



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 11:15 pm BST

PDB ID : 3PWE
Title : Crystal structure of the E. coli beta clamp mutant R103C, I305C, C260S, C333S at 2.2Å resolution
Authors : Marzahn, M.R.; Robbins, A.H.; McKenna, R.; Bloom, L.B.
Deposited on : 2010-12-08
Resolution : 2.20 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 ⓘ](#)) 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

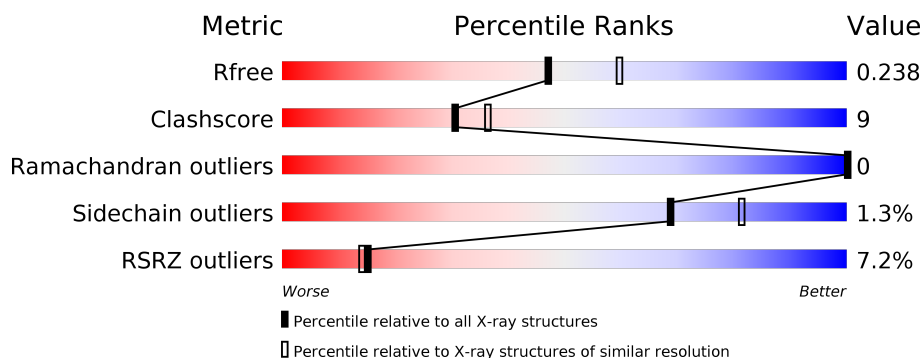
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 $\leq 5\%$. 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	366	<div> <div>6%</div> <div> <div></div> <div>81%</div> <div>19%</div> </div> </div>
1	B	366	<div> <div>8%</div> <div> <div></div> <div>81%</div> <div>19%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6026 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 DNA polymerase III subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	366	Total	C	N	O	S	0	1	0
			2844	1785	496	544	19			
1	B	366	Total	C	N	O	S	0	1	0
			2847	1786	498	544	19			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	103	CYS	ARG	ENGINEERED MUTATION	UNP P0A988
A	260	SER	CYS	ENGINEERED MUTATION	UNP P0A988
A	305	CYS	ILE	ENGINEERED MUTATION	UNP P0A988
A	333	SER	CYS	ENGINEERED MUTATION	UNP P0A988
B	103	CYS	ARG	ENGINEERED MUTATION	UNP P0A988
B	260	SER	CYS	ENGINEERED MUTATION	UNP P0A988
B	305	CYS	ILE	ENGINEERED MUTATION	UNP P0A988
B	333	SER	CYS	ENGINEERED MUTATION	UNP P0A988

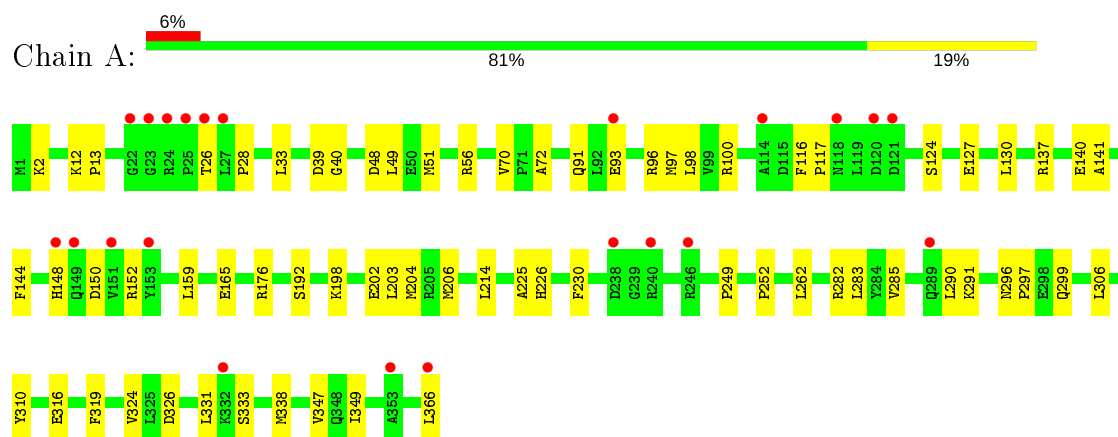
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	160	Total	O	0	0
			160	160		
2	B	175	Total	O	0	0
			175	175		

