



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 06:21 PM GMT

PDB ID : 1V0B  
Title : CRYSTAL STRUCTURE OF THE T198A MUTANT OF PFPK5  
Authors : Holton, S.; Merckx, A.; Burgess, D.; Doerig, C.; Noble, M.; Endicott, J.  
Deposited on : 2004-03-26  
Resolution : 2.20 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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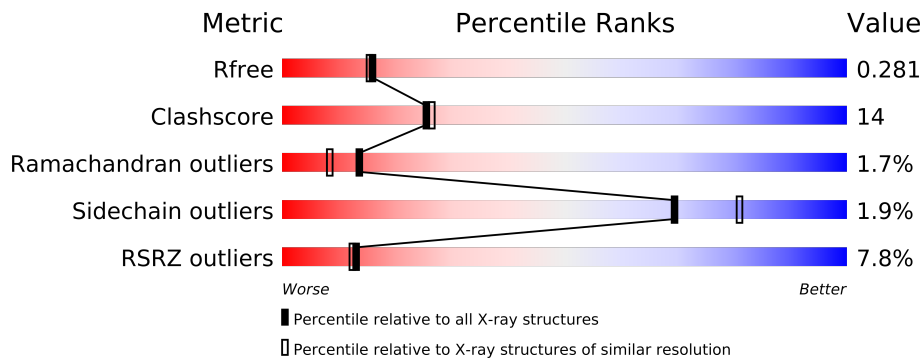
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

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))
$R_{free}$	66092	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	288	
1	B	288	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4728 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CONTROL PROTEIN 2 HOMOLOG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	0	1
			2310	1491	391	420	8			
1	B	262	Total	C	N	O	S	0	0	1
			2104	1363	356	378	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	198	ALA	THR	ENGINEERED MUTATION	UNP Q07785
B	198	ALA	THR	ENGINEERED MUTATION	UNP Q07785

- Molecule 2 is water.

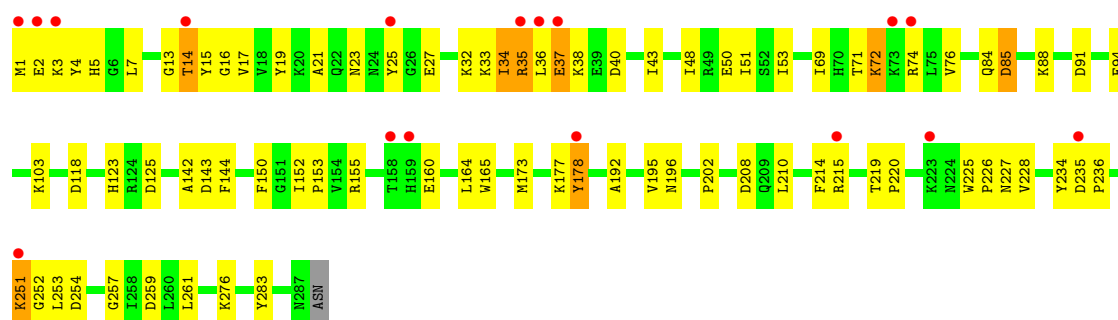
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	168	Total	O	0	0
			168	168		
2	B	146	Total	O	0	0
			146	146		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

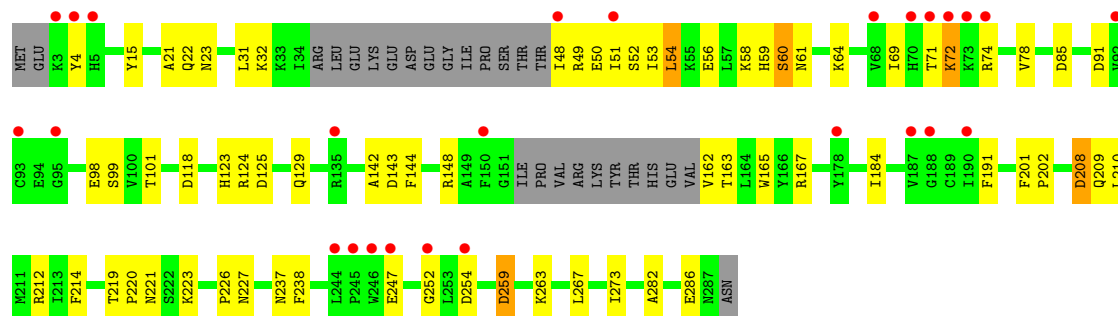
#### • Molecule 1: CELL DIVISION CONTROL PROTEIN 2 HOMOLOG

