



# Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 03:34 AM GMT

PDB ID : 1V7O  
Title : Alanyl-tRNA synthetase editing domain homologue protein from *Pyrococcus horikoshii*  
Authors : Okada, A.; Yao, M.; Sokabe, M.; Tanaka, I.  
Deposited on : 2003-12-18  
Resolution : 2.62 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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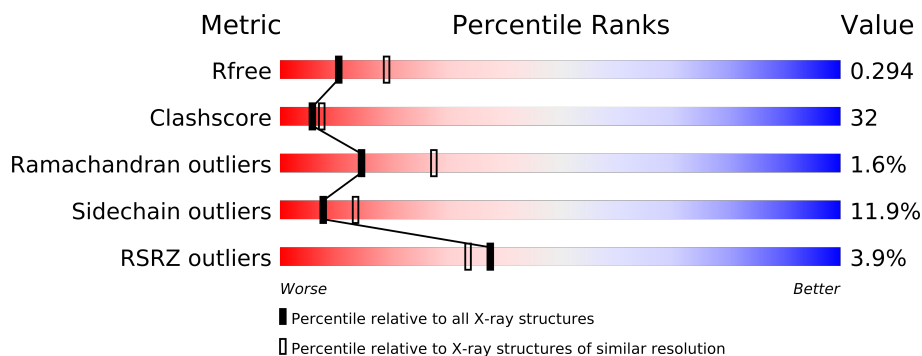
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.62 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	1940 (2.64-2.60)
Clashscore	79885	2404 (2.64-2.60)
Ramachandran outliers	78287	2360 (2.64-2.60)
Sidechain outliers	78261	2360 (2.64-2.60)
RSRZ outliers	66119	1939 (2.64-2.60)

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	165	
1	B	165	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2620 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 alanyl-tRNA synthetase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	155	Total	C	N	O	S	Se	0	0	0
			1265	811	217	233	1	3			
1	B	153	Total	C	N	O	S	Se	0	0	0
			1248	802	214	228	1	3			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP O58307
A	85	MSE	MET	MODIFIED RESIDUE	UNP O58307
A	90	MSE	MET	MODIFIED RESIDUE	UNP O58307
A	158	LEU	-	EXPRESSION TAG	UNP O58307
A	159	GLU	-	EXPRESSION TAG	UNP O58307
A	160	HIS	-	EXPRESSION TAG	UNP O58307
A	161	HIS	-	EXPRESSION TAG	UNP O58307
A	162	HIS	-	EXPRESSION TAG	UNP O58307
A	163	HIS	-	EXPRESSION TAG	UNP O58307
A	164	HIS	-	EXPRESSION TAG	UNP O58307
A	165	HIS	-	EXPRESSION TAG	UNP O58307
B	1	MSE	MET	MODIFIED RESIDUE	UNP O58307
B	85	MSE	MET	MODIFIED RESIDUE	UNP O58307
B	90	MSE	MET	MODIFIED RESIDUE	UNP O58307
B	158	LEU	-	EXPRESSION TAG	UNP O58307
B	159	GLU	-	EXPRESSION TAG	UNP O58307
B	160	HIS	-	EXPRESSION TAG	UNP O58307
B	161	HIS	-	EXPRESSION TAG	UNP O58307
B	162	HIS	-	EXPRESSION TAG	UNP O58307
B	163	HIS	-	EXPRESSION TAG	UNP O58307
B	164	HIS	-	EXPRESSION TAG	UNP O58307
B	165	HIS	-	EXPRESSION TAG	UNP O58307

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	54	Total 54	O 54	0	0
2	B	53	Total 53	O 53	0	0



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	34.19Å 88.74Å 110.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.62 24.30 – 2.62	Depositor EDS
% Data completeness (in resolution range)	99.2 (15.00-2.62) 99.4 (24.30-2.62)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.61 (at 2.64Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.235 , 0.273 0.250 , 0.294	Depositor DCC
$R_{free}$ test set	681 reflections (6.50%)	DCC
Wilson B-factor (Å <sup>2</sup> )	53.2	Xtriage
Anisotropy	0.269	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 35.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 10524 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	2620	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.59% 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	0.49	0/1286	0.66	0/1727
1	B	0.66	1/1269 (0.1%)	0.69	2/1704 (0.1%)
All	All	0.58	1/2555 (0.0%)	0.67	2/3431 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	151	LEU	C-N	-16.03	0.97	1.34

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	151	LEU	C-N-CA	5.32	134.99	121.70
1	B	153	LEU	N-CA-C	5.11	124.79	111.00

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	1265	0	1305	85	0
1	B	1248	0	1292	77	0
2	A	54	0	0	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	53	0	0	7	0
All	All	2620	0	2597	161	0

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 32.

