:A:31:LEU:HD23	1:A:31:LEU:HA	1.92	0.41
1:B:191:ASN:C	1:B:191:ASN:HD22	2.15	0.41
1:A:116:THR:CG2	1:A:121:LEU:HD21	2.50	0.41
1:B:50:LEU:CD1	1:B:57:VAL:HG22	2.50	0.41
1:B:33:VAL:HG22	1:B:65:GLY:HA3	2.02	0.41
1:B:52:GLU:HG2	2:B:260:HOH:O	2.20	0.40
1:A:35:GLY:O	1:A:62:ARG:HD3	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	208/224 (93%)	199 (96%)	7 (3%)	2 (1%)	19	21
1	B	208/224 (93%)	199 (96%)	7 (3%)	2 (1%)	19	21
All	All	416/448 (93%)	398 (96%)	14 (3%)	4 (1%)	19	21

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	ASN
1	B	67	ASN
1	A	125	MSE
1	B	125	MSE

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	179/185 (97%)	162 (90%)	17 (10%)	11	12
1	B	179/185 (97%)	163 (91%)	16 (9%)	12	14
All	All	358/370 (97%)	325 (91%)	33 (9%)	11	13

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	25	VAL
1	A	27	LEU
1	A	60	MSE
1	A	64	ARG
1	A	71	ASN
1	A	83	LEU
1	A	84	ARG
1	A	101	ASN
1	A	122	VAL
1	A	135	GLN
1	A	137	LYS
1	A	161	ILE
1	A	163	GLU
1	A	166	LYS
1	A	191	ASN
1	A	202	LYS
1	B	25	VAL
1	B	27	LEU
1	B	60	MSE

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Mol	Chain	Res	Type
1	B	64	ARG
1	B	71	ASN
1	B	83	LEU
1	B	84	ARG
1	B	101	ASN
1	B	122	VAL
1	B	135	GLN
1	B	137	LYS
1	B	161	ILE
1	B	163	GLU
1	B	166	LYS
1	B	191	ASN
1	B	202	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	47	ASN
1	A	67	ASN
1	A	71	ASN
1	A	102	ASN
1	A	117	GLN
1	A	135	GLN
1	A	141	GLN
1	A	167	HIS
1	A	191	ASN
1	A	197	GLN
1	B	10	ASN
1	B	47	ASN
1	B	67	ASN
1	B	71	ASN
1	B	102	ASN
1	B	117	GLN
1	B	141	GLN
1	B	167	HIS
1	B	191	ASN
1	B	197	GLN

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	205/224 (91%)	0.10	7 (3%) 49 58	13, 21, 36, 50	0
1	B	205/224 (91%)	0.17	3 (1%) 76 81	12, 20, 35, 50	0
All	All	410/448 (91%)	0.13	10 (2%) 62 71	12, 20, 36, 50	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	16	ARG	4.9
1	B	16	ARG	3.8
1	A	157	ASN	3.6
1	A	20	ALA	2.9
1	A	18	ASP	2.8
1	B	55	ILE	2.4
1	A	210	THR	2.3
1	B	208	ARG	2.3
1	A	208	ARG	2.2
1	A	209	ILE	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 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.