Chain A: 



#### • Molecule 1: CELL DIVISION CONTROL PROTEIN 2 HOMOLOG

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	55.29Å 94.65Å 133.23Å 90.00° 99.19° 90.00°	Depositor
Resolution (Å)	129.10 – 2.20 25.97 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.8 (129.10-2.20) 96.9 (25.97-2.20)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.45 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.223 , 0.275 0.226 , 0.281	Depositor DCC
$R_{free}$ test set	1679 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	28.8	Xtriage
Anisotropy	0.615	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 41.0	EDS
Estimated twinning fraction	0.019 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 33486 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4728	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.46% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 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	0.86	1/2361 (0.0%)	0.82	7/3191 (0.2%)
1	B	0.64	0/2149	0.77	7/2901 (0.2%)
All	All	0.76	1/4510 (0.0%)	0.80	14/6092 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	34	ILE	C-N	-8.55	1.14	1.34

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	208	ASP	CB-CG-OD2	8.12	125.61	118.30
1	B	259	ASP	CB-CG-OD2	6.50	124.15	118.30
1	A	35	ARG	CB-CA-C	6.21	122.83	110.40
1	A	208	ASP	CB-CG-OD2	6.21	123.89	118.30
1	A	91	ASP	CB-CG-OD2	6.11	123.80	118.30
1	A	40	ASP	CB-CG-OD2	5.97	123.67	118.30
1	B	85	ASP	CB-CG-OD2	5.56	123.31	118.30
1	B	91	ASP	CB-CG-OD2	5.55	123.30	118.30
1	B	118	ASP	CB-CG-OD2	5.53	123.27	118.30
1	A	85	ASP	CB-CG-OD2	5.40	123.16	118.30
1	A	259	ASP	CB-CG-OD2	5.23	123.01	118.30
1	B	254	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	143	ASP	CB-CG-OD2	5.17	122.95	118.30
1	B	143	ASP	CB-CG-OD2	5.01	122.81	118.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2310	0	2356	70	7
1	B	2104	0	2147	51	0
2	A	168	0	0	4	0
2	B	146	0	0	11	1
All	All	4728	0	4503	121	8