All (161) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:77:LEU:HB3	1:A:78:PRO:HD2	1.23	1.13
1:A:78:PRO:HG3	1:A:81:GLU:HB2	1.11	1.06
1:B:121:THR:HG21	1:B:126:GLU:HB2	1.36	1.05
1:A:77:LEU:CB	1:A:78:PRO:HD2	1.97	0.95
1:A:77:LEU:HB3	1:A:78:PRO:CD	2.02	0.90
1:A:83:GLU:HA	1:A:90:MSE:HE1	1.54	0.89
1:A:59:GLU:OE2	1:A:131:LYS:HA	1.74	0.85
1:A:78:PRO:CG	1:A:81:GLU:HB2	2.04	0.83
1:A:79:ARG:NH1	1:A:100:VAL:O	2.13	0.81
1:B:12:LEU:HD13	1:B:33:THR:HG21	1.62	0.81
1:A:101:ARG:HD3	1:A:101:ARG:N	1.96	0.80
1:A:55:ILE:HD12	1:A:135:VAL:HG21	1.64	0.79
1:B:100:VAL:O	1:B:101:ARG:HD2	1.84	0.77
1:A:153:LEU:HD23	1:A:153:LEU:H	1.49	0.76
1:B:131:LYS:HB2	1:B:151:LEU:HD11	1.68	0.75
1:A:76:GLU:HG3	1:A:76:GLU:O	1.85	0.74
1:B:9:HIS:HD2	1:B:35:VAL:HG21	1.52	0.73
1:B:16:LYS:HE3	1:B:30:THR:OG1	1.88	0.73
1:A:78:PRO:HG3	1:A:81:GLU:CB	2.05	0.73
1:A:104:LYS:CE	1:A:117:ASN:HD22	2.03	0.72
1:A:9:HIS:NE2	1:A:118:LYS:HD2	2.04	0.72
1:B:73:LYS:NZ	1:B:109:GLU:OE2	2.21	0.71
1:A:101:ARG:HG3	1:A:101:ARG:HH11	1.58	0.69
1:A:87:GLY:O	1:A:90:MSE:HE2	1.92	0.69
1:A:7:ARG:HG2	1:A:7:ARG:HH11	1.56	0.69
1:B:77:LEU:N	1:B:77:LEU:HD22	2.09	0.68
1:B:151:LEU:N	1:B:151:LEU:HD12	2.07	0.68
1:B:121:THR:CG2	1:B:126:GLU:HB2	2.20	0.68
1:A:9:HIS:CE1	1:A:118:LYS:HB2	2.30	0.67
1:A:78:PRO:HG2	1:A:79:ARG:H	1.59	0.66
1:A:104:LYS:HE2	1:A:117:ASN:HD22	1.61	0.66
1:A:85:MSE:N	2:A:166:HOH:O	2.29	0.65
1:A:81:GLU:HA	1:A:81:GLU:OE1	1.97	0.64
1:B:33:THR:HG22	2:B:171:HOH:O	1.97	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:7:ARG:HG2	1:A:7:ARG:NH1	2.13	0.64
1:A:86:PHE:CE2	1:A:105:VAL:HG11	2.34	0.63
1:A:86:PHE:HE2	1:A:105:VAL:HG11	1.62	0.63
1:B:77:LEU:HD22	1:B:77:LEU:H	1.64	0.63
1:B:69:ASN:ND2	1:B:123:THR:HB	2.14	0.62
1:A:84:LYS:NZ	1:A:84:LYS:HB2	2.15	0.62
1:B:34:TYR:CE1	1:B:36:LYS:HB2	2.34	0.62
1:A:59:GLU:HB2	1:A:132:ILE:HD12	1.83	0.61
1:B:5:GLU:HB2	2:B:175:HOH:O	2.01	0.60
1:A:84:LYS:N	2:A:166:HOH:O	2.35	0.59
1:B:55:ILE:HD12	1:B:135:VAL:HG21	1.84	0.59
1:A:155:ASN:HA	2:A:203:HOH:O	2.02	0.59
1:A:9:HIS:HB2	2:A:171:HOH:O	2.02	0.58
1:B:59:GLU:HB2	1:B:132:ILE:HD12	1.84	0.58
1:A:79:ARG:O	1:A:83:GLU:HB2	2.02	0.58
1:B:1:MSE:N	1:B:1:MSE:HE3	2.18	0.58
1:B:7:ARG:NH1	1:B:126:GLU:O	2.37	0.58
1:B:77:LEU:O	1:B:102:ILE:HD13	2.04	0.57
1:B:86:PHE:CE2	1:B:105:VAL:HG11	2.40	0.57
1:B:152:GLU:C	1:B:153:LEU:HD12	2.25	0.57
1:B:78:PRO:HD2	1:B:81:GLU:OE1	2.05	0.56
