



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 02:26 AM GMT

PDB ID : 2F4N  
Title : Crystal structure of protein MJ1651 from Methanococcus jannaschii DSM 2661, Pfam DUF62  
Authors : Rao, K.N.; Swaminathan, S.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)  
Deposited on : 2005-11-23  
Resolution : 2.50 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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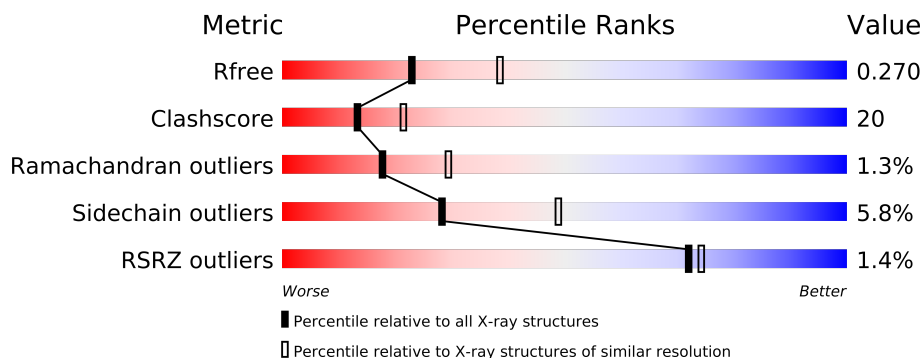
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.50 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	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (2.50-2.50)

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	273	
1	B	273	
1	C	273	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5570 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 Hypothetical protein MJ1651.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	Se	0	0	0
			1835	1189	291	349	2	4			
1	B	239	Total	C	N	O	S	Se	0	0	0
			1864	1211	303	344	2	4			
1	C	239	Total	C	N	O	S	Se	0	0	0
			1840	1194	293	348	2	3			

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	CLONING ARTIFACT	UNP Q59045
A	2	SER	-	CLONING ARTIFACT	UNP Q59045
A	3	LEU	-	CLONING ARTIFACT	UNP Q59045
A	7	MSE	MET	MODIFIED RESIDUE	UNP Q59045
A	31	MSE	MET	MODIFIED RESIDUE	UNP Q59045
A	165	MSE	MET	MODIFIED RESIDUE	UNP Q59045
A	202	MSE	MET	MODIFIED RESIDUE	UNP Q59045
A	246	MSE	MET	MODIFIED RESIDUE	UNP Q59045
A	264	GLU	-	CLONING ARTIFACT	UNP Q59045
A	265	GLY	-	CLONING ARTIFACT	UNP Q59045
A	266	GLY	-	CLONING ARTIFACT	UNP Q59045
A	267	SER	-	CLONING ARTIFACT	UNP Q59045
A	268	HIS	-	EXPRESSION TAG	UNP Q59045
A	269	HIS	-	EXPRESSION TAG	UNP Q59045
A	270	HIS	-	EXPRESSION TAG	UNP Q59045
A	271	HIS	-	EXPRESSION TAG	UNP Q59045
A	272	HIS	-	EXPRESSION TAG	UNP Q59045
A	273	HIS	-	EXPRESSION TAG	UNP Q59045
B	1	MSE	-	CLONING ARTIFACT	UNP Q59045
B	2	SER	-	CLONING ARTIFACT	UNP Q59045
B	3	LEU	-	CLONING ARTIFACT	UNP Q59045
B	7	MSE	MET	MODIFIED RESIDUE	UNP Q59045
B	31	MSE	MET	MODIFIED RESIDUE	UNP Q59045