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (121) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:36:LEU:HD11	1:A:150:PHE:HB3	1.66	0.77
1:B:59:HIS:O	1:B:61:ASN:N	2.18	0.77
1:A:36:LEU:HG	1:A:37:GLU:H	1.49	0.76
1:A:2:GLU:HG3	1:A:69:ILE:CD1	2.16	0.75
1:B:50:GLU:O	1:B:54:LEU:HD13	1.89	0.72
1:A:14:THR:HB	1:A:155:ARG:HH11	1.58	0.68
1:A:50:GLU:O	1:A:53:ILE:HG22	1.94	0.68
1:A:51:ILE:HD11	1:A:150:PHE:CE2	2.28	0.68
1:B:48:ILE:N	2:B:2034:HOH:O	2.28	0.66
1:A:123:HIS:HD2	1:A:125:ASP:H	1.44	0.62
1:B:49:ARG:O	1:B:53:ILE:HG13	1.98	0.62
1:B:208:ASP:O	1:B:212:ARG:HG2	2.00	0.61
1:A:36:LEU:HG	1:A:37:GLU:N	2.17	0.59
1:B:165:TRP:CD1	1:B:202:PRO:HA	2.38	0.59
1:A:1:MET:HE1	1:A:76:VAL:HG21	1.85	0.59
1:A:2:GLU:HG3	1:A:69:ILE:HD11	1.85	0.59
1:B:237:ASN:O	1:B:238:PHE:C	2.40	0.59
1:A:36:LEU:CG	1:A:37:GLU:H	2.16	0.58
1:B:22:GLN:NE2	2:B:2017:HOH:O	2.36	0.58
1:A:1:MET:CE	1:A:76:VAL:HG21	2.34	0.58
1:B:23:ASN:ND2	2:B:2019:HOH:O	2.36	0.58
1:A:25:TYR:HB2	1:A:27:GLU:HG2	1.86	0.58
1:B:49:ARG:NH1	2:B:2037:HOH:O	2.37	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:219:THR:HG23	1:A:220:PRO:HD2	1.86	0.57
1:A:123:HIS:HE1	1:A:142:ALA:O	1.88	0.57
1:A:215:ARG:NH1	2:A:2123:HOH:O	2.38	0.57
1:A:173:MET:HG3	1:A:210:LEU:HD11	1.86	0.56
1:B:53:ILE:HG22	2:B:2038:HOH:O	2.06	0.56
1:A:36:LEU:HD22	1:A:150:PHE:CD2	2.41	0.56
1:B:129:GLN:HB2	2:B:2076:HOH:O	2.06	0.55
1:A:1:MET:O	1:A:3:LYS:N	2.39	0.55
1:B:31:LEU:CD2	1:B:78:VAL:HG13	2.37	0.55
1:A:254:ASP:OD1	1:A:257:GLY:HA3	2.07	0.55
1:B:282:ALA:O	1:B:286:GLU:HG3	2.07	0.55
1:A:13:GLY:HA2	2:A:2006:HOH:O	2.05	0.54
1:A:226:PRO:O	1:A:227:ASN:HB2	2.06	0.54
1:B:123:HIS:HE1	1:B:142:ALA:O	1.90	0.54
1:A:192:ALA:O	1:A:195:VAL:HG22	2.07	0.54
1:A:196:ASN:OD1	1:A:251:LYS:HG3	2.07	0.54
1:A:15:TYR:HB2	1:A:32:LYS:HE3	1.89	0.54
1:A:219:THR:CG2	1:A:220:PRO:HD2	2.37	0.54
1:A:173:MET:CG	1:A:210:LEU:HD11	2.37	0.54
1:A:71:THR:O	1:A:74:ARG:HB2	2.08	0.53
1:A:2:GLU:HG3	1:A:69:ILE:HG12	1.91	0.53
1:B:31:LEU:HD22	1:B:78:VAL:HG22	1.91	0.52
1:B:148:ARG:HH11	1:B:148:ARG:HG3	1.74	0.52
1:A:7:LEU:HD22	1:A:21:ALA:HB2	1.91	0.52
1:B:71:THR:O	1:B:74:ARG:HB2	2.10	0.52
1:B:226:PRO:O	1:B:227:ASN:HB2	2.09	0.52
1:B:123:HIS:HD2	1:B:125:ASP:H	1.57	0.51
1:A:2:GLU:HG3	1:A:69:ILE:CG1	2.41	0.50
1:A:254:ASP:OD1	1:A:257:GLY:N	2.44	0.50
1:B:162:VAL:N	2:B:2087:HOH:O	2.45	0.50
1:A:84:GLN:HG3	1:A:85:ASP:N	2.27	0.50
1:A:234:TYR:O	1:A:235:ASP:OD1	2.30	0.50
1:B:210:LEU:HD22	1:B:214:PHE:CZ	2.47	0.50
1:A:43:ILE:HG21	1:A:48:ILE:HD11	1.93	0.49
1:A:165:TRP:CD1	1:A:202:PRO:HA	2.47	0.49
1:A:2:GLU:CG	1:A:69:ILE:CD1	2.90	0.48