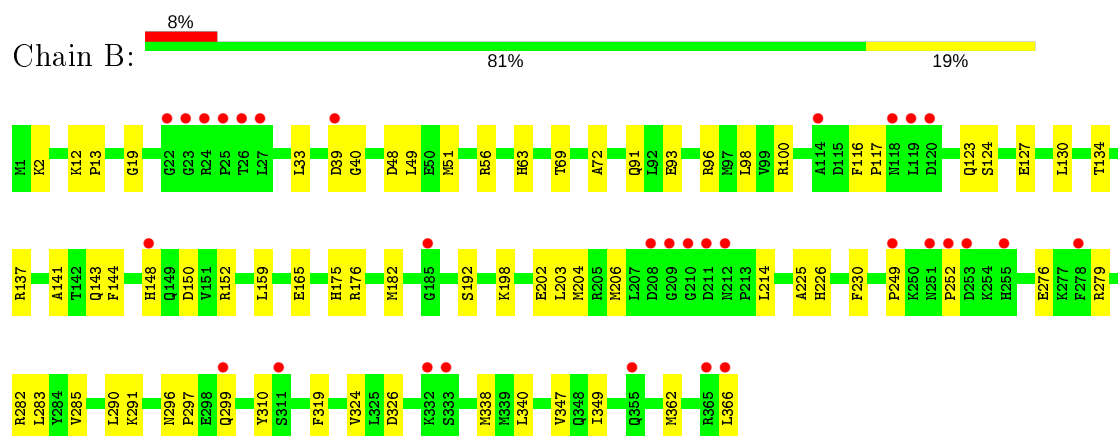
3 Residue-property plots [i](#)

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: DNA polymerase III subunit beta



- Molecule 1: DNA polymerase III subunit beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.80 Å 67.36 Å 80.74 Å 90.00° 114.16° 90.00°	Depositor
Resolution (Å)	28.50 – 2.20 28.50 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.6 (28.50-2.20) 99.4 (28.50-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.61 (at 2.20 Å)	Xtriage
Refinement program	PHENIX 1.6.3_467	Depositor
R, R_{free}	0.204 , 0.245 0.198 , 0.238	Depositor DCC
R_{free} test set	2006 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.546	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 54.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6026	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.52% 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, 0.333 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	0.35	0/2893	0.51	0/3916
1	B	0.35	0/2897	0.50	0/3920
All	All	0.35	0/5790	0.50	0/7836

There are no bond length outliers.

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	2844	0	2855	55	0
1	B	2847	0	2859	53	0
2	A	160	0	0	3	0
2	B	175	0	0	3	0
All	All	6026	0	5714	105	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 9.

