



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:43 PM UTC

PDB ID : 2VH1 / pdb_00002vh1
Title : Crystal structure of bacterial cell division protein FtsQ from E.coli
Authors : van den Ent, F.; Vinkenvleugel, T.; Ind, A.; West, P.; Veprintsev, D.; Naninga, N.; den Blaauwen, T.; Lowe, J.
Deposited on : 2007-11-16
Resolution : 2.70 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

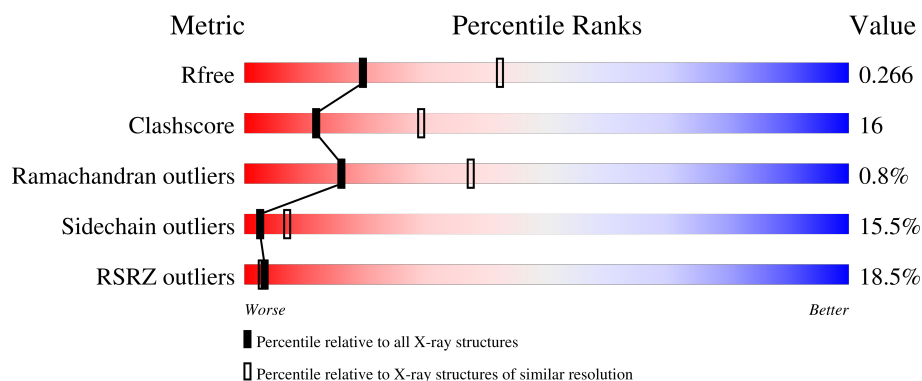
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 of 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	220	
1	B	220	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3234 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.

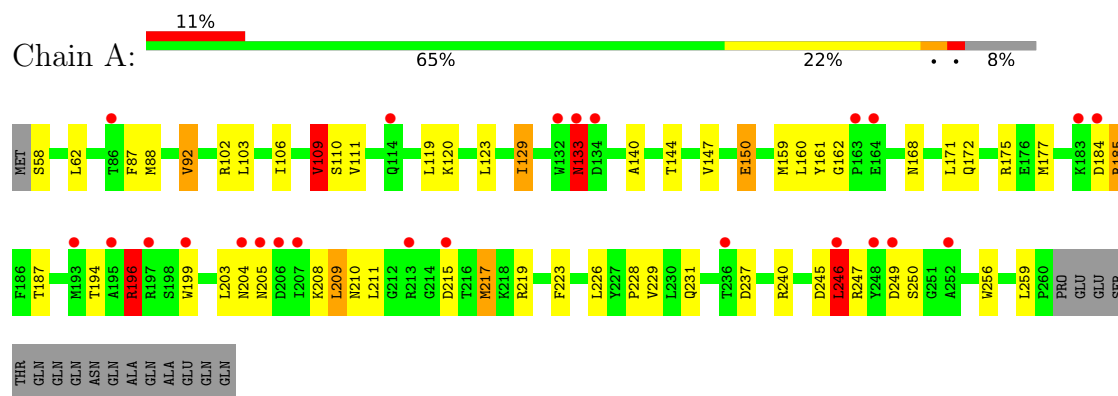
- Molecule 1 is a protein called CELL DIVISION PROTEIN FTSQ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	0	0	0
			1640	1034	296	303	7			
1	B	198	Total	C	N	O	S	0	0	1
			1594	1005	287	295	7			

