



# Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 12:31 AM GMT

PDB ID : 2GUJ  
Title : Three dimensional structure of the protein P54332 from Bacillus Subtilis.  
Northeast Structural Genomics Consortium target sr353.  
Authors : Kuzin, A.P.; Zhou, W.; Seetharaman, J.; Cunningham, K.; Janjua, H.;  
Konover, K.; Ma, L.C.; Xiao, R.; Acton, T.; Montelione, G.; Tong, L.; Hunt,  
J.F.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2006-04-30  
Resolution : 3.00 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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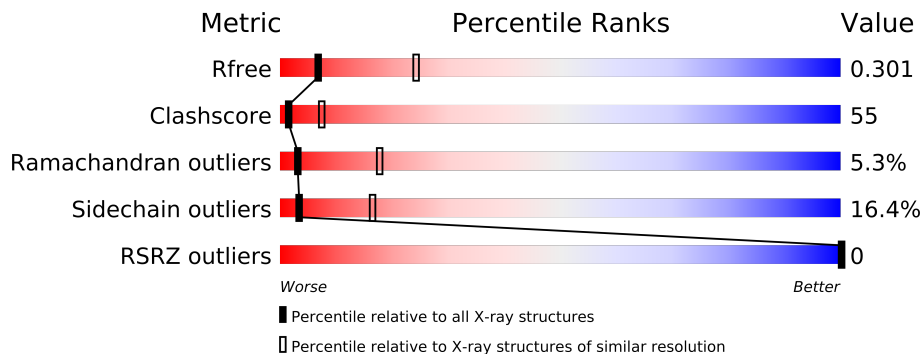
The following versions of software and data (see [references](#)) were used in the production of this report:

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 3.00 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))
$R_{free}$	66092	1216 (3.00-3.00)
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)

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 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	155	
1	B	155	

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2116 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 Phage-like element PBSX protein xkdM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	133	Total	C	N	O	Se	0	0	0
			1042	654	172	211	5			
1	B	137	Total	C	N	O	Se	0	0	0
			1074	674	176	219	5			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP P54332
A	23	MSE	MET	MODIFIED RESIDUE	UNP P54332
A	43	MSE	MET	MODIFIED RESIDUE	UNP P54332
A	47	MSE	MET	MODIFIED RESIDUE	UNP P54332
A	74	MSE	MET	MODIFIED RESIDUE	UNP P54332
A	75	MSE	MET	MODIFIED RESIDUE	UNP P54332
A	148	LEU	-	EXPRESSION TAG	UNP P54332
A	149	GLU	-	EXPRESSION TAG	UNP P54332
A	150	HIS	-	EXPRESSION TAG	UNP P54332
A	151	HIS	-	EXPRESSION TAG	UNP P54332
A	152	HIS	-	EXPRESSION TAG	UNP P54332
A	153	HIS	-	EXPRESSION TAG	UNP P54332
A	154	HIS	-	EXPRESSION TAG	UNP P54332
A	155	HIS	-	EXPRESSION TAG	UNP P54332
B	1	MSE	MET	MODIFIED RESIDUE	UNP P54332
B	23	MSE	MET	MODIFIED RESIDUE	UNP P54332
B	43	MSE	MET	MODIFIED RESIDUE	UNP P54332
B	47	MSE	MET	MODIFIED RESIDUE	UNP P54332
B	74	MSE	MET	MODIFIED RESIDUE	UNP P54332
B	75	MSE	MET	MODIFIED RESIDUE	UNP P54332
B	148	LEU	-	EXPRESSION TAG	UNP P54332
B	149	GLU	-	EXPRESSION TAG	UNP P54332
B	150	HIS	-	EXPRESSION TAG	UNP P54332
B	151	HIS	-	EXPRESSION TAG	UNP P54332
B	152	HIS	-	EXPRESSION TAG	UNP P54332

