



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:14 PM UTC

PDB ID : 3FK4 / pdb_00003fk4
Title : Crystal structure of RuBisCO-like protein from Bacillus Cereus ATCC 14579
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Deposited on : 2008-12-15
Resolution : 2.00 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

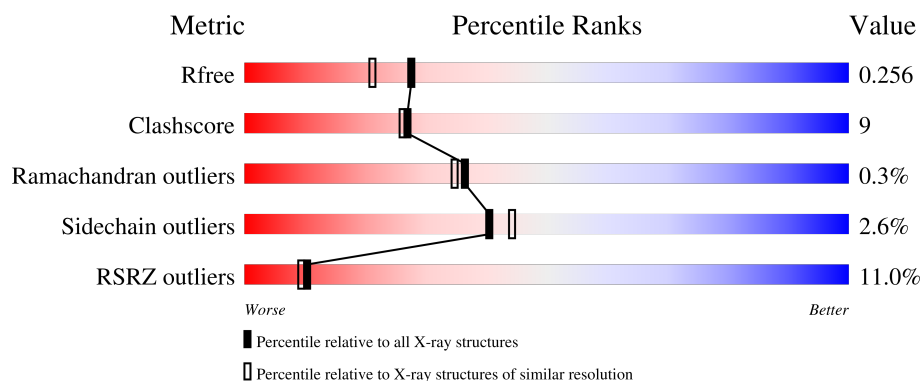
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	414	8% 76% 17% • 6%
1	B	414	13% 72% 20% • 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6206 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 RuBisCO-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	388	Total	C	N	O	S	0	0	0
			3009	1937	511	558	3			
1	B	391	Total	C	N	O	S	0	0	0
			3037	1959	516	559	3			

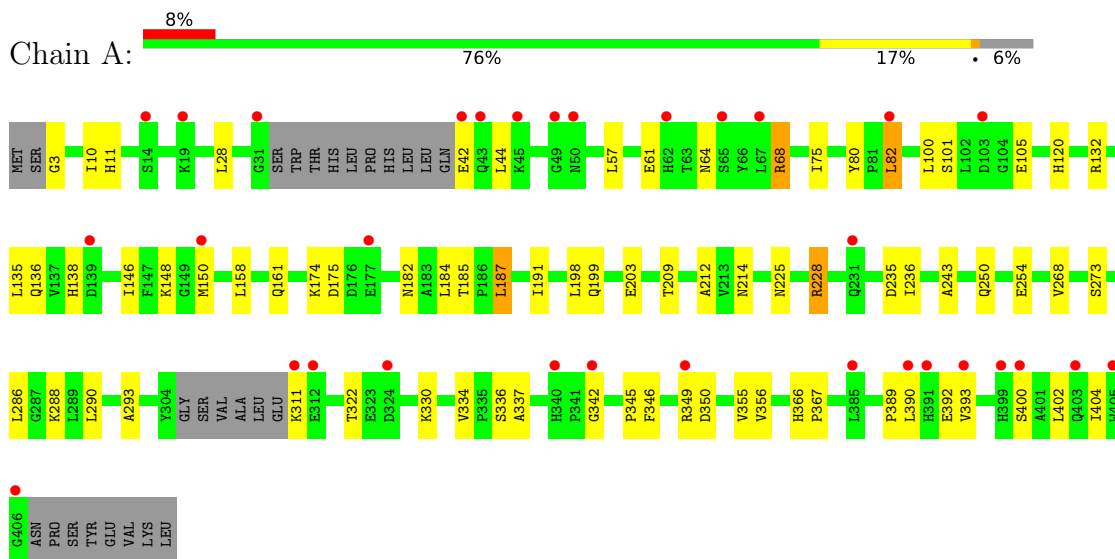
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	111	Total	O	0	0
			111	111		
2	B	49	Total	O	0	0
			49	49		