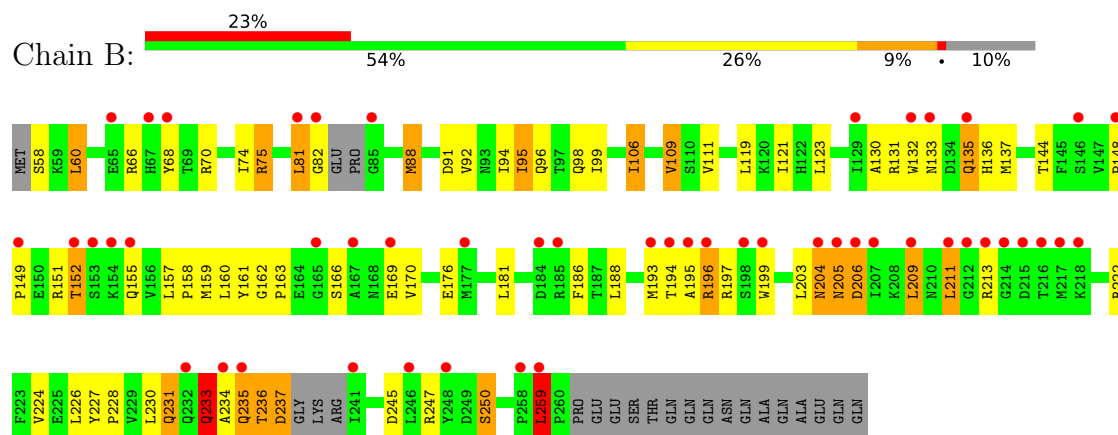
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: CELL DIVISION PROTEIN FTSQ



• Molecule 1: CELL DIVISION PROTEIN FTSQ



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	147.84Å 147.84Å 69.26Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.35 – 2.70 36.35 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.0 (36.35-2.70) 97.8 (36.35-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.3.0037	Depositor
R, R_{free}	0.237 , 0.288 0.225 , 0.266	Depositor DCC
R_{free} test set	6606 reflections (28.24%)	wwPDB-VP
Wilson B-factor (Å ²)	64.6	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 63.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.038 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	3234	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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 [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	0.73	0/1674	1.04	5/2268 (0.2%)
1	B	0.88	4/1624 (0.2%)	1.04	2/2198 (0.1%)
All	All	0.81	4/3298 (0.1%)	1.04	7/4466 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	82	GLY	C-O	17.48	1.58	1.23
1	B	233	GLN	CD-NE2	9.13	1.52	1.33
1	B	259	LEU	C-N	-6.44	1.24	1.34
1	B	233	GLN	CD-OE1	6.32	1.35	1.23

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	109	VAL	CB-CA-C	-7.16	99.47	110.50
1	A	196	ARG	N-CA-C	-6.87	101.46	110.53
1	A	133	ASN	CA-C-N	5.37	131.37	121.70
1	A	133	ASN	C-N-CA	5.37	131.37	121.70
1	B	205	ASN	N-CA-C	5.32	122.14	110.80
1	A	246	LEU	CA-CB-CG	5.17	134.41	116.30
1	B	233	GLN	OE1-CD-NE2	5.04	127.64	122.60

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	1640	0	1643	42	0
1	B	1594	0	1592	62	0
All	All	3234	0	3235	102	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 16.