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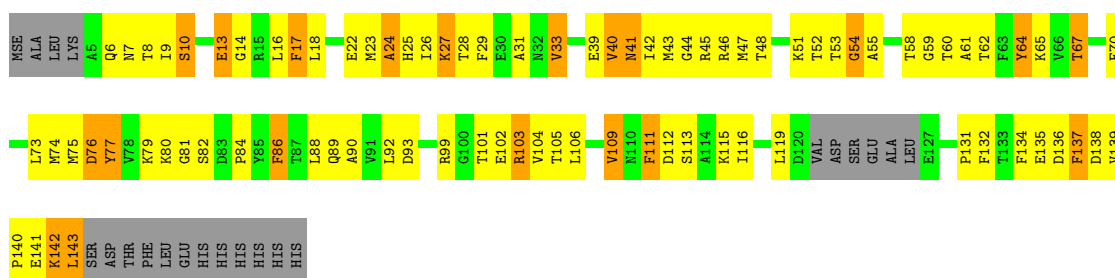
Chain	Residue	Modelled	Actual	Comment	Reference
B	153	HIS	-	EXPRESSION TAG	UNP P54332
B	154	HIS	-	EXPRESSION TAG	UNP P54332
B	155	HIS	-	EXPRESSION TAG	UNP P54332

### 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.

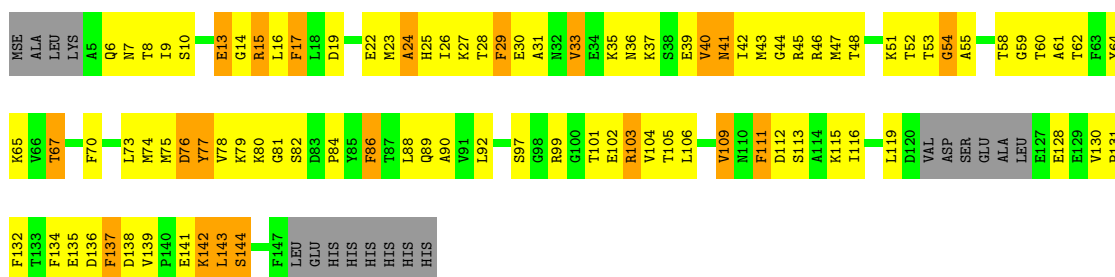
- Molecule 1: Phage-like element PBSX protein xkdM

Chain A:



- Molecule 1: Phage-like element PBSX protein xkdM

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.25Å 84.25Å 46.69Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.75 – 3.00 28.75 – 2.99	Depositor EDS
% Data completeness (in resolution range)	73.1 (28.75-3.00) 91.4 (28.75-2.99)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.16 (at 3.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.253 , 0.285 0.277 , 0.301	Depositor DCC
$R_{free}$ test set	329 reflections (4.85%)	DCC
Wilson B-factor (Å <sup>2</sup> )	47.3	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 20.4	EDS
Estimated twinning fraction	0.032 for -h,-k,l 0.458 for h,-h-k,-l 0.032 for -k,-h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 13213 reflections	Xtriage
$F_o, F_c$ correlation	0.81	EDS
Total number of atoms	2116	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.55% 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	0.47	0/1053	0.71	0/1406
1	B	0.44	0/1086	0.69	0/1451
All	All	0.45	0/2139	0.70	0/2857

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	1042	0	1012	112	0
1	B	1074	0	1037	117	0
All	All	2116	0	2049	229	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 55.

All (229) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:40:VAL:HG13	1:A:41:ASN:H	1.09	1.12
1:B:40:VAL:HG13	1:B:41:ASN:H	1.05	1.09