1:B:56:ARG:NH1	1:B:56:ARG:HG3	2.20	0.56
1:B:73:LYS:HA	2:B:215:HOH:O	2.05	0.56
1:A:9:HIS:HE1	1:A:116:CYS:SG	2.29	0.56
1:A:9:HIS:CD2	1:A:118:LYS:HD2	2.41	0.55
1:A:80:GLU:HA	1:A:83:GLU:HB2	1.87	0.55
1:B:129:PRO:O	1:B:151:LEU:HD13	2.07	0.55
1:A:51:SER:OG	1:A:54:GLU:HG3	2.07	0.55
1:A:7:ARG:NH1	1:A:127:ILE:O	2.40	0.55
1:B:56:ARG:HH11	1:B:56:ARG:HG3	1.71	0.54
1:A:17:GLY:HA3	1:A:111:TRP:CD1	2.42	0.54
1:B:129:PRO:HG2	1:B:151:LEU:HB2	1.90	0.54
1:B:151:LEU:N	1:B:151:LEU:CD1	2.70	0.54
1:B:140:SER:HB3	2:B:207:HOH:O	2.08	0.53
1:A:13:HIS:CE1	1:A:114:ASN:HB3	2.43	0.53
1:A:7:ARG:HB3	1:A:150:LEU:CD2	2.38	0.53
1:A:68:GLU:O	1:A:68:GLU:HG2	2.07	0.53
1:A:133:ARG:NH1	1:A:149:GLU:OE1	2.39	0.53
1:A:101:ARG:CG	1:A:101:ARG:HH11	2.22	0.53
1:A:81:GLU:C	2:A:166:HOH:O	2.47	0.53
1:B:102:ILE:HG23	1:B:102:ILE:O	2.10	0.52
1:B:134:LYS:NZ	1:B:134:LYS:HB3	2.24	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:129:PRO:HG3	1:A:153:LEU:HD21	1.91	0.52
1:B:121:THR:CG2	1:B:126:GLU:OE1	2.58	0.52
1:B:121:THR:HG21	1:B:126:GLU:CB	2.25	0.51
1:A:52:ASP:O	1:A:56:ARG:HG3	2.10	0.51
1:A:128:GLY:HA3	1:A:155:ASN:OD1	2.11	0.51
1:B:22:VAL:HG22	1:B:57:GLU:HG3	1.93	0.50
1:A:100:VAL:C	1:A:101:ARG:HD3	2.31	0.50
1:A:13:HIS:HE1	1:A:114:ASN:HB3	1.76	0.50
1:B:110:ASP:OD1	1:B:110:ASP:O	2.30	0.50
1:A:154:GLU:HA	1:A:154:GLU:OE1	2.12	0.50
1:A:60:ARG:HD2	1:A:60:ARG:N	2.27	0.50
1:B:69:ASN:HD22	1:B:125:GLY:H	1.60	0.49
1:B:122:LYS:CD	1:B:123:THR:HG23	2.42	0.49
1:B:34:TYR:HE1	1:B:36:LYS:HB2	1.76	0.49
1:B:12:LEU:CD1	1:B:33:THR:HG21	2.38	0.48
1:B:63:ASN:OD1	1:B:129:PRO:HA	2.12	0.48
1:A:9:HIS:CE1	1:A:116:CYS:SG	3.06	0.48
1:A:84:LYS:HZ2	1:A:84:LYS:HB2	1.76	0.48
1:A:80:GLU:HA	1:A:83:GLU:CB	2.44	0.48
1:B:138:ARG:HD2	1:B:145:GLU:OE2	2.14	0.48
1:B:106:VAL:HG21	1:B:120:HIS:CD2	2.49	0.47
1:A:76:GLU:CG	1:A:76:GLU:O	2.57	0.47
1:B:102:ILE:CG2	1:B:102:ILE:O	2.63	0.47
1:A:86:PHE:CD2	1:A:105:VAL:HG21	2.50	0.47
1:A:86:PHE:HB3	1:A:113:VAL:HG11	1.96	0.47
1:A:96:VAL:CG1	1:A:97:PRO:HD2	2.45	0.47
1:B:121:THR:HG22	1:B:126:GLU:OE1	2.14	0.47
1:A:153:LEU:HD23	1:A:153:LEU:N	2.24	0.47
1:B:51:SER:O	1:B:55:ILE:HG12	2.15	0.47
1:B:75:TYR:HB2	1:B:105:VAL:HG13	1.96	0.47
1:A:23:LEU:CD1	1:A:23:LEU:N	2.78	0.46
1:B:77:LEU:H	1:B:77:LEU:CD2	2.28	0.46
1:B:15:VAL:O	1:B:19:VAL:HG23	2.16	0.46
1:A:53:GLU:CD	1:A:53:GLU:C	2.73	0.46
1:A:77:LEU:CB	1:A:78:PRO:CD	2.74	0.46
