



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 04:03 PM BST

PDB ID : 6KRR  
Title : Peroxiredoxin from *Aeropyrum pernix* K1 (ApPrx) 0Cys W210A mutant  
Authors : Himiyama, T.; Nakamura, T.  
Deposited on : 2019-08-22  
Resolution : 2.15 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](mailto: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.

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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	:	2.13.1
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)	:	2.13.1

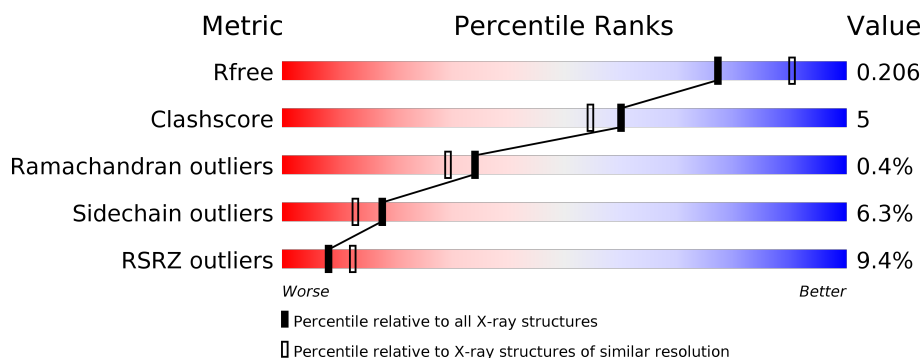
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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  $\geq 3$ , 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	244	<div> <div>9%</div> <div>87%</div> <div>12%</div> </div>
1	B	244	<div> <div>9%</div> <div>85%</div> <div>13%</div> </div>
1	C	244	<div> <div>9%</div> <div>82%</div> <div>15%</div> </div>
1	D	244	<div> <div>9%</div> <div>87%</div> <div>11%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8186 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 Peroxiredoxin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	244	Total	C	N	O	S	0	0	0
			1964	1260	346	354	4			
1	B	244	Total	C	N	O	S	0	0	0
			1964	1260	346	354	4			
1	C	244	Total	C	N	O	S	0	0	0
			1964	1260	346	354	4			
1	D	244	Total	C	N	O	S	0	0	0
			1964	1260	346	354	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	50	SER	CYS	engineered mutation	UNP Q9Y9L0
A	207	SER	CYS	engineered mutation	UNP Q9Y9L0
A	210	ALA	TRP	engineered mutation	UNP Q9Y9L0
A	213	SER	CYS	engineered mutation	UNP Q9Y9L0
B	50	SER	CYS	engineered mutation	UNP Q9Y9L0
B	207	SER	CYS	engineered mutation	UNP Q9Y9L0
B	210	ALA	TRP	engineered mutation	UNP Q9Y9L0
B	213	SER	CYS	engineered mutation	UNP Q9Y9L0
C	50	SER	CYS	engineered mutation	UNP Q9Y9L0
C	207	SER	CYS	engineered mutation	UNP Q9Y9L0
C	210	ALA	TRP	engineered mutation	UNP Q9Y9L0
C	213	SER	CYS	engineered mutation	UNP Q9Y9L0
D	50	SER	CYS	engineered mutation	UNP Q9Y9L0
D	207	SER	CYS	engineered mutation	UNP Q9Y9L0
D	210	ALA	TRP	engineered mutation	UNP Q9Y9L0
D	213	SER	CYS	engineered mutation	UNP Q9Y9L0

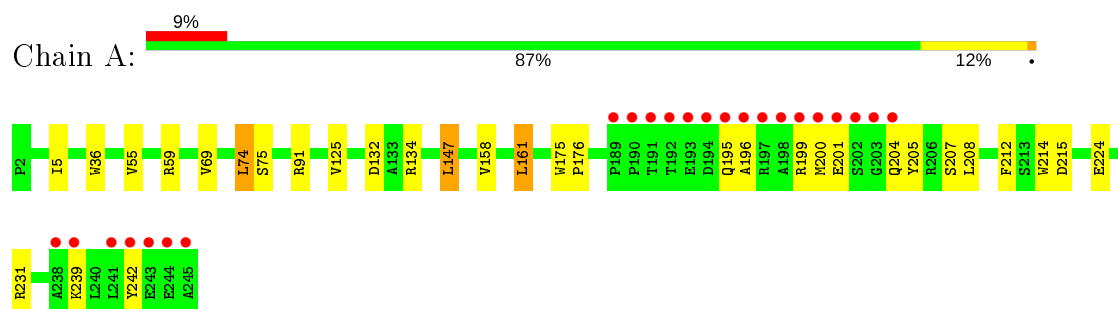
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	90	Total 90	O 90	0	0
2	B	87	Total 87	O 87	0	0
2	C	80	Total 80	O 80	0	0
2	D	73	Total 73	O 73	0	0