1:B:49:ARG:HA	1:B:52:SER:OG	2.13	0.48
1:A:160:GLU:O	1:A:164:LEU:HD13	2.13	0.48
1:A:2:GLU:N	2:A:2002:HOH:O	2.27	0.48
1:B:51:ILE:HD12	1:B:51:ILE:H	1.78	0.47
1:A:103:LYS:NZ	1:A:283:TYR:O	2.45	0.47
1:B:247:GLU:CD	1:B:247:GLU:H	2.17	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:59:HIS:O	1:B:60:SER:C	2.53	0.47
1:A:23:ASN:ND2	1:A:27:GLU:HG3	2.30	0.47
1:A:71:THR:O	1:A:72:LYS:C	2.53	0.46
1:A:210:LEU:HD22	1:A:214:PHE:CZ	2.51	0.46
1:B:51:ILE:O	1:B:54:LEU:HB2	2.16	0.46
1:B:167:ARG:NE	2:B:2090:HOH:O	2.46	0.46
1:B:221:ASN:OD1	1:B:223:LYS:N	2.43	0.45
1:B:69:ILE:HG22	1:B:71:THR:HG23	1.98	0.45
1:A:225:TRP:CG	1:A:228:VAL:HB	2.52	0.45
1:B:123:HIS:O	1:B:124:ARG:HB2	2.17	0.45
1:A:72:LYS:O	1:A:74:ARG:HG3	2.17	0.45
1:B:220:PRO:HG3	1:B:267:LEU:HD13	1.98	0.45
1:A:254:ASP:OD1	1:A:257:GLY:CA	2.65	0.44
1:B:201:PHE:HB2	1:B:209:GLN:HE22	1.82	0.44
1:B:148:ARG:NH1	1:B:148:ARG:HG3	2.32	0.44
1:A:36:LEU:CD2	1:A:150:PHE:CD2	3.00	0.44
1:A:123:HIS:HD2	1:A:125:ASP:N	2.14	0.44
1:A:3:LYS:HD3	1:A:3:LYS:N	2.33	0.44
1:A:36:LEU:CG	1:A:37:GLU:N	2.79	0.44
1:B:53:ILE:O	1:B:53:ILE:HG22	2.16	0.44
1:A:17:VAL:HG13	1:A:19:TYR:CE1	2.53	0.44
1:A:23:ASN:HD21	1:A:27:GLU:HG3	1.83	0.44
1:A:16:GLY:HA3	1:A:33:LYS:O	2.18	0.43
1:A:226:PRO:O	1:A:227:ASN:CB	2.65	0.43
1:B:221:ASN:OD1	1:B:221:ASN:C	2.55	0.43
1:B:4:TYR:CD1	1:B:4:TYR:N	2.86	0.43
1:A:5:HIS:HE1	1:A:23:ASN:O	2.02	0.43
1:A:123:HIS:CD2	1:A:125:ASP:H	2.29	0.43
1:B:98:GLU:O	1:B:99:SER:C	2.53	0.43
1:A:34:ILE:HG21	1:A:150:PHE:CZ	2.54	0.43
1:A:235:ASP:HA	1:A:236:PRO:HD3	1.92	0.43
1:A:36:LEU:CD1	1:A:150:PHE:HB3	2.41	0.43
1:B:64:LYS:NZ	2:B:2046:HOH:O	2.31	0.42
1:A:32:LYS:HB2	2:A:2005:HOH:O	2.19	0.42
1:B:259:ASP:O	1:B:263:LYS:HG2	2.20	0.42
1:A:118:ASP:OD2	1:A:276:LYS:HE3	2.18	0.42
1:B:4:TYR:HB3	1:B:21:ALA:HB1	2.01	0.42
1:A:219:THR:CG2	1:A:220:PRO:CD	2.97	0.42
1:A:36:LEU:HD12	1:A:36:LEU:HA	1.73	0.42
1:A:17:VAL:HG13	1:A:19:TYR:HE1	1.83	0.42
1:B:98:GLU:O	1:B:101:THR:N	2.46	0.42
1:B:219:THR:HA	1:B:220:PRO:HD3	1.90	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:15:TYR:H	1:B:32:LYS:NZ	2.18	0.42
1:B:53:ILE:O	1:B:56:GLU:HG2	2.20	0.42
1:A:152:ILE:HA	1:A:153:PRO:HD3	1.78	0.42
1:B:58:LYS:NZ	2:B:2044:HOH:O	2.41	0.41
1:B:71:THR:O	1:B:72:LYS:C	2.58	0.41
1:B:59:HIS:C	1:B:61:ASN:N	2.74	0.41
1:B:184:ILE:HG21	1:B:273:ILE:O	2.21	0.41
1:A:15:TYR:HB2	1:A:32:LYS:CE	2.50	0.41
1:A:177:LYS:HG2	1:A:178:TYR:N	2.36	0.41
1:A:2:GLU:HB2	1:A:4:TYR:CE1	2.56	0.41
1:A:32:LYS:HG2	1:A:34:ILE:HD13	2.02	0.40
1:A:25:TYR:HB2	1:A:27:GLU:CG	2.49	0.40
1:B:162:VAL:CG2	2:B:2016:HOH:O	2.69	0.40
1:B:221:ASN:OD1	1:B:223:LYS:HB2	2.21	0.40