1:A:27:ALA:O	1:A:44:VAL:HG21	2.15	0.46
1:B:150:LEU:C	1:B:151:LEU:HD12	2.36	0.46
1:A:93:LEU:HD23	1:A:93:LEU:HA	1.71	0.46
1:B:87:GLY:HA3	2:B:188:HOH:O	2.15	0.46
1:A:104:LYS:N	1:A:104:LYS:HD3	2.31	0.46
1:A:8:THR:HG22	1:A:35:VAL:HG22	1.98	0.45
1:B:3:SER:O	1:B:7:ARG:HD3	2.16	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:8:THR:HG21	1:B:35:VAL:HG13	1.98	0.45
1:B:122:LYS:HD2	1:B:123:THR:HG23	1.98	0.45
1:B:6:VAL:HG13	1:B:119:GLU:O	2.16	0.45
1:A:101:ARG:CG	1:A:101:ARG:NH1	2.79	0.45
1:B:50:PRO:HB2	1:B:55:ILE:HD11	1.98	0.45
1:B:33:THR:CG2	1:B:34:TYR:N	2.80	0.45
1:A:83:GLU:O	1:A:87:GLY:HA2	2.17	0.45
1:B:77:LEU:N	1:B:77:LEU:CD2	2.78	0.44
1:B:43:ILE:HG12	1:B:145:GLU:HG2	2.00	0.44
1:A:94:PHE:HB2	1:B:94:PHE:CZ	2.52	0.44
1:B:1:MSE:HE3	1:B:1:MSE:H3	1.82	0.44
1:B:136:ARG:NH2	1:B:138:ARG:HD2	2.33	0.44
1:A:7:ARG:HB3	1:A:150:LEU:HD21	1.98	0.44
1:A:3:SER:O	1:A:7:ARG:HD3	2.18	0.44
1:A:137:PHE:CE2	1:A:139:LYS:HA	2.53	0.44
1:B:12:LEU:HD13	1:B:33:THR:CG2	2.42	0.43
1:A:104:LYS:HE2	1:A:117:ASN:ND2	2.32	0.43
1:B:2:TYR:HE1	1:B:126:GLU:OE2	2.01	0.43
1:B:135:VAL:CG1	1:B:144:LEU:HD11	2.49	0.43
1:B:8:THR:CG2	1:B:35:VAL:HG13	2.49	0.43
1:A:32:SER:HA	2:A:204:HOH:O	2.18	0.43
1:B:21:LYS:HD3	1:B:21:LYS:HA	1.85	0.43
1:A:84:LYS:CB	1:A:84:LYS:NZ	2.82	0.43
1:A:9:HIS:ND1	1:A:120:HIS:CE1	2.86	0.43
1:B:118:LYS:NZ	2:B:171:HOH:O	2.52	0.42
1:B:56:ARG:HH11	1:B:56:ARG:CG	2.33	0.42
1:B:131:LYS:HD3	1:B:132:ILE:N	2.35	0.42
1:B:134:LYS:HB3	1:B:134:LYS:HZ3	1.84	0.42
1:A:22:VAL:HG12	1:A:23:LEU:HD12	2.01	0.42
1:A:78:PRO:HG2	1:A:79:ARG:N	2.31	0.42
1:A:96:VAL:HG12	1:A:97:PRO:HD2	2.02	0.42
1:A:23:LEU:HD12	1:A:23:LEU:N	2.35	0.41
1:B:93:LEU:HA	1:B:93:LEU:HD23	1.89	0.41
1:B:79:ARG:HD2	1:B:101:ARG:CZ	2.50	0.41
1:B:45:LYS:NZ	2:B:178:HOH:O	2.53	0.41
1:A:106:VAL:HG21	1:A:120:HIS:CD2	2.56	0.41
1:A:136:ARG:HD2	2:A:188:HOH:O	2.19	0.41
1:B:12:LEU:HD11	1:B:35:VAL:CG2	2.51	0.40
1:B:67:LYS:HD2	1:B:67:LYS:HA	1.98	0.40
1:B:71:PRO:HA	1:B:123:THR:HG22	2.03	0.40
1:A:29:TRP:HB2	2:A:168:HOH:O	2.20	0.40
1:A:59:GLU:OE2	1:A:131:LYS:CA	2.59	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:57:GLU:HA	1:A:57:GLU:OE1	2.21	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	153/165 (93%)	136 (89%)	13 (8%)	4 (3%)	8	13
1	B	151/165 (92%)	139 (92%)	11 (7%)	1 (1%)	30	57
All	All	304/330 (92%)	275 (90%)	24 (8%)	5 (2%)	14	27