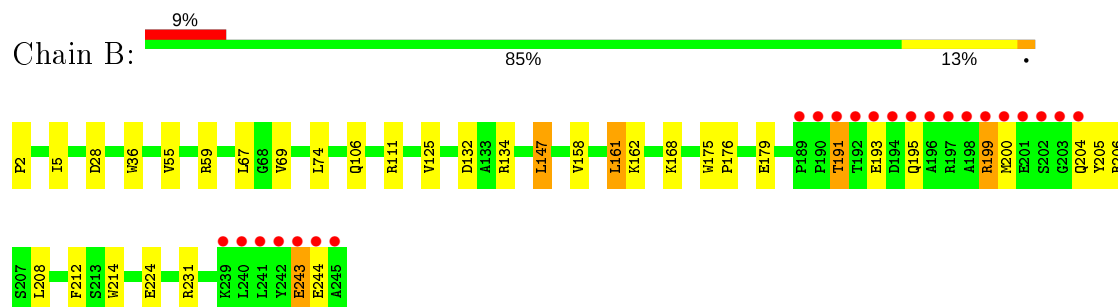
### 3 Residue-property plots [i](#)

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.

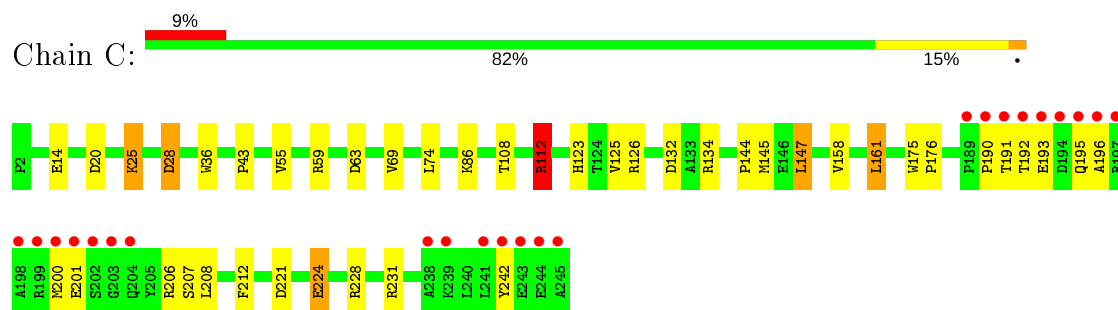
#### • Molecule 1: Peroxiredoxin



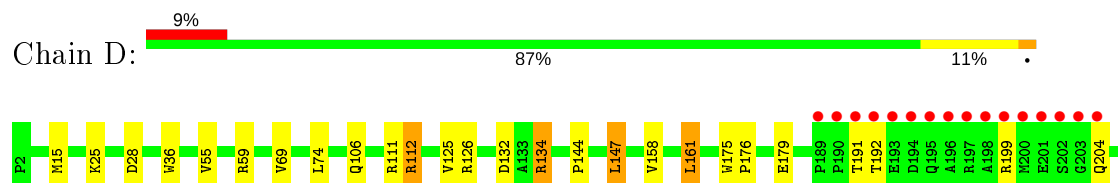
#### • Molecule 1: Peroxiredoxin

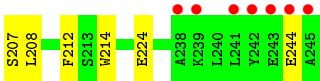


#### • Molecule 1: Peroxiredoxin



#### • Molecule 1: Peroxiredoxin





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	98.50Å 98.71Å 194.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.40 – 2.15 49.36 – 2.15	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.40-2.15) 100.0 (49.36-2.15)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.77 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.8	Depositor
R, $R_{free}$	0.169 , 0.199 0.180 , 0.206	Depositor DCC
$R_{free}$ test set	5106 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	33.0	Xtriage
Anisotropy	0.060	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 51.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.008 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8186	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.20% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.79	0/2017	0.96	1/2741 (0.0%)
1	B	0.79	0/2017	0.97	1/2741 (0.0%)
1	C	0.79	0/2017	0.94	4/2741 (0.1%)
1	D	0.78	0/2017	0.97	2/2741 (0.1%)
All	All	0.79	0/8068	0.96	8/10964 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	126	ARG	NE-CZ-NH1	-6.89	116.86	120.30
1	C	112	ARG	NE-CZ-NH1	6.71	123.65	120.30
1	B	134	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	C	126	ARG	NE-CZ-NH2	5.48	123.04	120.30
1	A	134	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	C	134	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	D	134	ARG	NE-CZ-NH1	5.17	122.88	120.30
1	C	231	ARG	CG-CD-NE	-5.09	101.10	111.80

