



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 07:50 PM GMT

PDB ID : 3VPA  
Title : Staphylococcus aureus FtsZ apo-form  
Authors : Matsui, T.; Yamane, J.; Mogi, N.; Yao, M.; Tanaka, I.  
Deposited on : 2012-02-28  
Resolution : 2.49 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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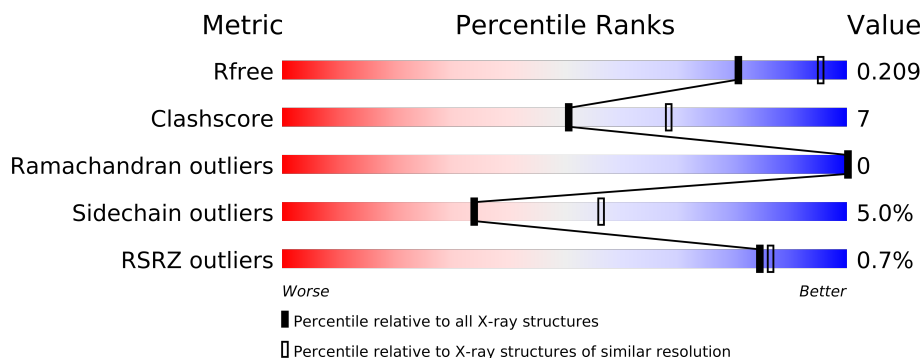
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.49 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	3277 (2.50-2.46)
Clashscore	79885	4136 (2.50-2.46)
Ramachandran outliers	78287	4052 (2.50-2.46)
Sidechain outliers	78261	4054 (2.50-2.46)
RSRZ outliers	66119	3279 (2.50-2.46)

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	308	
1	B	308	
1	C	308	
1	D	308	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8492 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 protein FtsZ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2067	1284	349	423	11			
1	B	280	Total	C	N	O	S	0	0	0
			2008	1246	337	414	11			
1	C	300	Total	C	N	O	S	0	0	0
			2163	1340	370	442	11			
1	D	295	Total	C	N	O	S	0	0	0
			2115	1311	359	433	12			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	GLY	-	EXPRESSION TAG	UNP P0A029
A	10	HIS	-	EXPRESSION TAG	UNP P0A029
A	11	MET	-	EXPRESSION TAG	UNP P0A029
B	9	GLY	-	EXPRESSION TAG	UNP P0A029
B	10	HIS	-	EXPRESSION TAG	UNP P0A029
B	11	MET	-	EXPRESSION TAG	UNP P0A029
C	9	GLY	-	EXPRESSION TAG	UNP P0A029
C	10	HIS	-	EXPRESSION TAG	UNP P0A029
C	11	MET	-	EXPRESSION TAG	UNP P0A029
D	9	GLY	-	EXPRESSION TAG	UNP P0A029
D	10	HIS	-	EXPRESSION TAG	UNP P0A029
D	11	MET	-	EXPRESSION TAG	UNP P0A029

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	36	Total	O	0	0
			36	36		
2	B	14	Total	O	0	0
			14	14		

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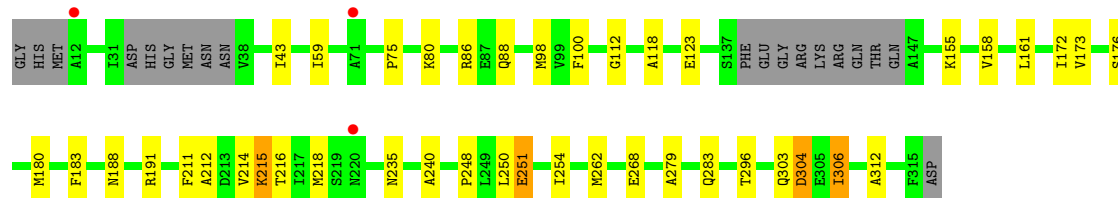
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	34	Total 34	O 34	0	0
2	D	55	Total 55	O 55	0	0