All (105) 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:283:LEU:CD1	1:A:290:LEU:HD11	1.95	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:LEU:CD1	1:B:290:LEU:HD11	1.97	0.95
1:A:283:LEU:HD11	1:A:290:LEU:HD11	1.56	0.87
1:B:137:ARG:NH1	1:B:141:ALA:HB2	1.90	0.86
1:B:283:LEU:HD11	1:B:290:LEU:HD11	1.59	0.85
1:A:137:ARG:NH1	1:A:141:ALA:HB2	1.94	0.83
1:B:282:ARG:CG	1:B:282:ARG:HH11	1.93	0.82
1:A:338:MET:HE3	1:A:347:VAL:HG11	1.62	0.81
1:A:282:ARG:HH11	1:A:282:ARG:HG3	1.46	0.80
1:B:338:MET:HE3	1:B:347:VAL:HG11	1.64	0.80
1:B:282:ARG:HG3	1:B:282:ARG:HH11	1.46	0.79
1:B:214:LEU:HD11	1:B:225:ALA:HB1	1.64	0.79
1:A:282:ARG:HH11	1:A:282:ARG:CG	1.94	0.78
1:B:39:ASP:CG	1:B:40:GLY:H	1.84	0.77
1:A:214:LEU:HD11	1:A:225:ALA:HB1	1.67	0.77
1:A:39:ASP:CG	1:A:40:GLY:H	1.86	0.77
1:B:150:ASP:OD1	1:B:152:ARG:HD3	1.93	0.69
1:A:283:LEU:HD12	1:A:290:LEU:HD11	1.73	0.68
1:B:39:ASP:CG	1:B:40:GLY:N	2.47	0.68
1:B:283:LEU:HD12	1:B:290:LEU:HD11	1.76	0.67
1:A:39:ASP:CG	1:A:40:GLY:N	2.49	0.66
1:B:175:HIS:HB3	1:B:362:MET:HE2	1.79	0.65
1:A:150:ASP:OD1	1:A:152:ARG:HD3	1.98	0.64
1:A:140:GLU:HG3	1:A:204:MET:CE	2.26	0.64
1:A:51:MET:CE	1:A:202:GLU:HG3	2.28	0.63
1:A:282:ARG:HG3	1:A:282:ARG:NH1	2.13	0.63
1:B:51:MET:CE	1:B:202:GLU:HG3	2.28	0.63
1:A:319:PHE:CE2	1:A:324:VAL:HG21	2.34	0.63
1:B:51:MET:HE1	1:B:198:LYS:HG2	1.81	0.61
1:B:319:PHE:CE2	1:B:324:VAL:HG21	2.36	0.61
1:B:282:ARG:NH1	1:B:282:ARG:HG3	2.13	0.60
1:A:51:MET:HE1	1:A:198:LYS:HG2	1.85	0.58
1:A:51:MET:HE2	1:A:198:LYS:HE3	1.86	0.58
1:B:51:MET:HE2	1:B:198:LYS:HE3	1.87	0.56
1:A:159:LEU:HD11	1:A:192:SER:HB3	1.89	0.55
1:A:96:ARG:HH21	1:B:299:GLN:HB2	1.71	0.55
1:A:165:GLU:O	1:A:165:GLU:HG3	2.07	0.54
1:B:143:GLN:NE2	1:B:204:MET:SD	2.80	0.54
1:B:159:LEU:HD11	1:B:192:SER:HB3	1.89	0.53
1:B:165:GLU:O	1:B:165:GLU:HG3	2.09	0.52
1:A:2:LYS:HG3	2:A:460:HOH:O	2.10	0.52
1:A:140:GLU:HG3	1:A:204:MET:HE2	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:LEU:O	1:B:206:MET:HG2	2.10	0.51
1:B:282:ARG:NH1	1:B:366:LEU:HB2	2.26	0.50
1:A:140:GLU:HG3	1:A:204:MET:HE1	1.94	0.49
1:B:276:GLU:HA	1:B:279[A]:ARG:HH21	1.78	0.49
1:A:282:ARG:NH1	1:A:366:LEU:HB2	2.28	0.48
1:B:296:ASN:HB2	1:B:297:PRO:CD	2.42	0.48
1:B:51:MET:HE3	1:B:202:GLU:HG3	1.94	0.48
1:A:12:LYS:HB3	1:A:13:PRO:HD3	1.95	0.48
1:B:134:THR:HG23	1:B:182:MET:HE2	1.96	0.48
1:B:282:ARG:HH11	1:B:366:LEU:HD12	1.79	0.48
1:A:296:ASN:HB2	1:A:297:PRO:CD	2.45	0.47
1:A:130:LEU:N	1:A:130:LEU:HD23	2.29	0.47
1:B:93:GLU:OE2	1:B:98:LEU:HD22	2.13	0.47
1:A:93:GLU:OE2	1:A:98:LEU:HD22	2.14	0.47
1:B:124:SER:HG	1:B:127:GLU:HG2	1.80	0.47
1:B:63:HIS:HE1	2:B:488:HOH:O	1.98	0.47
1:B:175:HIS:O	1:B:362:MET:HG2	2.15	0.47
1:B:338:MET:HG2	1:B:349:ILE:HG12	1.97	0.46
1:A:249:PRO:HB2	1:A:252:PRO:HG3	1.97	0.46
1:B:249:PRO:HB2	1:B:252:PRO:HG3	1.98	0.46
1:A:282:ARG:HH11	1:A:366:LEU:HD12	1.81	0.46
1:A:124:SER:HG	1:A:127:GLU:HG2	1.81	0.46
1:A:51:MET:HE1	1:A:202:GLU:HG3	1.98	0.45
1:A:203:LEU:O	1:A:206:MET:HG2	2.17	0.45
1:A:48:ASP:O	1:A:49:LEU:HB2	2.17	0.45
1:A:299:GLN:HB2	1:B:96:ARG:HH21	1.82	0.44
1:A:338:MET:HG2	1:A:349:ILE:HG12	2.00	0.44
1:A:12:LYS:HE2	2:A:491:HOH:O	2.17	0.43
1:A:331:LEU:HA	1:A:331:LEU:HD23	1.86	0.43
1:B:130:LEU:HD23	1:B:130:LEU:N	2.33	0.43
1:A:96:ARG:NH2	1:B:299:GLN:HB2	2.33	0.43
1:B:33:LEU:HG	1:B:72:ALA:HB2	2.01	0.43
1:A:285:VAL:HG21	1:A:310:TYR:CD1	2.53	0.43
1:A:70:VAL:HG11	1:A:97:MET:CE	2.49	0.43
1:A:51:MET:HE3	1:A:202:GLU:HG3	2.00	0.42
1:B:285:VAL:HG21	1:B:310:TYR:CD1	2.54	0.42
1:B:19:GLY:HA3	2:B:444:HOH:O	2.19	0.42
1:A:51:MET:CE	1:A:198:LYS:HG2	2.49	0.42
1:A:283:LEU:O	1:A:316:GLU:HA	2.19	0.42
1:B:124:SER:HG	1:B:127:GLU:CG	2.32	0.42
1:A:333:SER:HB2	2:A:414:HOH:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:LEU:HG	1:A:72:ALA:HB2	2.02	0.42
1:B:123:GLN:NE2	2:B:468:HOH:O	2.51	0.42
1:B:116:PHE:HA	1:B:117:PRO:HD3	1.89	0.42
1:B:12:LYS:HB3	1:B:13:PRO:HD3	2.02	0.42
1:B:144:PHE:CD2	1:B:326:ASP:HB3	2.54	0.42
1:A:98:LEU:HD23	1:A:100:ARG:NH2	2.35	0.41
1:A:116:PHE:HA	1:A:117:PRO:HD3	1.89	0.41
1:A:124:SER:HG	1:A:127:GLU:CG	2.33	0.41
1:B:33:LEU:O	1:B:69:THR:HA	2.19	0.41
1:A:26:THR:O	1:A:28:PRO:HD3	2.21	0.41
1:B:13:PRO:HA	1:B:230:PHE:HE1	1.84	0.41
1:B:290:LEU:HD12	1:B:291:LYS:N	2.35	0.41
1:A:2:LYS:NZ	1:A:91:GLN:HE21	2.19	0.41
1:A:290:LEU:HD12	1:A:291:LYS:N	2.36	0.41
1:A:144:PHE:CD2	1:A:326:ASP:HB3	2.56	0.41
1:B:48:ASP:O	1:B:49:LEU:HB2	2.21	0.41
1:B:2:LYS:NZ	1:B:91:GLN:HE21	2.19	0.41
1:A:262:LEU:HG	1:A:306:LEU:HD11	2.02	0.41
1:B:98:LEU:HD23	1:B:100:ARG:NH2	2.35	0.40
1:B:51:MET:CE	1:B:198:LYS:HG2	2.48	0.40
1:A:13:PRO:HA	1:A:230:PHE:HE1	1.86	0.40
1:B:340:LEU:HD23	1:B:347:VAL:HG22	2.02	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	365/366 (100%)	356 (98%)	9 (2%)	0	100	100
1	B	365/366 (100%)	356 (98%)	9 (2%)	0	100	100
All	All	730/732 (100%)	712 (98%)	18 (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	314/313 (100%)	310 (99%)	4 (1%)	69	81
1	B	314/313 (100%)	310 (99%)	4 (1%)	69	81
All	All	628/626 (100%)	620 (99%)	8 (1%)	69	81