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:74:MSE:HE2	1:A:116:ILE:HD12	1.40	1.00
1:B:42:ILE:HG22	1:B:51:LYS:NZ	1.78	0.99
1:A:42:ILE:HG22	1:A:51:LYS:NZ	1.81	0.95
1:B:112:ASP:HB3	1:B:115:LYS:HD3	1.47	0.94
1:B:39:GLU:HG2	1:B:42:ILE:HG12	1.52	0.91
1:B:40:VAL:HG13	1:B:41:ASN:N	1.87	0.89
1:B:58:THR:HG22	1:B:59:GLY:H	1.40	0.87
1:A:39:GLU:HG2	1:A:42:ILE:HG12	1.55	0.87
1:B:42:ILE:HG22	1:B:51:LYS:HZ1	1.37	0.86
1:A:33:VAL:HG21	1:A:43:MSE:HB2	1.58	0.86
1:A:88:LEU:C	1:A:88:LEU:HD23	1.95	0.86
1:A:40:VAL:HG13	1:A:41:ASN:N	1.91	0.85
1:B:76:ASP:HA	1:B:79:LYS:HD2	1.59	0.84
1:A:42:ILE:HG22	1:A:51:LYS:HZ1	1.41	0.84
1:B:33:VAL:HG21	1:B:43:MSE:HB2	1.60	0.83
1:B:27:LYS:HD3	1:B:64:TYR:HB2	1.60	0.83
1:A:76:ASP:HA	1:A:79:LYS:HD2	1.62	0.82
1:A:112:ASP:HB3	1:A:115:LYS:HD3	1.61	0.81
1:A:27:LYS:HD3	1:A:64:TYR:HB2	1.63	0.81
1:B:74:MSE:HE2	1:B:116:ILE:HD12	1.61	0.81
1:A:75:MSE:O	1:A:79:LYS:HG3	1.81	0.80
1:B:75:MSE:O	1:B:79:LYS:HG3	1.83	0.79
1:B:58:THR:HG22	1:B:59:GLY:N	1.98	0.78
1:B:25:HIS:O	1:B:64:TYR:HB3	1.83	0.78
1:B:88:LEU:HD23	1:B:88:LEU:C	2.04	0.77
1:A:13:GLU:OE1	1:A:27:LYS:HE3	1.85	0.77
1:B:40:VAL:CG1	1:B:41:ASN:H	1.90	0.77
1:A:40:VAL:CG1	1:A:41:ASN:H	1.93	0.76
1:A:25:HIS:O	1:A:64:TYR:HB3	1.85	0.76
1:A:92:LEU:HD13	1:A:104:VAL:HG21	1.65	0.75
1:A:102:GLU:HB3	1:A:141:GLU:O	1.86	0.75
1:B:143:LEU:HD13	1:B:144:SER:H	1.53	0.74
1:B:52:THR:HG23	1:B:54:GLY:H	1.53	0.74
1:B:89:GLN:HG3	1:B:105:THR:HG22	1.71	0.72
1:B:42:ILE:HG22	1:B:51:LYS:HZ2	1.55	0.72
1:A:67:THR:HG23	1:A:70:PHE:HB2	1.72	0.71
1:B:45:ARG:HG3	1:B:45:ARG:HH11	1.57	0.70
1:B:102:GLU:HB3	1:B:141:GLU:O	1.91	0.70
1:B:67:THR:HG23	1:B:70:PHE:HB2	1.74	0.69
1:A:89:GLN:HG3	1:A:105:THR:HG22	1.73	0.69
1:A:58:THR:HG22	1:A:59:GLY:N	2.08	0.68
1:A:42:ILE:HG22	1:A:51:LYS:HZ2	1.57	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:45:ARG:NH1	1:B:45:ARG:HG3	2.10	0.67
1:B:86:PHE:CD1	1:B:86:PHE:N	2.63	0.67
1:A:16:LEU:HD21	1:A:24:ALA:HB3	1.77	0.67
1:A:106:LEU:O	1:A:109:VAL:HG12	1.95	0.67
1:B:64:TYR:CE2	1:B:65:LYS:HE3	2.31	0.66
1:B:16:LEU:HD21	1:B:24:ALA:HB3	1.76	0.66
1:B:13:GLU:OE1	1:B:27:LYS:HE3	1.95	0.66
1:A:116:ILE:HA	1:A:119:LEU:HD13	1.78	0.66
1:A:58:THR:HG22	1:A:59:GLY:H	1.59	0.65
1:B:112:ASP:CB	1:B:115:LYS:HD3	2.25	0.64
1:B:86:PHE:HD1	1:B:86:PHE:H	1.44	0.64
1:A:84:PRO:HG2	1:A:86:PHE:CZ	2.33	0.64
1:A:64:TYR:CE2	1:A:65:LYS:HE3	2.33	0.63
1:B:29:PHE:HB2	1:B:61:ALA:HB2	1.80	0.63
1:A:115:LYS:HE3	1:A:131:PRO:HD2	1.80	0.63