All (102) 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:B:205:ASN:H	1:B:206:ASP:HA	1.12	1.10
1:B:205:ASN:N	1:B:206:ASP:HA	1.79	0.95
1:B:132:TRP:O	1:B:136:HIS:HB2	1.72	0.90
1:B:204:ASN:H	1:B:204:ASN:HD22	1.22	0.82
1:B:233:GLN:HB3	1:B:237:ASP:HB2	1.61	0.82
1:A:129:ILE:HD11	1:A:171:LEU:HD11	1.62	0.82
1:B:169:GLU:CD	1:B:197:ARG:HH12	1.87	0.82
1:A:133:ASN:ND2	1:A:162:GLY:H	1.82	0.77
1:B:133:ASN:HD21	1:B:162:GLY:H	1.33	0.76
1:B:133:ASN:ND2	1:B:162:GLY:H	1.88	0.71
1:A:223:PHE:HB2	1:A:246:LEU:HD21	1.72	0.70
1:B:194:THR:HG22	1:B:195:ALA:N	2.07	0.70
1:B:66:ARG:HD3	1:B:123:LEU:HD12	1.73	0.69
1:B:245:ASP:OD2	1:B:247:ARG:HB3	1.92	0.69
1:A:133:ASN:HD22	1:A:162:GLY:H	1.40	0.68
1:A:194:THR:O	1:A:196:ARG:O	2.12	0.68
1:A:106:ILE:HG21	1:A:109:VAL:HG22	1.74	0.68
1:A:92:VAL:HG22	1:A:111:VAL:HG12	1.75	0.67
1:B:194:THR:HG22	1:B:195:ALA:H	1.61	0.65
1:A:129:ILE:CD1	1:A:171:LEU:HD11	2.26	0.65
1:B:204:ASN:HD22	1:B:204:ASN:N	1.96	0.63
1:B:206:ASP:OD1	1:B:206:ASP:N	2.32	0.61
1:A:109:VAL:HG13	1:A:123:LEU:HD23	1.84	0.60
1:B:132:TRP:O	1:B:136:HIS:CB	2.49	0.59
1:A:187:THR:HB	1:A:204:ASN:HB3	1.85	0.58
1:A:245:ASP:OD1	1:A:247:ARG:HB2	2.04	0.58
1:B:106:ILE:HG21	1:B:109:VAL:HG22	1.85	0.58
1:A:219:ARG:HD3	1:A:246:LEU:O	2.04	0.57
1:B:205:ASN:N	1:B:206:ASP:CA	2.63	0.57
1:A:140:ALA:O	1:A:175:ARG:NH1	2.38	0.57
1:A:177:MET:CG	1:A:217:MET:HE1	2.35	0.56
1:A:177:MET:HG3	1:A:217:MET:HE1	1.85	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:GLN:HG2	1:B:234:ALA:HB2	1.88	0.56
1:B:227:TYR:HA	1:B:230:LEU:HD12	1.87	0.56
1:B:169:GLU:OE2	1:B:197:ARG:NH1	2.39	0.55
1:A:129:ILE:HD11	1:A:171:LEU:CD1	2.36	0.54
1:A:209:LEU:HD21	1:A:246:LEU:CD2	2.36	0.54
1:B:135:GLN:HG2	1:B:136:HIS:CD2	2.43	0.54
1:A:133:ASN:HD22	1:A:161:TYR:HA	1.73	0.53
1:A:177:MET:HE1	1:A:199:TRP:CE3	2.44	0.53
1:A:184:ASP:O	1:A:185:ARG:HB2	2.09	0.52
1:A:133:ASN:HD22	1:A:162:GLY:N	2.08	0.51
1:B:166:SER:O	1:B:170:VAL:HG23	2.11	0.51
1:B:106:ILE:HG21	1:B:109:VAL:CG2	2.41	0.51
1:B:235:GLN:HG2	1:B:259:LEU:O	2.11	0.51
1:B:70:ARG:HH21	1:B:151:ARG:HD3	1.76	0.50
1:B:169:GLU:OE1	1:B:197:ARG:NH1	2.37	0.50
1:A:87:PHE:CE2	1:A:88:MET:HE1	2.47	0.50
1:B:194:THR:CG2	1:B:195:ALA:N	2.75	0.49
1:A:109:VAL:HG13	1:A:123:LEU:CD2	2.43	0.48
1:A:58:SER:HA	1:A:88:MET:HE3	1.95	0.48
1:B:194:THR:C	1:B:196:ARG:O	2.57	0.48
1:B:195:ALA:C	1:B:196:ARG:O	2.55	0.48