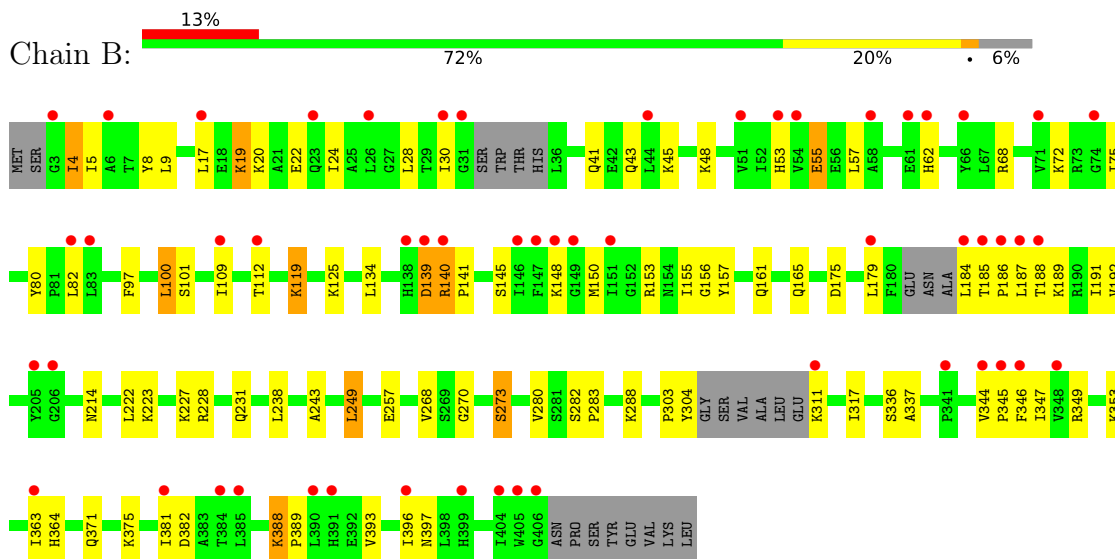
3 Residue-property plots

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: RuBisCO-like protein



• Molecule 1: RuBisCO-like protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.92Å 115.67Å 116.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.67 – 2.00 24.67 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (24.67-2.00) 99.6 (24.67-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.235 , 0.257 0.235 , 0.256	Depositor DCC
R_{free} test set	3572 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.012 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6206	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.05% 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	0.40	0/3072	0.88	6/4154 (0.1%)
1	B	0.35	0/3101	0.87	5/4193 (0.1%)
All	All	0.38	0/6173	0.88	11/8347 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	138	HIS	N-CA-C	8.53	123.80	113.23
1	A	273	SER	N-CA-C	7.18	122.22	113.38
1	B	273	SER	N-CA-C	6.81	122.43	113.30
1	B	243	ALA	N-CA-C	-6.38	105.31	113.23
1	A	243	ALA	N-CA-C	-5.78	106.23	113.28
1	A	288	LYS	N-CA-C	5.55	117.02	110.97
1	B	139	ASP	N-CA-C	5.52	119.56	111.34
1	B	222	LEU	N-CA-C	5.45	116.90	111.07
1	A	10	ILE	N-CA-C	5.44	116.32	108.48
1	B	288	LYS	N-CA-C	5.31	116.75	110.97
1	A	228	ARG	N-CA-C	-5.08	105.83	111.36