1:A:76:ASP:HA	1:A:79:LYS:CD	2.29	0.63
1:A:43:MSE:CG	1:A:47:MSE:HE2	2.28	0.62
1:A:67:THR:HG23	1:A:70:PHE:CB	2.30	0.62
1:A:52:THR:HG23	1:A:54:GLY:H	1.65	0.61
1:A:29:PHE:HB2	1:A:61:ALA:HB2	1.81	0.61
1:B:39:GLU:CG	1:B:42:ILE:HG12	2.29	0.60
1:A:43:MSE:SE	1:A:47:MSE:HE2	2.51	0.60
1:B:92:LEU:HD13	1:B:104:VAL:HG21	1.84	0.59
1:B:106:LEU:O	1:B:109:VAL:HG12	2.02	0.59
1:B:116:ILE:HA	1:B:119:LEU:HD13	1.84	0.59
1:B:52:THR:C	1:B:54:GLY:H	2.06	0.59
1:A:92:LEU:HB2	1:A:104:VAL:CG2	2.33	0.59
1:A:86:PHE:CD1	1:A:86:PHE:N	2.70	0.59
1:B:16:LEU:H	1:B:16:LEU:HD23	1.67	0.59
1:B:9:ILE:HG23	1:B:9:ILE:O	2.01	0.59
1:A:43:MSE:HG2	1:A:47:MSE:HE2	1.84	0.58
1:B:76:ASP:HA	1:B:79:LYS:CD	2.32	0.58
1:A:112:ASP:CB	1:A:115:LYS:HD3	2.32	0.58
1:A:136:ASP:OD1	1:A:137:PHE:N	2.36	0.58
1:A:46:ARG:NH2	1:A:136:ASP:OD2	2.35	0.58
1:A:39:GLU:CG	1:A:42:ILE:HG12	2.30	0.57
1:B:67:THR:HG23	1:B:70:PHE:CB	2.34	0.57
1:A:16:LEU:HD23	1:A:16:LEU:H	1.69	0.57
1:A:88:LEU:C	1:A:88:LEU:CD2	2.69	0.56
1:A:101:THR:O	1:A:143:LEU:HB2	2.05	0.56
1:B:136:ASP:OD1	1:B:137:PHE:N	2.38	0.56
1:B:58:THR:CG2	1:B:59:GLY:H	2.15	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:77:TYR:O	1:A:80:LYS:HG2	2.05	0.56
1:A:73:LEU:HD13	1:A:73:LEU:C	2.26	0.56
1:B:84:PRO:HG2	1:B:86:PHE:CZ	2.40	0.56
1:A:23:MSE:O	1:A:24:ALA:HB2	2.05	0.56
1:A:45:ARG:HG3	1:A:45:ARG:HH11	1.70	0.55
1:B:45:ARG:O	1:B:51:LYS:HD3	2.07	0.55
1:B:17:PHE:N	1:B:17:PHE:CD1	2.74	0.55
1:A:52:THR:C	1:A:54:GLY:H	2.08	0.55
1:A:17:PHE:CD1	1:A:17:PHE:N	2.74	0.55
1:A:89:GLN:HG2	1:A:103:ARG:HD2	1.88	0.55
1:B:101:THR:O	1:B:143:LEU:HB2	2.07	0.55
1:A:26:ILE:HD11	1:A:61:ALA:HB1	1.89	0.55
1:B:111:PHE:CD2	1:B:132:PHE:HB3	2.42	0.54
1:A:74:MSE:CE	1:A:113:SER:HA	2.37	0.54
1:A:74:MSE:HE3	1:A:113:SER:HA	1.89	0.54
1:B:14:GLY:HA3	1:B:90:ALA:HB2	1.89	0.54
1:B:46:ARG:HH22	1:B:136:ASP:CG	2.11	0.54
1:A:45:ARG:NH1	1:A:45:ARG:HG3	2.23	0.54
1:B:77:TYR:O	1:B:80:LYS:HG2	2.07	0.53
1:A:46:ARG:HD3	1:A:51:LYS:HG2	1.89	0.53
1:A:55:ALA:HB1	1:A:137:PHE:HD2	1.73	0.53
1:A:136:ASP:OD1	1:A:137:PHE:O	2.26	0.53
1:A:16:LEU:CD2	1:A:24:ALA:HB3	2.39	0.52
1:A:109:VAL:HA	1:A:134:PHE:HA	1.91	0.52
1:B:143:LEU:HD13	1:B:144:SER:N	2.21	0.52
1:B:43:MSE:O	1:B:44:GLY:C	2.46	0.52
1:B:29:PHE:HB2	1:B:61:ALA:CB	2.39	0.52
1:B:115:LYS:HE3	1:B:131:PRO:HD2	1.91	0.52
1:B:88:LEU:C	1:B:88:LEU:CD2	2.77	0.52
1:A:89:GLN:OE1	1:A:103:ARG:NE	2.43	0.52
1:B:16:LEU:CD2	1:B:24:ALA:HB3	2.40	0.51