1:B:245:ASP:OD2	1:B:247:ARG:CB	2.62	0.48
1:A:133:ASN:HB3	1:A:161:TYR:CD2	2.49	0.47
1:B:181:LEU:HB3	1:B:186:PHE:HB2	1.95	0.47
1:B:60:LEU:HD21	1:B:121:ILE:HD12	1.96	0.47
1:A:194:THR:C	1:A:196:ARG:O	2.58	0.47
1:A:237:ASP:CG	1:B:222:ARG:HH12	2.23	0.47
1:B:155:GLN:NE2	1:B:157:LEU:HD21	2.29	0.47
1:A:249:ASP:O	1:A:250:SER:OG	2.27	0.47
1:B:194:THR:O	1:B:196:ARG:O	2.33	0.47
1:A:203:LEU:O	1:A:205:ASN:O	2.34	0.47
1:A:185:ARG:HG2	1:A:185:ARG:O	2.15	0.46
1:B:58:SER:HA	1:B:88:MET:HE3	1.98	0.46
1:A:256:TRP:CE2	1:B:250:SER:HA	2.51	0.46
1:A:211:LEU:O	1:A:219:ARG:NH2	2.46	0.45
1:B:91:ASP:O	1:B:95:ILE:HG23	2.16	0.45
1:B:163:PRO:HD2	1:B:193:MET:HE2	1.98	0.45
1:B:95:ILE:HD12	1:B:99:ILE:CD1	2.46	0.45
1:B:130:ALA:HB2	1:B:158:PRO:HG2	1.99	0.45
1:B:91:ASP:HB3	1:B:94:ILE:HD12	1.99	0.44
1:B:109:VAL:HG13	1:B:123:LEU:HD23	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:PRO:O	1:A:231:GLN:HB3	2.18	0.44
1:A:87:PHE:CZ	1:A:88:MET:CE	3.00	0.43
1:B:68:TYR:HE1	1:B:152:THR:HB	1.82	0.43
1:A:184:ASP:O	1:A:185:ARG:CB	2.65	0.43
1:B:163:PRO:CD	1:B:193:MET:HE2	2.48	0.43
1:B:209:LEU:HD22	1:B:211:LEU:HD13	1.99	0.43
1:A:210:ASN:OD1	1:A:247:ARG:HD3	2.17	0.43
1:B:227:TYR:O	1:B:228:PRO:C	2.61	0.43
1:B:132:TRP:O	1:B:136:HIS:CA	2.67	0.43
1:B:194:THR:CG2	1:B:195:ALA:H	2.28	0.43
1:B:170:VAL:HA	1:B:199:TRP:HH2	1.84	0.43
1:A:102:ARG:O	1:A:103:LEU:HD23	2.18	0.42
1:A:87:PHE:CE2	1:A:88:MET:CE	3.03	0.42
1:B:222:ARG:O	1:B:226:LEU:HB2	2.20	0.42
1:B:111:VAL:HG13	1:B:121:ILE:HG12	2.01	0.41
1:B:148:PRO:HA	1:B:149:PRO:HD3	1.86	0.41
1:B:81:LEU:HD11	1:B:98:GLN:HG3	2.01	0.41
1:B:95:ILE:HG13	1:B:96:GLN:N	2.35	0.41
1:B:133:ASN:HD22	1:B:161:TYR:HA	1.85	0.41
1:B:74:ILE:O	1:B:75:ARG:C	2.63	0.41
1:B:203:LEU:HB3	1:B:204:ASN:H	1.58	0.41
1:A:133:ASN:ND2	1:A:162:GLY:N	2.59	0.41
1:B:148:PRO:HD2	1:B:151:ARG:HG3	2.02	0.41
1:A:102:ARG:C	1:A:103:LEU:HD23	2.46	0.41
1:B:131:ARG:HG2	1:B:137:MET:HE3	2.02	0.41
1:B:235:GLN:HG2	1:B:259:LEU:C	2.46	0.41
1:A:150:GLU:H	1:A:150:GLU:CD	2.29	0.40
1:B:81:LEU:HD21	1:B:94:ILE:HG22	2.02	0.40
1:A:249:ASP:O	1:A:250:SER:CB	2.69	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	201/220 (91%)	187 (93%)	12 (6%)	2 (1%)	12	32
1	B	192/220 (87%)	170 (88%)	21 (11%)	1 (0%)	24	48
All	All	393/440 (89%)	357 (91%)	33 (8%)	3 (1%)	16	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	185	ARG
1	A	133	ASN
1	B	236	THR