### 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.

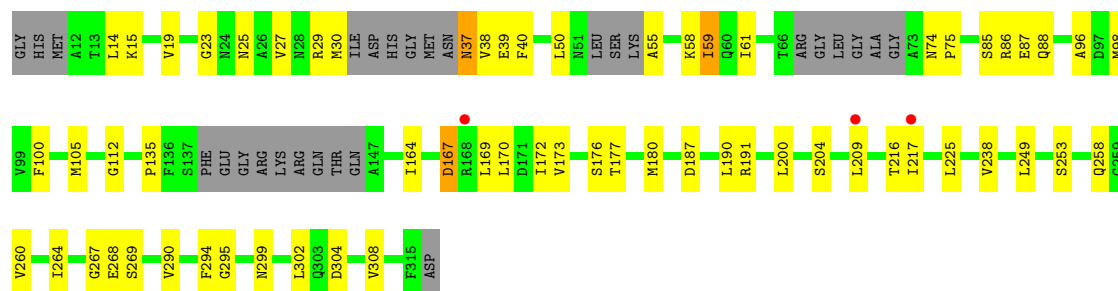
- Molecule 1: Cell division protein FtsZ

Chain A: 



- Molecule 1: Cell division protein FtsZ

Chain B: 



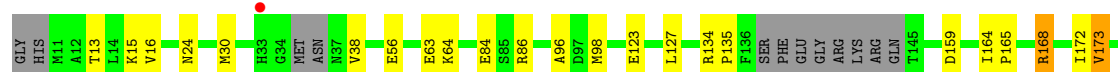
- Molecule 1: Cell division protein FtsZ

Chain C: 



- Molecule 1: Cell division protein FtsZ

Chain D: 



M180	L200	L201	S204	N208	L209	D210	D213	V214	K215	T216	I217	N235	E239	K242	S246	L249	T252	S253	I254	V255	M262	G267	E268	K279	D280	I281	N299	P300	F315	ASP
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.48Å 81.06Å 225.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.49 – 2.49 44.73 – 2.49	Depositor EDS
% Data completeness (in resolution range)	98.7 (35.49-2.49) 98.7 (44.73-2.49)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.25 (at 2.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, $R_{free}$	0.210 , 0.258 0.207 , 0.209	Depositor DCC
$R_{free}$ test set	2220 reflections (4.66%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.5	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 36.2	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 47666 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8492	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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	0.48	0/2079	0.61	0/2807
1	B	0.44	0/2018	0.61	0/2724
1	C	0.47	0/2177	0.59	0/2938
1	D	0.51	0/2128	0.63	0/2873
All	All	0.47	0/8402	0.61	0/11342

There are no bond length outliers.

There are no bond angle outliers.

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	2067	0	2101	28	0
1	B	2008	0	2027	42	0
1	C	2163	0	2192	28	0
1	D	2115	0	2140	30	0
2	A	36	0	0	0	0
2	B	14	0	0	0	0
2	C	34	0	0	1	0
2	D	55	0	0	1	0
All	All	8492	0	8460	124	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 7.