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	3009	0	3048	48	0
1	B	3037	0	3085	66	0
2	A	111	0	0	2	0
2	B	49	0	0	1	0
All	All	6206	0	6133	112	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 (112) 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:9:LEU:HB2	1:B:109:ILE:HD11	1.29	1.09
1:B:148:LYS:H	1:B:161:GLN:HE22	1.04	0.93
1:A:148:LYS:H	1:A:161:GLN:HE22	1.31	0.78
1:A:349:ARG:HG3	1:A:350:ASP:N	1.99	0.78
1:B:19:LYS:HE3	1:B:19:LYS:HA	1.66	0.75
1:B:283:PRO:HA	1:B:317:ILE:HG12	1.72	0.72
1:B:381:ILE:HD12	1:B:382:ASP:N	2.07	0.70
1:B:303:PRO:HB2	1:B:311:LYS:HG2	1.75	0.69
1:B:150:MET:HE3	1:B:153:ARG:HB2	1.75	0.68
1:A:182:ASN:HD21	1:A:185:THR:H	1.42	0.67
1:A:61:GLU:HA	1:A:64:ASN:HD22	1.61	0.66
1:B:184:LEU:O	1:B:184:LEU:HD23	1.96	0.65
1:B:148:LYS:H	1:B:161:GLN:NE2	1.88	0.64
1:A:182:ASN:ND2	1:A:185:THR:H	1.97	0.63
1:A:28:LEU:HD11	1:A:100:LEU:HD13	1.80	0.62
1:B:57:LEU:HD11	1:B:75:ILE:HG12	1.82	0.61
1:B:140:ARG:HG3	1:B:381:ILE:HD13	1.84	0.59
1:A:198:LEU:HD22	1:A:209:THR:HB	1.85	0.58
1:A:57:LEU:HD11	1:A:75:ILE:HG13	1.86	0.57
1:A:235:ASP:O	1:A:236:ILE:HD13	2.05	0.57
1:B:139:ASP:CG	1:B:140:ARG:H	2.12	0.57
1:B:388:LYS:HB2	1:B:388:LYS:NZ	2.19	0.57
1:B:282:SER:N	1:B:283:PRO:HD2	2.20	0.57
1:B:5:ILE:HB	1:B:112:THR:OG1	2.05	0.56
1:B:140:ARG:HG2	1:B:141:PRO:O	2.07	0.55
1:B:150:MET:HE1	1:B:157:TYR:HD2	1.72	0.55
1:B:388:LYS:HB2	1:B:388:LYS:HZ2	1.71	0.55
1:B:119:LYS:HB2	1:B:119:LYS:NZ	2.22	0.54
1:A:187:LEU:HD22	1:A:191:ILE:CD1	2.38	0.53
1:B:148:LYS:N	1:B:161:GLN:HE22	1.88	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:400:SER:O	1:A:404:ILE:HG13	2.09	0.52
1:A:349:ARG:HG3	1:A:350:ASP:H	1.74	0.52
1:B:68:ARG:HG3	1:B:68:ARG:HH11	1.75	0.52
1:A:389:PRO:HB2	1:A:392:GLU:HG3	1.93	0.51
1:B:336:SER:O	1:B:337:ALA:HB3	2.11	0.51
1:A:28:LEU:HD11	1:A:100:LEU:CD1	2.41	0.51
1:B:4:ILE:HD13	1:B:5:ILE:N	2.26	0.51
1:B:227:LYS:HE2	1:B:257:GLU:HB2	1.92	0.51
1:A:11:HIS:HB2	1:A:105:GLU:HB3	1.93	0.51
1:A:199:GLN:HE21	1:A:203:GLU:HG3	1.76	0.51
1:B:17:LEU:HG	1:B:72:LYS:HD2	1.92	0.50
1:B:249:LEU:HD22	1:B:249:LEU:O	2.12	0.50
1:A:212:ALA:HA	1:A:236:ILE:HB	1.94	0.50
1:B:139:ASP:CG	1:B:140:ARG:N	2.70	0.50
1:A:120:HIS:HD2	2:A:477:HOH:O	1.95	0.49
1:A:182:ASN:HD22	1:A:184:LEU:H	1.60	0.49
1:A:68:ARG:HD2	1:B:62:HIS:ND1	2.28	0.49
1:B:186:PRO:O	1:B:189:LYS:HB3	2.13	0.49
1:B:303:PRO:HB3	1:B:311:LYS:HA	1.95	0.48
1:A:334:VAL:HG22	1:A:356:VAL:HB	1.96	0.48
1:B:363:ILE:HG13	1:B:364:HIS:N	2.29	0.48
1:A:346:PHE:HA	1:A:349:ARG:HG2	1.96	0.47
1:B:53:HIS:NE2	1:B:55:GLU:HB2	2.29	0.47
1:A:336:SER:O	1:A:337:ALA:HB3	2.14	0.47
1:B:30:ILE:O	1:B:30:ILE:HD12	2.14	0.47
1:A:355:VAL:HG22	1:A:356:VAL:N	2.29	0.47