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	1964	0	1954	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1964	0	1954	22	0
1	C	1964	0	1954	25	0
1	D	1964	0	1954	18	0
2	A	90	0	0	3	0
2	B	87	0	0	1	1
2	C	80	0	0	2	0
2	D	73	0	0	1	1
All	All	8186	0	7816	72	2

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 5.

All (72) 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:69:VAL:HG21	1:B:158:VAL:HG11	1.47	0.94
1:C:14:GLU:OE1	1:C:25:LYS:NZ	2.05	0.90
1:B:191:THR:H	1:B:195:GLN:HE21	1.21	0.88
1:C:69:VAL:HG21	1:C:158:VAL:HG11	1.57	0.86
1:C:112:ARG:HG2	1:C:112:ARG:HH11	1.43	0.84
1:D:69:VAL:HG21	1:D:158:VAL:HG11	1.60	0.81
1:A:69:VAL:HG21	1:A:158:VAL:HG11	1.67	0.76
1:A:161:LEU:HD13	1:B:147:LEU:HG	1.76	0.67
1:D:208:LEU:HD13	1:D:214:TRP:HZ3	1.60	0.67
1:B:106:GLN:O	1:B:111:ARG:NH2	2.28	0.67
1:C:191:THR:HG22	1:C:192:THR:HG23	1.77	0.66
1:A:147:LEU:HG	1:B:161:LEU:HD13	1.77	0.65
1:C:242:TYR:CD1	1:D:214:TRP:HH2	2.19	0.60
1:D:191:THR:HG23	1:D:192:THR:HG23	1.84	0.59
1:B:199:ARG:HG3	1:B:205:TYR:CE2	2.38	0.58
1:D:208:LEU:HD13	1:D:214:TRP:CZ3	2.39	0.57
1:C:112:ARG:CG	1:C:112:ARG:HH11	2.16	0.56
1:C:161:LEU:HD13	1:D:147:LEU:HG	1.88	0.55
1:D:132:ASP:OD2	1:D:134:ARG:NH1	2.37	0.55
1:C:147:LEU:HG	1:D:161:LEU:HD13	1.89	0.54
1:A:69:VAL:CG2	1:A:158:VAL:HG11	2.38	0.53
1:B:69:VAL:CG2	1:B:158:VAL:HG11	2.31	0.53
1:C:55:VAL:O	1:C:59:ARG:HG2	2.08	0.52
1:C:25:LYS:HE3	1:C:28:ASP:OD2	2.10	0.52
1:D:36:TRP:CD2	1:D:132:ASP:HA	2.44	0.52
1:D:199:ARG:NH2	2:D:301:HOH:O	2.38	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:55:VAL:O	1:D:59:ARG:HG2	2.11	0.51
1:A:199:ARG:HB3	1:A:205:TYR:HE2	1.77	0.50
1:B:67:LEU:O	1:B:162:LYS:HE2	2.12	0.50
1:A:196:ALA:O	1:A:200:MET:HB3	2.11	0.50
1:B:55:VAL:O	1:B:59:ARG:HG2	2.12	0.50
1:A:231:ARG:NH1	2:A:304:HOH:O	2.45	0.49
1:A:91:ARG:NH2	2:A:303:HOH:O	2.43	0.49
1:B:162:LYS:HE2	2:B:381:HOH:O	2.12	0.49
1:B:36:TRP:CD2	1:B:132:ASP:HA	2.48	0.48
1:C:36:TRP:CD2	1:C:132:ASP:HA	2.49	0.48
1:A:36:TRP:CD2	1:A:132:ASP:HA	2.49	0.48
1:A:242:TYR:CD1	1:B:214:TRP:HH2	2.32	0.48
1:C:123:HIS:HE1	2:C:309:HOH:O	1.97	0.47
1:C:175:TRP:CG	1:C:176:PRO:HA	2.49	0.47
1:C:193:GLU:HA	1:C:196:ALA:HB3	1.97	0.47
2:A:318:HOH:O	1:B:231:ARG:HD3	2.15	0.47
1:C:63:ASP:OD2	1:C:228:ARG:NH2	2.48	0.47
1:D:15:MET:O	1:D:25:LYS:HD3	2.14	0.47
1:A:74:LEU:HD23	1:A:75:SER:N	2.31	0.46
1:C:59:ARG:HG3	1:D:179:GLU:OE2	2.14	0.46
1:B:199:ARG:HG2	1:B:199:ARG:O	2.15	0.46
1:C:112:ARG:HG2	1:C:112:ARG:NH1	2.18	0.46
1:A:175:TRP:CG	1:A:176:PRO:HA	2.51	0.45
1:C:224:GLU:HG3	2:C:301:HOH:O	2.15	0.45
1:A:5:ILE:HD13	1:B:5:ILE:HD13	1.98	0.45
1:B:191:THR:N	1:B:195:GLN:HE21	2.02	0.45
1:C:43:PRO:HG3	1:C:145:MET:SD	2.57	0.44
1:A:55:VAL:O	1:A:59:ARG:HG2	2.18	0.44
1:A:59:ARG:HG3	1:B:179:GLU:OE2	2.16	0.44
1:D:112:ARG:CG	1:D:112:ARG:HH11	2.30	0.44
1:B:191:THR:HB	1:B:195:GLN:HE22	1.81	0.44
1:A:208:LEU:HD13	1:A:214:TRP:HZ3	1.83	0.43
1:C:191:THR:H	1:C:195:GLN:NE2	2.16	0.43
1:C:63:ASP:CG	1:C:228:ARG:HH22	2.22	0.43
1:B:191:THR:HB	1:B:195:GLN:NE2	2.34	0.42
1:B:175:TRP:CG	1:B:176:PRO:HA	2.55	0.42
1:B:191:THR:H	1:B:195:GLN:NE2	2.02	0.42
1:C:190:PRO:CB	1:C:196:ALA:HB2	2.50	0.42
1:C:20:ASP:OD2	1:C:86:LYS:NZ	2.40	0.41
1:D:106:GLN:O	1:D:111:ARG:NH2	2.53	0.41
1:C:69:VAL:CG2	1:C:158:VAL:HG11	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:ARG:CG	1:D:112:ARG:NH1	2.82	0.41
1:A:215:ASP:HB2	1:C:221:ASP:OD2	2.21	0.41
1:D:59:ARG:HA	1:D:59:ARG:HD2	1.93	0.40
1:B:243:GLU:H	1:B:243:GLU:HG2	1.74	0.40
1:D:175:TRP:CG	1:D:176:PRO:HA	2.56	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:379:HOH:O	2:B:379:HOH:O[2_555]	1.87	0.33
2:D:335:HOH:O	2:D:335:HOH:O[2_655]	2.13	0.07