1:A:105:THR:C	1:A:106:LEU:HD22	2.31	0.51
1:A:90:ALA:O	1:A:103:ARG:HG3	2.10	0.51
1:B:64:TYR:CD2	1:B:65:LYS:HG2	2.46	0.51
1:A:22:GLU:O	1:A:23:MSE:HG2	2.11	0.51
1:B:58:THR:CG2	1:B:59:GLY:N	2.69	0.51
1:B:138:ASP:OD1	1:B:139:VAL:N	2.43	0.51
1:B:142:LYS:O	1:B:143:LEU:HB2	2.11	0.51
1:A:6:GLN:C	1:A:8:THR:H	2.13	0.51
1:B:55:ALA:HB1	1:B:137:PHE:HD2	1.76	0.51
1:A:88:LEU:HD23	1:A:88:LEU:O	2.11	0.51
1:B:23:MSE:O	1:B:24:ALA:HB2	2.11	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:92:LEU:HB2	1:B:104:VAL:CG2	2.40	0.51
1:B:116:ILE:O	1:B:119:LEU:HD13	2.10	0.50
1:A:29:PHE:HB2	1:A:61:ALA:CB	2.41	0.50
1:A:51:LYS:HG3	1:A:52:THR:O	2.11	0.50
1:B:43:MSE:SE	1:B:47:MSE:HE2	2.62	0.50
1:B:26:ILE:HD11	1:B:61:ALA:HB1	1.93	0.50
1:B:6:GLN:C	1:B:8:THR:H	2.14	0.50
1:B:46:ARG:HD3	1:B:51:LYS:HG2	1.93	0.49
1:A:9:ILE:O	1:A:9:ILE:HG23	2.12	0.49
1:A:43:MSE:O	1:A:44:GLY:C	2.48	0.49
1:B:86:PHE:O	1:B:109:VAL:HG13	2.13	0.49
1:B:109:VAL:HA	1:B:134:PHE:HA	1.94	0.49
1:B:51:LYS:HG3	1:B:52:THR:O	2.13	0.49
1:A:79:LYS:C	1:A:81:GLY:H	2.15	0.48
1:B:79:LYS:C	1:B:81:GLY:H	2.16	0.48
1:B:78:VAL:HB	1:B:113:SER:HB2	1.94	0.48
1:B:73:LEU:HD13	1:B:73:LEU:C	2.33	0.48
1:A:111:PHE:CD2	1:A:132:PHE:HB3	2.49	0.48
1:A:14:GLY:HA3	1:A:90:ALA:HB2	1.96	0.48
1:A:99:ARG:NH2	1:A:102:GLU:OE2	2.47	0.48
1:A:64:TYR:HE2	1:A:65:LYS:HE3	1.77	0.48
1:B:102:GLU:HA	1:B:142:LYS:O	2.14	0.47
1:A:138:ASP:OD1	1:A:139:VAL:N	2.46	0.47
1:A:45:ARG:O	1:A:51:LYS:HD3	2.14	0.47
1:B:15:ARG:HG2	1:B:22:GLU:OE2	2.14	0.47
1:B:64:TYR:HE2	1:B:65:LYS:HE3	1.79	0.47
1:A:18:LEU:HD23	1:A:18:LEU:C	2.35	0.47
1:B:41:ASN:HD21	1:B:44:GLY:H	1.63	0.47
1:A:116:ILE:O	1:A:119:LEU:HD13	2.14	0.47
1:A:28:THR:HG23	1:A:62:THR:OG1	2.14	0.47
1:A:58:THR:CG2	1:A:59:GLY:N	2.78	0.47
1:A:132:PHE:C	1:A:132:PHE:CD1	2.88	0.47
1:B:41:ASN:HD22	1:B:41:ASN:C	2.18	0.47
1:A:46:ARG:HH22	1:A:136:ASP:CG	2.18	0.46
1:B:132:PHE:CD1	1:B:132:PHE:C	2.89	0.46
1:B:10:SER:HB3	1:B:30:GLU:HG3	1.97	0.46
1:A:23:MSE:HE2	1:A:70:PHE:HD1	1.81	0.46
1:A:22:GLU:C	1:A:23:MSE:HG2	2.36	0.46
1:B:74:MSE:CE	1:B:113:SER:HA	2.46	0.46
1:B:43:MSE:CG	1:B:47:MSE:HE2	2.46	0.46
1:B:28:THR:HG23	1:B:62:THR:OG1	2.16	0.46
1:A:115:LYS:CE	1:A:131:PRO:HD2	2.45	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:43:MSE:HE3	1:A:47:MSE:CE	2.46	0.45
1:A:31:ALA:HB1	1:A:134:PHE:CE1	2.51	0.45
1:A:89:GLN:HG2	1:A:103:ARG:CD	2.46	0.45
1:B:52:THR:HG23	1:B:54:GLY:N	2.27	0.45
1:B:143:LEU:O	1:B:144:SER:C	2.55	0.45