1:B:30:ILE:HG13	1:B:48:LYS:HA	1.96	0.47
1:B:396:ILE:HG13	1:B:397:ASN:N	2.29	0.47
1:B:214:ASN:HA	1:B:238:LEU:HB3	1.97	0.46
1:B:184:LEU:HD22	1:B:185:THR:HG23	1.98	0.46
1:B:43:GLN:NE2	2:B:453:HOH:O	2.48	0.46
1:B:311:LYS:O	1:B:311:LYS:HD3	2.16	0.46
1:B:41:GLN:O	1:B:45:LYS:HG2	2.16	0.45
1:B:344:VAL:HB	1:B:345:PRO:HD3	1.98	0.45
1:A:225:ASN:HD22	1:A:228:ARG:HH11	1.65	0.45
1:A:250:GLN:O	1:A:254:GLU:HG3	2.17	0.45
1:A:342:GLY:O	1:A:345:PRO:HD2	2.16	0.45
1:B:228:ARG:NH1	1:B:231:GLN:HE22	2.15	0.45
1:A:150:MET:HE1	1:A:175:ASP:OD1	2.16	0.45
1:B:304:TYR:CE2	1:B:347:ILE:HG12	2.52	0.45
1:B:186:PRO:O	1:B:189:LYS:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:304:TYR:HE2	1:B:347:ILE:HG12	1.83	0.44
1:B:20:LYS:O	1:B:24:ILE:HG13	2.17	0.44
1:A:28:LEU:HD12	1:A:80:TYR:OH	2.18	0.44
1:B:28:LEU:HD12	1:B:80:TYR:OH	2.17	0.44
1:A:367:PRO:HD3	1:A:404:ILE:HD12	1.99	0.44
1:B:145:SER:OG	1:B:165:GLN:NE2	2.51	0.44
1:A:290:LEU:HD11	2:A:498:HOH:O	2.18	0.44
1:A:366:HIS:CG	1:A:367:PRO:HD2	2.53	0.44
1:B:45:LYS:O	1:B:48:LYS:HG2	2.17	0.44
1:B:184:LEU:HD23	1:B:184:LEU:C	2.42	0.44
1:A:390:LEU:HB3	1:A:402:LEU:HD11	1.99	0.43
1:A:42:GLU:C	1:A:44:LEU:H	2.27	0.43
1:B:155:ILE:HG23	1:B:156:GLY:N	2.33	0.43
1:B:8:TYR:CE1	1:B:100:LEU:HD23	2.54	0.43
1:B:82:LEU:O	1:B:82:LEU:HD23	2.17	0.43
1:B:175:ASP:HB3	1:B:179:LEU:HD11	2.00	0.43
1:A:132:ARG:O	1:A:136:GLN:N	2.52	0.42
1:A:225:ASN:ND2	1:A:228:ARG:HH11	2.17	0.42
1:B:165:GLN:HG2	1:B:363:ILE:CD1	2.50	0.42
1:B:97:PHE:O	1:B:101:SER:HB3	2.19	0.42
1:B:388:LYS:HE3	1:B:393:VAL:HG22	2.01	0.42
1:A:146:ILE:HG12	1:A:174:LYS:HG2	2.02	0.42
1:A:389:PRO:O	1:A:393:VAL:HG23	2.20	0.42
1:B:371:GLN:O	1:B:375:LYS:HG3	2.20	0.42
1:B:22:GLU:HA	1:B:22:GLU:OE1	2.20	0.42
1:A:158:LEU:HD23	1:A:158:LEU:C	2.44	0.41
1:B:228:ARG:NH1	1:B:231:GLN:NE2	2.68	0.41
1:B:388:LYS:HA	1:B:389:PRO:HD3	1.95	0.41
1:B:187:LEU:O	1:B:191:ILE:HG13	2.20	0.41
1:A:187:LEU:HD22	1:A:191:ILE:HD11	2.01	0.41
1:A:286:LEU:HA	1:A:290:LEU:HD12	2.02	0.41
1:A:322:THR:O	1:A:330:LYS:HE2	2.20	0.41
1:B:188:THR:O	1:B:192:VAL:HG23	2.20	0.41
1:B:273:SER:HB3	1:B:280:VAL:O	2.21	0.41
1:A:250:GLN:HA	1:A:293:ALA:O	2.21	0.41
1:A:346:PHE:O	1:A:349:ARG:HG2	2.21	0.41
1:A:290:LEU:HD23	1:A:290:LEU:HA	1.95	0.40
1:A:3:GLY:HA2	1:A:82:LEU:HD23	2.02	0.40
1:A:101:SER:O	1:B:270:GLY:HA3	2.21	0.40
1:A:135:LEU:O	1:A:136:GLN:HB2	2.22	0.40
1:B:346:PHE:O	1:B:349:ARG:HB3	2.21	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	382/414 (92%)	366 (96%)	15 (4%)	1 (0%)	36	35
1	B	383/414 (92%)	367 (96%)	15 (4%)	1 (0%)	36	35
All	All	765/828 (92%)	733 (96%)	30 (4%)	2 (0%)	36	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	268	VAL
1	B	268	VAL