## 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	242/244 (99%)	236 (98%)	5 (2%)	1 (0%)	34	29
1	B	242/244 (99%)	232 (96%)	9 (4%)	1 (0%)	34	29
1	C	242/244 (99%)	234 (97%)	7 (3%)	1 (0%)	34	29
1	D	242/244 (99%)	234 (97%)	7 (3%)	1 (0%)	34	29
All	All	968/976 (99%)	936 (97%)	28 (3%)	4 (0%)	34	29

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	125	VAL
1	D	125	VAL
1	B	125	VAL
1	A	125	VAL

### 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	209/209 (100%)	199 (95%)	10 (5%)	25	22
1	B	209/209 (100%)	192 (92%)	17 (8%)	11	7
1	C	209/209 (100%)	194 (93%)	15 (7%)	14	9
1	D	209/209 (100%)	198 (95%)	11 (5%)	22	19
All	All	836/836 (100%)	783 (94%)	53 (6%)	18	13

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	LEU
1	A	147	LEU
1	A	161	LEU
1	A	195	GLN
1	A	201	GLU
1	A	204	GLN
1	A	207	SER
1	A	212	PHE
1	A	224	GLU
1	A	239	LYS
1	B	2	PRO
1	B	28	ASP
1	B	74	LEU
1	B	147	LEU
1	B	161	LEU
1	B	168	LYS
1	B	191	THR
1	B	193	GLU
1	B	199	ARG
1	B	200	MET
1	B	204	GLN
1	B	206	ARG
1	B	208	LEU
1	B	212	PHE