All (8) 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	Distance(Å)	Clash(Å)
1:A:35:ARG:NE	1:A:35:ARG:NH2[2_655]	1.29	0.91
1:A:35:ARG:NE	1:A:35:ARG:CZ[2_655]	1.35	0.85
1:A:35:ARG:CD	1:A:35:ARG:NH2[2_655]	1.39	0.81
1:A:35:ARG:NE	1:A:35:ARG:NE[2_655]	1.39	0.81
2:B:2023:HOH:O	2:B:2128:HOH:O[3_455]	1.86	0.34
1:A:35:ARG:CZ	1:A:35:ARG:CZ[2_655]	1.94	0.26
1:A:35:ARG:CG	1:A:35:ARG:NH2[2_655]	1.98	0.22
1:A:94:GLU:O	1:A:94:GLU:O[2_555]	2.08	0.12

## 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	285/288 (99%)	265 (93%)	15 (5%)	5 (2%)	13	7
1	B	256/288 (89%)	234 (91%)	18 (7%)	4 (2%)	14	9
All	All	541/576 (94%)	499 (92%)	33 (6%)	9 (2%)	14	8

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	253	LEU
1	B	60	SER
1	A	72	LYS
1	A	251	LYS
1	A	252	GLY
1	A	37	GLU
1	B	163	THR
1	B	252	GLY
1	B	72	LYS

### 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	253/255 (99%)	247 (98%)	6 (2%)	61	73
1	B	229/255 (90%)	226 (99%)	3 (1%)	80	89
All	All	482/510 (94%)	473 (98%)	9 (2%)	69	81

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	THR
1	A	38	LYS
1	A	88	LYS
1	A	144	PHE
1	A	178	TYR
1	A	261	LEU
1	B	54	LEU
1	B	144	PHE
1	B	191	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	HIS

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Mol	Chain	Res	Type
1	A	61	ASN
1	A	123	HIS
1	A	130	ASN
1	B	61	ASN
1	B	123	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

There are no ligands in this entry.

### 5.7 Other polymers ⓘ

There are no such residues in this entry.

### 5.8 Polymer linkage issues

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	287/288 (99%)	0.23	17 (5%) 22 21	17, 29, 52, 69	0
1	B	262/288 (90%)	0.65	26 (9%) 8 7	21, 39, 60, 78	0
All	All	549/576 (95%)	0.43	43 (7%) 13 12	17, 34, 58, 78	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	73	LYS	11.6
1	B	72	LYS	7.9
1	B	135	ARG	6.9
1	B	74	ARG	6.6
1	B	70	HIS	6.2
1	B	252	GLY	4.8
1	A	235	ASP	4.6
1	B	150	PHE	4.5
1	B	51	ILE	4.5
1	A	2	GLU	4.0
1	B	71	THR	3.8
1	B	247	GLU	3.7
1	A	36	LEU	3.7
1	A	74	ARG	3.7
1	B	48	ILE	3.6
1	A	25	TYR	3.6
1	A	178	TYR	3.5
1	A	73	LYS	3.5
1	A	158	THR	3.5
1	A	3	LYS	3.3
1	B	178	TYR	3.2
1	A	37	GLU	3.2
1	B	3	LYS	3.1
1	B	93	CYS	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	35	ARG	3.0
1	B	92	VAL	2.9
1	A	1	MET	2.7
1	A	14	THR	2.6
1	B	187	VAL	2.6
1	B	245	PRO	2.5
1	A	251	LYS	2.5
1	B	244	LEU	2.4
1	B	68	VAL	2.4
1	A	215	ARG	2.3
1	B	188	GLY	2.3
1	A	159	HIS	2.3
1	B	190	ILE	2.3
1	B	5	HIS	2.2
1	A	223	LYS	2.1
1	B	246	TRP	2.1
1	B	254	ASP	2.1
1	B	4	TYR	2.0
1	B	95	GLY	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 ⓘ

There are no such residues in this entry.