All (124) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:159:ASP:HB2	1:C:190:LEU:HD13	1.72	0.72
1:C:86:ARG:NH1	1:C:90:GLU:OE1	2.24	0.70
1:A:250:LEU:O	1:A:254:ILE:HG12	1.96	0.65
1:C:287:ASP:N	1:C:287:ASP:OD1	2.29	0.63
1:D:30:MET:HE1	1:D:172:ILE:HD11	1.80	0.63
1:A:248:PRO:HA	1:A:251:GLU:OE1	1.99	0.62
1:A:43:ILE:HD13	1:A:59:ILE:HB	1.82	0.62
1:A:211:PHE:CZ	1:A:215:LYS:HE3	2.34	0.61
1:B:59:ILE:HD13	1:B:88:GLN:HG2	1.81	0.61
1:C:207:VAL:HG12	1:C:209:LEU:HG	1.83	0.60
1:D:86:ARG:NH2	1:D:123:GLU:OE1	2.23	0.60
1:B:164:ILE:HG22	1:B:204:SER:HB3	1.82	0.60
1:B:100:PHE:CD2	1:B:172:ILE:HD12	2.39	0.58
1:B:264:ILE:HD12	1:B:294:PHE:HE1	1.69	0.57
1:B:15:LYS:NZ	1:B:39:GLU:OE1	2.28	0.57
1:D:235:ASN:O	1:D:239:GLU:HG2	2.05	0.57
1:B:55:ALA:HB3	1:B:58:LYS:HE3	1.87	0.57
1:B:50:LEU:HD23	1:B:58:LYS:HB3	1.87	0.56
1:B:98:MET:HE1	1:B:180:MET:HA	1.88	0.56
1:B:190:LEU:HD11	1:B:200:LEU:HB3	1.87	0.56
1:A:86:ARG:NH2	1:A:123:GLU:OE1	2.31	0.56
1:B:169:LEU:O	1:B:172:ILE:HG12	2.06	0.56
1:D:267:GLY:HA2	1:D:299:ASN:O	2.05	0.56
1:D:63:GLU:H	1:D:63:GLU:CD	2.09	0.55
1:A:155:LYS:HE3	1:A:218:MET:HG2	1.86	0.55
1:D:98:MET:SD	1:D:180:MET:HG2	2.47	0.55
1:A:191:ARG:NE	1:B:191:ARG:HD3	2.22	0.55
1:B:27:VAL:HG22	1:B:40:PHE:HD1	1.72	0.54
1:A:211:PHE:CE1	1:A:215:LYS:HE3	2.42	0.54
1:B:37:ASN:N	1:B:37:ASN:HD22	2.06	0.54
1:B:299:ASN:HB3	1:B:302:LEU:HD13	1.91	0.53
1:B:87:GLU:OE1	1:B:87:GLU:N	2.42	0.53
1:C:198:SER:N	1:C:222:GLY:O	2.36	0.53
1:A:188:ASN:OD1	1:A:191:ARG:NH1	2.42	0.52
1:B:167:ASP:N	1:B:167:ASP:OD1	2.26	0.52
1:D:86:ARG:HD2	2:D:427:HOH:O	2.08	0.52
1:A:176:SER:O	1:A:180:MET:HG3	2.10	0.52
1:B:264:ILE:HG12	1:B:308:VAL:HG22	1.90	0.52
1:C:60:GLN:NE2	1:C:63:GLU:OE2	2.44	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:252:THR:O	1:D:255:VAL:HG22	2.10	0.51
1:B:100:PHE:HD2	1:B:172:ILE:HD12	1.74	0.51
1:C:73:ALA:N	2:C:422:HOH:O	2.35	0.51
1:C:218:MET:HE1	1:C:223:SER:HB2	1.93	0.51
1:D:165:PRO:HD2	1:D:168:ARG:HB2	1.91	0.51
1:C:75:PRO:O	1:C:112:GLY:HA3	2.09	0.51
1:A:254:ILE:HD13	1:A:312:ALA:HB2	1.92	0.50
1:B:74:ASN:HB2	1:C:69:LEU:HD21	1.93	0.50
1:D:210:ASP:OD1	1:D:213:ASP:N	2.45	0.50
1:C:45:THR:HG21	1:C:109:THR:HG21	1.93	0.49
1:A:262:MET:HE1	1:A:279:ALA:HA	1.95	0.49
1:D:201:ILE:HD11	1:D:217:ILE:HG21	1.95	0.49
1:A:283:GLN:NE2	1:A:283:GLN:O	2.44	0.49
1:A:240:ALA:HB3	1:A:306:ILE:HD11	1.94	0.49
1:A:75:PRO:O	1:A:112:GLY:HA3	2.12	0.49
1:C:59:ILE:HD11	1:C:92:ALA:CB	2.43	0.49
1:D:16:VAL:HG23	1:D:38:VAL:HG21	1.95	0.48
1:A:250:LEU:HD12	1:A:254:ILE:HD11	1.95	0.48
1:D:15:LYS:HD2	1:D:96:ALA:HB2	1.94	0.48
1:A:172:ILE:HG13	1:A:173:VAL:N	2.29	0.48
1:A:212:ALA:C	1:A:214:VAL:H	2.17	0.48
1:D:208:ASN:ND2	1:D:208:ASN:O	2.46	0.47
1:B:187:ASP:OD1	1:B:191:ARG:HD2	2.15	0.47
1:D:262:MET:SD	1:D:279:ALA:HA	2.55	0.47
1:C:175:LYS:HD2	1:C:230:VAL:HG21	1.96	0.47
1:C:292:MET:HG2	1:C:294:PHE:CE2	2.50	0.46
1:A:191:ARG:HD3	1:B:187:ASP:OD2	2.16	0.45
1:B:61:ILE:HG22	1:B:85:SER:HB2	1.98	0.45
1:D:209:LEU:HD22	1:D:213:ASP:HB3	1.99	0.45