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	322/346 (93%)	317 (98%)	5 (2%)	55	62
1	B	326/346 (94%)	314 (96%)	12 (4%)	30	30
All	All	648/692 (94%)	631 (97%)	17 (3%)	40	44

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	68	ARG

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Mol	Chain	Res	Type
1	A	82	LEU
1	A	187	LEU
1	A	214	ASN
1	A	311	LYS
1	B	4	ILE
1	B	19	LYS
1	B	55	GLU
1	B	100	LEU
1	B	119	LYS
1	B	125	LYS
1	B	134	LEU
1	B	140	ARG
1	B	223	LYS
1	B	249	LEU
1	B	353	LYS
1	B	388	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	47	HIS
1	A	62	HIS
1	A	120	HIS
1	A	133	ASN
1	A	161	GLN
1	A	182	ASN
1	A	199	GLN
1	A	214	ASN
1	A	225	ASN
1	A	391	HIS
1	A	397	ASN
1	B	23	GLN
1	B	38	HIS
1	B	41	GLN
1	B	46	GLN
1	B	50	ASN
1	B	133	ASN
1	B	138	HIS
1	B	161	GLN
1	B	165	GLN
1	B	214	ASN

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Mol	Chain	Res	Type
1	B	231	GLN
1	B	386	GLN
1	B	397	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	388/414 (93%)	0.43	32 (8%) 17 16	23, 37, 68, 77	0
1	B	391/414 (94%)	0.93	54 (13%) 6 6	26, 49, 71, 80	0
All	All	779/828 (94%)	0.68	86 (11%) 10 9	23, 44, 70, 80	0

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	184	LEU	6.4
1	B	406	GLY	5.1
1	B	151	ILE	4.8
1	B	31	GLY	4.7
1	A	42	GLU	4.6
1	B	139	ASP	4.0
1	A	177	GLU	3.9
1	A	390	LEU	3.8
1	B	185	THR	3.8
1	A	406	GLY	3.7
1	B	17	LEU	3.5
1	B	51	VAL	3.5
1	B	109	ILE	3.3
1	B	405	TRP	3.3
1	B	205	TYR	3.2
1	B	381	ILE	3.1
1	B	390	LEU	3.1
1	B	138	HIS	3.0
1	B	149	GLY	3.0
1	A	349	ARG	2.9
1	B	54	VAL	2.9
1	A	405	TRP	2.9
1	B	62	HIS	2.9
1	B	30	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	112	THR	2.8
1	B	82	LEU	2.8
1	A	311	LYS	2.7
1	B	391	HIS	2.7
1	B	3	GLY	2.7
1	A	231	GLN	2.6
1	B	346	PHE	2.6
1	B	363	ILE	2.6
1	A	67	LEU	2.6
1	B	188	THR	2.6
1	A	14	SER	2.6
1	B	348	VAL	2.5
1	B	66	TYR	2.5
1	A	150	MET	2.5
1	B	385	LEU	2.5
1	A	45	LYS	2.5
1	A	31	GLY	2.5
1	B	71	VAL	2.4
1	B	23	GLN	2.4
1	B	341	PRO	2.4
1	B	187	LEU	2.4
1	B	396	ILE	2.4
1	A	65	SER	2.4
1	B	140	ARG	2.4
1	B	186	PRO	2.4
1	B	58	ALA	2.3
1	B	26	LEU	2.3
1	B	206	GLY	2.3
1	A	342	GLY	2.3
1	A	43	GLN	2.3
1	A	50	ASN	2.3
1	A	139	ASP	2.3
1	B	61	GLU	2.3
1	B	74	GLY	2.3
1	B	399	HIS	2.2
1	B	147	PHE	2.2
1	B	345	PRO	2.2
1	B	311	LYS	2.2
1	B	179	LEU	2.2
1	A	19	LYS	2.2
1	B	44	LEU	2.2
1	A	340	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	404	ILE	2.2
1	A	393	VAL	2.2
1	B	6	ALA	2.2
1	A	399	HIS	2.2
1	B	344	VAL	2.1
1	B	53	HIS	2.1
1	A	82	LEU	2.1
1	A	324	ASP	2.1
1	B	146	ILE	2.1
1	A	391	HIS	2.1
1	B	148	LYS	2.1
1	A	400	SER	2.1
1	A	403	GLN	2.1
1	A	62	HIS	2.1
1	A	385	LEU	2.1
1	A	312	GLU	2.0
1	B	384	THR	2.0
1	A	49	GLY	2.0
1	B	83	LEU	2.0
1	A	103	ASP	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.