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Mol	Chain	Res	Type
1	B	224	GLU
1	B	243	GLU
1	B	244	GLU
1	C	25	LYS
1	C	28	ASP
1	C	74	LEU
1	C	108	THR
1	C	112	ARG
1	C	144	PRO
1	C	147	LEU
1	C	161	LEU
1	C	200	MET
1	C	201	GLU
1	C	206	ARG
1	C	207	SER
1	C	208	LEU
1	C	212	PHE
1	C	224	GLU
1	D	28	ASP
1	D	74	LEU
1	D	112	ARG
1	D	144	PRO
1	D	147	LEU
1	D	161	LEU
1	D	204	GLN
1	D	207	SER
1	D	212	PHE
1	D	224	GLU
1	D	244	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	106	GLN
1	B	123	HIS
1	B	195	GLN
1	C	106	GLN
1	C	123	HIS
1	C	195	GLN
1	D	123	HIS
1	D	195	GLN
1	D	204	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 monosaccharides 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	244/244 (100%)	0.30	23 (9%) 8 12	28, 42, 108, 143	0
1	B	244/244 (100%)	0.26	23 (9%) 8 12	28, 41, 110, 152	0
1	C	244/244 (100%)	0.23	23 (9%) 8 12	28, 41, 117, 153	0
1	D	244/244 (100%)	0.29	23 (9%) 8 12	28, 43, 122, 146	0
All	All	976/976 (100%)	0.27	92 (9%) 8 12	28, 42, 119, 153	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	192	THR	10.6
1	D	245	ALA	10.4
1	A	192	THR	10.2
1	D	192	THR	9.2
1	B	197	ARG	8.9
1	C	200	MET	8.6
1	D	191	THR	8.5
1	B	194	ASP	8.3
1	A	245	ALA	8.3
1	A	197	ARG	8.1
1	C	201	GLU	8.0
1	A	200	MET	7.7
1	B	193	GLU	7.3
1	C	194	ASP	7.2
1	D	200	MET	7.1
1	A	242	TYR	7.0
1	D	193	GLU	7.0
1	C	193	GLU	7.0
1	D	244	GLU	6.9
1	B	242	TYR	6.9
1	A	195	GLN	6.9

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Mol	Chain	Res	Type	RSRZ
1	A	194	ASP	6.9
1	B	195	GLN	6.9
1	B	192	THR	6.8
1	B	198	ALA	6.7
1	C	245	ALA	6.6
1	B	200	MET	6.6
1	A	244	GLU	6.6
1	B	191	THR	6.5
1	A	201	GLU	6.3
1	C	191	THR	6.3
1	B	196	ALA	6.1
1	C	197	ARG	6.1
1	C	195	GLN	6.1
1	C	202	SER	6.1
1	D	196	ALA	6.1
1	D	197	ARG	5.9
1	D	194	ASP	5.9
1	C	196	ALA	5.8
1	A	193	GLU	5.6
1	A	191	THR	5.6
1	D	195	GLN	5.5
1	C	244	GLU	5.5
1	C	198	ALA	5.5
1	B	244	GLU	5.5
1	D	242	TYR	5.4
1	C	242	TYR	5.3
1	D	190	PRO	5.3
1	D	198	ALA	5.0
1	B	201	GLU	4.9
1	A	190	PRO	4.9
1	B	190	PRO	4.9
1	B	203	GLY	4.8
1	D	201	GLU	4.8
1	C	243	GLU	4.7
1	A	198	ALA	4.6
1	A	202	SER	4.6
1	C	190	PRO	4.5
1	C	203	GLY	4.4
1	A	196	ALA	4.1
1	B	239	LYS	4.0
1	B	245	ALA	4.0
1	B	199	ARG	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	202	SER	3.9
1	B	202	SER	3.8
1	A	199	ARG	3.7
1	D	239	LYS	3.6
1	C	189	PRO	3.6
1	D	243	GLU	3.6
1	B	241	LEU	3.5
1	B	243	GLU	3.4
1	B	189	PRO	3.4
1	A	243	GLU	3.4
1	D	204	GLN	3.3
1	D	189	PRO	3.2
1	C	238	ALA	3.1
1	C	199	ARG	3.1
1	B	204	GLN	3.1
1	A	204	GLN	2.8
1	A	203	GLY	2.7
1	C	204	GLN	2.7
1	A	238	ALA	2.6
1	D	199	ARG	2.5
1	D	203	GLY	2.5
1	A	241	LEU	2.4
1	A	239	LYS	2.4
1	B	240	LEU	2.3
1	C	241	LEU	2.3
1	C	239	LYS	2.2
1	D	241	LEU	2.2
1	A	189	PRO	2.1
1	D	238	ALA	2.1

## 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 monosaccharides in this entry.

## 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.