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ARG
1	A	148	HIS
1	A	176	ARG
1	A	226	HIS
1	B	56	ARG
1	B	148	HIS
1	B	176	ARG
1	B	226	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	143	GLN
1	B	91	GLN
1	B	123	GLN

5.3.3 RNA ⓘ

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 ⓘ

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, 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	366/366 (100%)	0.31	22 (6%)	21 20	16, 28, 51, 96	0
1	B	366/366 (100%)	0.38	31 (8%)	10 9	15, 28, 51, 93	0
All	All	732/732 (100%)	0.34	53 (7%)	15 14	15, 28, 51, 96	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	23	GLY	14.0
1	A	24	ARG	9.4
1	B	366	LEU	8.9
1	A	26	THR	7.8
1	A	25	PRO	6.5
1	B	251	ASN	6.1
1	A	240	ARG	5.8
1	B	365	ARG	5.7
1	B	211	ASP	5.7
1	A	22	GLY	5.4
1	B	23	GLY	5.0
1	B	26	THR	5.0
1	B	210	GLY	4.8
1	B	148	HIS	4.5
1	B	25	PRO	4.3
1	B	208	ASP	4.2
1	B	311	SER	4.2
1	B	24	ARG	4.1
1	A	149	GLN	4.0
1	A	120	ASP	4.0
1	B	355	GLN	3.8
1	B	212	ASN	3.8
1	B	249	PRO	3.8
1	A	153	TYR	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	148	HIS	3.7
1	A	118	ASN	3.7
1	B	22	GLY	3.6
1	A	114	ALA	3.3
1	A	27	LEU	3.3
1	B	120	ASP	3.3
1	B	253	ASP	3.2
1	B	27	LEU	3.1
1	A	151	VAL	3.0
1	A	353	ALA	3.0
1	B	209	GLY	2.9
1	A	366	LEU	2.9
1	B	119	LEU	2.8
1	A	238	ASP	2.8
1	A	289	GLN	2.7
1	A	246	ARG	2.7
1	A	121	ASP	2.6
1	B	252	PRO	2.5
1	B	185	GLY	2.5
1	B	118	ASN	2.5
1	A	332	LYS	2.4
1	B	333	SER	2.3
1	B	255	HIS	2.3
1	B	39	ASP	2.2
1	B	278	PHE	2.1
1	B	114	ALA	2.1
1	B	332	LYS	2.1
1	B	299	GLN	2.1
1	A	93	GLU	2.0

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.