1:B:225:LEU:HD22	1:B:249:LEU:HD21	1.99	0.44
1:B:267:GLY:HA2	1:B:299:ASN:O	2.17	0.44
1:A:80:LYS:HD3	1:A:80:LYS:HA	1.76	0.44
1:C:159:ASP:OD1	1:C:160:THR:HG22	2.17	0.44
1:A:100:PHE:HD2	1:A:172:ILE:HD12	1.81	0.44
1:B:260:VAL:HG23	1:B:290:VAL:HG21	1.99	0.44
1:B:176:SER:O	1:B:180:MET:HG3	2.17	0.44
1:C:114:ALA:HB3	1:C:115:PRO:HD3	2.00	0.44
1:B:225:LEU:HD12	1:B:253:SER:OG	2.17	0.44
1:C:138:PHE:HE2	1:C:217:ILE:HD13	1.82	0.44
1:B:19:VAL:N	1:B:23:GLY:HA3	2.32	0.44
1:D:38:VAL:HG11	1:D:173:VAL:HG12	2.00	0.44
1:C:86:ARG:HD2	1:C:120:ILE:HG12	1.99	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:100:PHE:CD2	1:A:172:ILE:HD12	2.52	0.43
1:B:105:MET:HE1	1:B:135:PRO:HD3	1.99	0.43
1:D:242:LYS:O	1:D:246:SER:HB3	2.18	0.43
1:C:66:THR:HA	1:C:69:LEU:HD12	2.00	0.43
1:B:61:ILE:HG22	1:B:85:SER:CB	2.49	0.43
1:A:98:MET:HE1	1:A:183:PHE:HB2	2.00	0.43
1:D:56:GLU:CD	1:D:56:GLU:H	2.21	0.43
1:D:64:LYS:HG3	1:D:84:GLU:OE1	2.17	0.43
1:A:296:THR:O	1:B:295:GLY:HA2	2.19	0.43
1:D:254:ILE:HG13	1:D:254:ILE:O	2.19	0.43
1:C:209:LEU:HB2	1:C:249:LEU:HD22	2.01	0.42
1:B:75:PRO:O	1:B:112:GLY:HA3	2.18	0.42
1:B:98:MET:HE3	1:B:100:PHE:HE1	1.84	0.42
1:D:164:ILE:HG22	1:D:204:SER:HB3	2.00	0.42
1:B:25:ASN:O	1:B:29:ARG:HG3	2.19	0.42
1:A:212:ALA:O	1:A:214:VAL:HG23	2.19	0.42
1:B:172:ILE:HG13	1:B:173:VAL:HG23	2.01	0.42
1:B:30:MET:SD	1:B:173:VAL:HG21	2.60	0.42
1:C:191:ARG:HB3	1:C:191:ARG:HE	1.62	0.42
1:B:15:LYS:HE2	1:B:96:ALA:HB2	2.02	0.42
1:D:168:ARG:NE	1:D:168:ARG:HA	2.34	0.41
1:C:17:ILE:HG12	1:C:41:ILE:HB	2.01	0.41
1:A:118:ALA:HB1	1:A:158:VAL:HG12	2.02	0.41
1:C:241:ALA:HA	1:C:308:VAL:HG11	2.02	0.41
1:D:268:GLU:OE1	1:D:268:GLU:N	2.52	0.41
1:A:98:MET:HE2	1:A:180:MET:HG2	2.02	0.41
1:C:38:VAL:HG21	1:C:173:VAL:HG22	2.03	0.41
1:A:303:GLN:HB2	1:A:304:ASP:H	1.59	0.41
1:B:55:ALA:O	1:B:58:LYS:HE3	2.21	0.41
1:D:98:MET:CE	1:D:127:LEU:HD23	2.50	0.41
1:B:19:VAL:H	1:B:23:GLY:HA3	1.85	0.41
1:D:127:LEU:HD12	1:D:159:ASP:OD2	2.21	0.41
1:D:249:LEU:HA	1:D:249:LEU:HD23	1.63	0.41
1:C:137:SER:HB3	1:C:142:LYS:HD3	2.02	0.41
1:C:242:LYS:O	1:C:246:SER:HB3	2.21	0.41
1:B:14:LEU:HD13	1:B:98:MET:HG2	2.03	0.40
1:C:74:ASN:HA	1:C:75:PRO:HD3	1.95	0.40
1:D:200:LEU:HD12	1:D:201:ILE:N	2.35	0.40
1:C:59:ILE:HD11	1:C:92:ALA:HB2	2.03	0.40
1:B:172:ILE:HG13	1:B:173:VAL:N	2.36	0.40
1:B:105:MET:CE	1:B:135:PRO:HD3	2.52	0.40
1:D:299:ASN:HA	1:D:300:PRO:HD3	1.89	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:134:ARG:HA	1:D:135:PRO:HD3	1.97	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	283/308 (92%)	279 (99%)	4 (1%)	0	100	100
1	B	270/308 (88%)	265 (98%)	5 (2%)	0	100	100
1	C	296/308 (96%)	292 (99%)	4 (1%)	0	100	100
1	D	289/308 (94%)	286 (99%)	3 (1%)	0	100	100
All	All	1138/1232 (92%)	1122 (99%)	16 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	217/233 (93%)	208 (96%)	9 (4%)	41	66
1	B	212/233 (91%)	197 (93%)	15 (7%)	21	36
1	C	227/233 (97%)	216 (95%)	11 (5%)	35	58
1	D	222/233 (95%)	213 (96%)	9 (4%)	41	66
All	All	878/932 (94%)	834 (95%)	44 (5%)	34	56