1:A:26:ILE:HD11	1:A:61:ALA:CB	2.46	0.45
1:A:88:LEU:HD23	1:A:89:GLN:N	2.31	0.45
1:A:27:LYS:HB3	1:A:28:THR:H	1.64	0.45
1:A:17:PHE:HD1	1:A:17:PHE:N	2.15	0.45
1:B:35:LYS:O	1:B:35:LYS:HG2	2.16	0.45
1:B:105:THR:C	1:B:106:LEU:HD22	2.37	0.45
1:B:88:LEU:HB2	1:B:109:VAL:HG11	1.99	0.45
1:B:89:GLN:HG2	1:B:103:ARG:HD2	1.98	0.45
1:A:140:PRO:HG2	1:A:141:GLU:H	1.81	0.44
1:A:86:PHE:HD1	1:A:86:PHE:H	1.60	0.44
1:B:31:ALA:HB1	1:B:134:PHE:CE1	2.52	0.44
1:A:46:ARG:NH1	1:A:136:ASP:OD2	2.48	0.44
1:A:142:LYS:O	1:A:143:LEU:HD23	2.17	0.44
1:B:130:VAL:HA	1:B:131:PRO:HD3	1.79	0.44
1:B:62:THR:HG22	1:B:128:GLU:C	2.37	0.44
1:A:10:SER:O	1:A:93:ASP:OD1	2.36	0.44
1:B:52:THR:C	1:B:54:GLY:N	2.70	0.43
1:B:42:ILE:CG2	1:B:51:LYS:NZ	2.65	0.43
1:A:112:ASP:HB3	1:A:115:LYS:CD	2.41	0.43
1:A:46:ARG:HB2	1:A:51:LYS:HD3	2.01	0.43
1:B:46:ARG:NH2	1:B:136:ASP:OD1	2.51	0.43
1:B:79:LYS:C	1:B:81:GLY:N	2.72	0.43
1:A:77:TYR:HA	1:A:80:LYS:HE3	2.00	0.43
1:A:86:PHE:O	1:A:109:VAL:HG13	2.19	0.43
1:B:23:MSE:HB3	1:B:70:PHE:CD1	2.54	0.43
1:B:14:GLY:C	1:B:15:ARG:HD2	2.38	0.43
1:B:106:LEU:HA	1:B:106:LEU:HD13	1.87	0.43
1:B:41:ASN:ND2	1:B:44:GLY:H	2.16	0.43
1:A:46:ARG:HD2	1:A:46:ARG:HA	1.83	0.43
1:B:97:SER:C	1:B:99:ARG:N	2.72	0.43
1:A:116:ILE:CA	1:A:119:LEU:HD13	2.48	0.43
1:B:42:ILE:CG2	1:B:51:LYS:HZ1	2.19	0.42
1:B:77:TYR:HA	1:B:80:LYS:HE3	2.01	0.42
1:B:99:ARG:NH2	1:B:102:GLU:OE2	2.53	0.42
1:B:46:ARG:HB2	1:B:51:LYS:HD3	2.00	0.42
1:A:51:LYS:HG3	1:A:52:THR:N	2.33	0.42
1:B:26:ILE:HD11	1:B:61:ALA:CB	2.50	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:36:ASN:ND2	1:B:37:LYS:HE3	2.34	0.42
1:B:17:PHE:N	1:B:17:PHE:HD1	2.16	0.42
1:A:9:ILE:HG23	1:A:31:ALA:HB3	2.02	0.42
1:B:27:LYS:HB3	1:B:28:THR:H	1.62	0.42
1:B:74:MSE:HE2	1:B:113:SER:HA	2.01	0.42
1:B:41:ASN:ND2	1:B:41:ASN:C	2.74	0.41
1:A:47:MSE:HG2	1:A:137:PHE:HZ	1.86	0.41
1:B:53:THR:O	1:B:53:THR:HG22	2.19	0.41
1:A:79:LYS:C	1:A:81:GLY:N	2.74	0.41
1:A:64:TYR:CD2	1:A:65:LYS:HG2	2.55	0.41
1:A:41:ASN:HD22	1:A:41:ASN:C	2.24	0.40
1:B:47:MSE:HG2	1:B:137:PHE:HZ	1.86	0.40
1:A:112:ASP:O	1:A:115:LYS:HB2	2.21	0.40
1:B:26:ILE:HG13	1:B:62:THR:O	2.22	0.40
1:B:89:GLN:HG2	1:B:103:ARG:CD	2.52	0.40
1:B:142:LYS:O	1:B:143:LEU:CB	2.69	0.40
1:A:88:LEU:HB2	1:A:109:VAL:HG11	2.04	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	129/155 (83%)	100 (78%)	22 (17%)	7 (5%)	3	17
1	B	133/155 (86%)	98 (74%)	28 (21%)	7 (5%)	3	18
All	All	262/310 (84%)	198 (76%)	50 (19%)	14 (5%)	3	18