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	177/192 (92%)	153 (86%)	24 (14%)	3	9
1	B	172/192 (90%)	142 (83%)	30 (17%)	2	5
All	All	349/384 (91%)	295 (84%)	54 (16%)	2	7

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	LEU
1	A	92	VAL
1	A	109	VAL
1	A	110	SER
1	A	119	LEU
1	A	120	LYS
1	A	129	ILE
1	A	144	THR
1	A	147	VAL
1	A	150	GLU
1	A	159	MET
1	A	160	LEU
1	A	168	ASN

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Mol	Chain	Res	Type
1	A	172	GLN
1	A	196	ARG
1	A	208	LYS
1	A	209	LEU
1	A	215	ASP
1	A	217	MET
1	A	226	LEU
1	A	229	VAL
1	A	240	ARG
1	A	246	LEU
1	A	259	LEU
1	B	60	LEU
1	B	75	ARG
1	B	81	LEU
1	B	88	MET
1	B	92	VAL
1	B	95	ILE
1	B	106	ILE
1	B	109	VAL
1	B	119	LEU
1	B	135	GLN
1	B	144	THR
1	B	152	THR
1	B	159	MET
1	B	160	LEU
1	B	176	GLU
1	B	188	LEU
1	B	196	ARG
1	B	204	ASN
1	B	206	ASP
1	B	209	LEU
1	B	211	LEU
1	B	213	ARG
1	B	224	VAL
1	B	231	GLN
1	B	233	GLN
1	B	235	GLN
1	B	236	THR
1	B	237	ASP
1	B	250	SER
1	B	259	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	HIS
1	A	90	GLN
1	A	108	GLN
1	A	122	HIS
1	A	133	ASN
1	A	200	GLN
1	B	71	ASN
1	B	76	GLN
1	B	98	GLN
1	B	122	HIS
1	B	133	ASN
1	B	135	GLN
1	B	136	HIS
1	B	204	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 oligosaccharides 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	203/220 (92%)	1.03	24 (11%) 9 7	49, 58, 63, 75	0
1	B	198/220 (90%)	1.36	50 (25%) 1 1	49, 58, 64, 74	0
All	All	401/440 (91%)	1.19	74 (18%) 3 3	49, 58, 64, 75	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	259	LEU	6.1
1	A	204	ASN	5.9
1	B	82	GLY	5.8
1	A	184	ASP	5.8
1	B	184	ASP	4.9
1	A	246	LEU	4.8
1	A	205	ASN	4.5
1	B	205	ASN	4.3
1	B	132	TRP	4.1
1	B	235	GLN	4.0
1	A	249	ASP	4.0
1	B	246	LEU	3.9
1	B	68	TYR	3.9
1	B	204	ASN	3.9
1	B	153	SER	3.8
1	B	214	GLY	3.8
1	B	213	ARG	3.8
1	B	217	MET	3.7
1	B	206	ASP	3.6
1	B	152	THR	3.6
1	B	258	PRO	3.6
1	A	206	ASP	3.5
1	B	133	ASN	3.5
1	B	65	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	155	GLN	3.5
1	B	185	ARG	3.3
1	B	193	MET	3.3
1	A	193	MET	3.2
1	B	194	THR	3.2
1	B	218	LYS	3.2
1	B	207	ILE	3.2
1	B	216	THR	3.1
1	B	177	MET	3.1
1	B	167	ALA	3.0
1	A	164	GLU	3.0
1	B	154	LYS	3.0
1	B	234	ALA	3.0
1	B	67	HIS	2.9
1	B	129	ILE	2.9
1	A	132	TRP	2.9
1	A	236	THR	2.8
1	B	209	LEU	2.8
1	A	252	ALA	2.8
1	B	85	GLY	2.7
1	B	135	GLN	2.7
1	B	148	PRO	2.7
1	B	211	LEU	2.7
1	B	81	LEU	2.6
1	A	215	ASP	2.6
1	A	195	ALA	2.6
1	A	213	ARG	2.5
1	B	146	SER	2.5
1	B	198	SER	2.4
1	A	207	ILE	2.4
1	A	248	TYR	2.4
1	A	86	THR	2.3
1	A	134	ASP	2.3
1	A	199	TRP	2.3
1	B	241	ILE	2.2
1	A	114	GLN	2.2
1	B	199	TRP	2.2
1	A	163	PRO	2.2
1	B	149	PRO	2.2
1	A	183	LYS	2.2
1	A	133	ASN	2.2
1	B	232	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	195	ALA	2.1
1	B	165	GLY	2.1
1	B	248	TYR	2.1
1	B	215	ASP	2.1
1	B	169	GLU	2.1
1	A	197	ARG	2.0
1	B	196	ARG	2.0
1	B	212	GLY	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 oligosaccharides 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.