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	78	PRO
1	A	79	ARG
1	B	102	ILE
1	A	102	ILE
1	A	95	PRO

### 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	140/147 (95%)	122 (87%)	18 (13%)	6	11
1	B	138/147 (94%)	123 (89%)	15 (11%)	9	16
All	All	278/294 (95%)	245 (88%)	33 (12%)	8	13

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ARG
1	A	44	VAL
1	A	53	GLU
1	A	59	GLU
1	A	60	ARG
1	A	81	GLU
1	A	83	GLU
1	A	84	LYS
1	A	85	MSE
1	A	93	LEU
1	A	98	GLU
1	A	101	ARG
1	A	102	ILE
1	A	107	VAL
1	A	109	GLU
1	A	117	ASN
1	A	136	ARG
1	A	153	LEU
1	B	1	MSE
1	B	56	ARG
1	B	60	ARG
1	B	61	LEU
1	B	69	ASN
1	B	77	LEU
1	B	78	PRO
1	B	79	ARG
1	B	83	GLU
1	B	93	LEU
1	B	99	ASP
1	B	101	ARG
1	B	102	ILE
1	B	131	LYS
1	B	152	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	112	ASN
1	A	117	ASN
1	A	120	HIS

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Mol	Chain	Res	Type
1	B	9	HIS
1	B	13	HIS
1	B	69	ASN

### 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	155/165 (93%)	0.06	5 (3%)	45 42	23, 47, 81, 100	0
1	B	153/165 (92%)	0.06	7 (4%)	31 27	23, 49, 75, 100	0
All	All	308/330 (93%)	0.06	12 (3%)	37 33	23, 48, 78, 100	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	153	LEU	11.6
1	B	1	MSE	6.0
1	B	152	GLU	5.5
1	A	153	LEU	4.8
1	A	154	GLU	3.4
1	A	102	ILE	3.1
1	A	155	ASN	2.6
1	B	84	LYS	2.4
1	B	67	LYS	2.3
1	A	100	VAL	2.2
1	B	2	TYR	2.1
1	B	4	ILE	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 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.