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	40	VAL
1	B	40	VAL
1	A	13	GLU

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Mol	Chain	Res	Type
1	A	24	ALA
1	A	54	GLY
1	B	13	GLU
1	B	19	ASP
1	B	24	ALA
1	B	54	GLY
1	B	144	SER
1	A	27	LYS
1	B	142	LYS
1	A	53	THR
1	A	142	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	114/128 (89%)	95 (83%)	19 (17%)	3	16
1	B	118/128 (92%)	99 (84%)	19 (16%)	3	17
All	All	232/256 (91%)	194 (84%)	38 (16%)	3	16

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	10	SER
1	A	17	PHE
1	A	33	VAL
1	A	41	ASN
1	A	48	THR
1	A	60	THR
1	A	64	TYR
1	A	67	THR
1	A	76	ASP
1	A	77	TYR
1	A	82	SER
1	A	86	PHE
1	A	103	ARG

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Mol	Chain	Res	Type
1	A	109	VAL
1	A	111	PHE
1	A	135	GLU
1	A	137	PHE
1	A	143	LEU
1	B	7	ASN
1	B	15	ARG
1	B	17	PHE
1	B	29	PHE
1	B	33	VAL
1	B	41	ASN
1	B	48	THR
1	B	60	THR
1	B	67	THR
1	B	76	ASP
1	B	77	TYR
1	B	82	SER
1	B	86	PHE
1	B	103	ARG
1	B	109	VAL
1	B	111	PHE
1	B	135	GLU
1	B	137	PHE
1	B	143	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	36	ASN
1	A	41	ASN
1	A	50	HIS
1	A	95	GLN
1	B	7	ASN
1	B	36	ASN
1	B	41	ASN
1	B	50	HIS
1	B	95	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	133/155 (85%)	0.03	0 100 100	10, 33, 69, 70	0
1	B	137/155 (88%)	0.03	0 100 100	10, 32, 69, 70	0
All	All	270/310 (87%)	0.03	0 100 100	10, 34, 69, 70	0

There are no RSRZ outliers to report.

### 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.