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	GLN
1	A	161	LEU
1	A	215	LYS
1	A	216	THR
1	A	235	ASN
1	A	251	GLU
1	A	268	GLU
1	A	304	ASP
1	A	306	ILE
1	B	37	ASN
1	B	38	VAL
1	B	59	ILE
1	B	86	ARG
1	B	167	ASP
1	B	170	LEU
1	B	177	THR
1	B	209	LEU
1	B	216	THR
1	B	217	ILE
1	B	238	VAL
1	B	258	GLN
1	B	268	GLU
1	B	269	SER
1	B	304	ASP
1	C	39	GLU
1	C	45	THR
1	C	91	ASP
1	C	168	ARG
1	C	172	ILE
1	C	173	VAL
1	C	230	VAL
1	C	238	VAL
1	C	254	ILE
1	C	277	GLU
1	C	304	ASP
1	D	13	THR
1	D	24	ASN
1	D	168	ARG
1	D	173	VAL
1	D	208	ASN
1	D	210	ASP
1	D	215	LYS
1	D	254	ILE

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Mol	Chain	Res	Type
1	D	281	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	303	GLN
1	C	60	GLN
1	D	146	GLN

### 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	289/308 (93%)	-0.08	3 (1%) 79 80	33, 53, 82, 100	0
1	B	280/308 (90%)	0.16	3 (1%) 77 78	36, 65, 86, 96	0
1	C	300/308 (97%)	0.01	1 (0%) 91 94	34, 59, 82, 95	0
1	D	295/308 (95%)	-0.10	1 (0%) 91 94	29, 46, 74, 89	0
All	All	1164/1232 (94%)	-0.00	8 (0%) 84 86	29, 55, 82, 100	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	33	HIS	2.7
1	A	12	ALA	2.6
1	A	71	ALA	2.3
1	A	220	ASN	2.2
1	C	144	GLN	2.2
1	B	209	LEU	2.2
1	B	217	ILE	2.2
1	B	168	ARG	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.