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Chain	Residue	Modelled	Actual	Comment	Reference
B	165	MSE	MET	MODIFIED RESIDUE	UNP Q59045
B	202	MSE	MET	MODIFIED RESIDUE	UNP Q59045
B	246	MSE	MET	MODIFIED RESIDUE	UNP Q59045
B	264	GLU	-	CLONING ARTIFACT	UNP Q59045
B	265	GLY	-	CLONING ARTIFACT	UNP Q59045
B	266	GLY	-	CLONING ARTIFACT	UNP Q59045
B	267	SER	-	CLONING ARTIFACT	UNP Q59045
B	268	HIS	-	EXPRESSION TAG	UNP Q59045
B	269	HIS	-	EXPRESSION TAG	UNP Q59045
B	270	HIS	-	EXPRESSION TAG	UNP Q59045
B	271	HIS	-	EXPRESSION TAG	UNP Q59045
B	272	HIS	-	EXPRESSION TAG	UNP Q59045
B	273	HIS	-	EXPRESSION TAG	UNP Q59045
C	1	MSE	-	CLONING ARTIFACT	UNP Q59045
C	2	SER	-	CLONING ARTIFACT	UNP Q59045
C	3	LEU	-	CLONING ARTIFACT	UNP Q59045
C	7	MSE	MET	MODIFIED RESIDUE	UNP Q59045
C	31	MSE	MET	MODIFIED RESIDUE	UNP Q59045
C	165	MSE	MET	MODIFIED RESIDUE	UNP Q59045
C	202	MSE	MET	MODIFIED RESIDUE	UNP Q59045
C	246	MSE	MET	MODIFIED RESIDUE	UNP Q59045
C	264	GLU	-	CLONING ARTIFACT	UNP Q59045
C	265	GLY	-	CLONING ARTIFACT	UNP Q59045
C	266	GLY	-	CLONING ARTIFACT	UNP Q59045
C	267	SER	-	CLONING ARTIFACT	UNP Q59045
C	268	HIS	-	EXPRESSION TAG	UNP Q59045
C	269	HIS	-	EXPRESSION TAG	UNP Q59045
C	270	HIS	-	EXPRESSION TAG	UNP Q59045
C	271	HIS	-	EXPRESSION TAG	UNP Q59045
C	272	HIS	-	EXPRESSION TAG	UNP Q59045
C	273	HIS	-	EXPRESSION TAG	UNP Q59045

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	12	Total O 12 12	0	0
2	B	7	Total O 7 7	0	0
2	C	12	Total O 12 12	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.

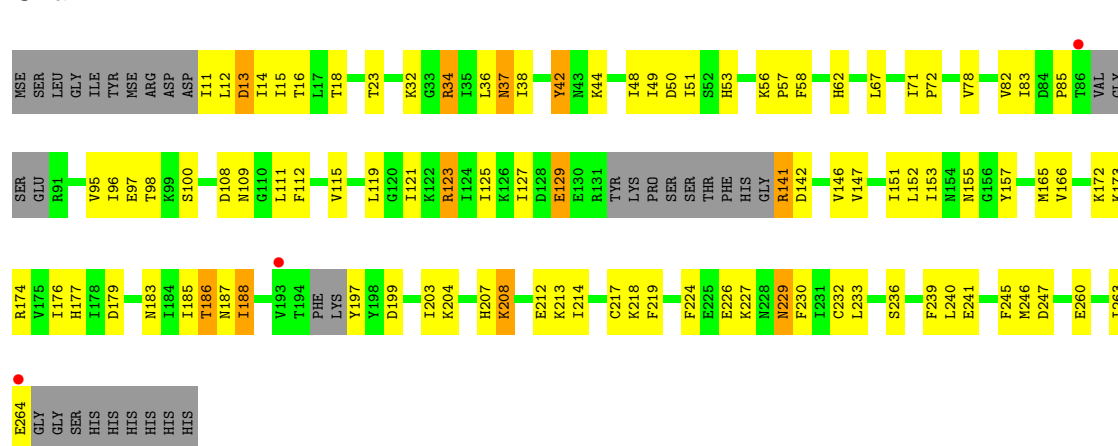
#### • Molecule 1: Hypothetical protein MJ1651

Chain A:



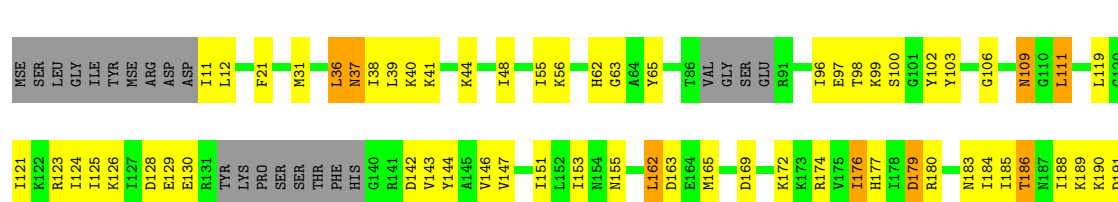
#### • Molecule 1: Hypothetical protein MJ1651

Chain B:



#### • Molecule 1: Hypothetical protein MJ1651

Chain C:



GLU	VAL	THR	F195	1201	1205	N209	G210	1211	K218	F219	S222	E225	E226	K227	C232	L233	1234	E241	1242	1261	F262	1263	E264	GLY	GLY	SER	HIS	HIS	HIS	HIS	HIS	HIS
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## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	126.84Å 110.08Å 75.82Å 90.00° 115.39° 90.00°	Depositor
Resolution (Å)	39.86 – 2.50 39.86 – 2.36	Depositor EDS
% Data completeness (in resolution range)	95.4 (39.86-2.50) 95.5 (39.86-2.36)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.20 (at 2.37Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.237 , 0.265 0.239 , 0.270	Depositor DCC
$R_{free}$ test set	941 reflections (3.03%)	DCC
Wilson B-factor (Å <sup>2</sup> )	40.8	Xtriage
Anisotropy	0.491	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 39.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 72852 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5570	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 5.12% 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	0.39	0/1859	0.63	0/2510
1	B	0.38	0/1890	0.63	0/2547
1	C	0.40	0/1866	0.67	1/2522 (0.0%)
All	All	0.39	0/5615	0.64	1/7579 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	176	ILE	CB-CA-C	-6.19	99.22	111.60

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	1835	0	1797	61	0
1	B	1864	0	1864	94	0
1	C	1840	0	1793	72	0
2	A	12	0	0	0	0
2	B	7	0	0	1	0
2	C	12	0	0	1	0
All	All	5570	0	5454	224	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 20.

All (224) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:141:ARG:HG3	1:B:142:ASP:H	1.22	1.02
1:A:229:ASN:N	1:A:229:ASN:HD22	1.56	1.01
1:B:11:ILE:HA	1:B:155:ASN:HD21	1.26	1.01
1:A:229:ASN:H	1:A:229:ASN:HD22	1.03	0.94
1:C:12:LEU:H	1:C:155:ASN:HD21	0.99	0.91
1:C:12:LEU:H	1:C:155:ASN:ND2	1.70	0.89
1:B:229:ASN:ND2	1:B:230:PHE:H	1.74	0.86
1:A:176:ILE:HG22	1:A:185:ILE:O	1.76	0.85
1:B:34:ARG:HA	1:B:34:ARG:CZ	2.07	0.84
1:C:12:LEU:N	1:C:155:ASN:HD21	1.76	0.84
1:A:202:MSE:HE2	1:A:216:LYS:HE2	1.60	0.83
1:A:194:THR:HG22	1:A:195:PHE:H	1.44	0.82
1:B:141:ARG:HG3	1:B:142:ASP:N	1.94	0.82
1:A:229:ASN:H	1:A:229:ASN:ND2	1.79	0.79
1:C:109:ASN:HD21	1:C:111:LEU:HB2	1.49	0.77
1:B:229:ASN:HD22	1:B:230:PHE:H	1.32	0.76
1:A:229:ASN:N	1:A:229:ASN:ND2	2.30	0.76
1:B:204:LYS:HG3	1:B:214:ILE:HG12	1.71	0.73
1:C:186:THR:HG23	1:C:188:ILE:H	1.51	0.73
1:B:98:THR:HG21	1:B:119:LEU:HB3	1.71	0.73
1:B:208:LYS:H	1:B:208:LYS:HD3	1.52	0.72
1:B:172:LYS:HE2	1:B:260:GLU:HG2	1.72	0.71
1:B:11:ILE:HA	1:B:155:ASN:ND2	2.04	0.71
1:C:56:LYS:HG3	1:C:62:HIS:ND1	2.06	0.70
1:C:11:ILE:HD11	1:C:44:LYS:NZ	2.06	0.70
1:B:204:LYS:HE2	1:B:212:GLU:CD	2.10	0.70
1:B:13:ASP:HB3	1:B:14:ILE:HD12	1.73	0.69
1:B:125:ILE:HD12	1:B:125:ILE:N	2.07	0.69
1:B:176:ILE:HG22	1:B:185:ILE:O	1.91	0.69
1:B:121:ILE:HD13	1:B:165:MSE:HE3	1.75	0.69
1:C:162:LEU:HD23	1:C:163:ASP:H	1.60	0.67
1:A:109:ASN:C	1:A:109:ASN:HD22	1.97	0.67
1:B:218:LYS:HD2	1:B:226:GLU:OE1	1.96	0.65
1:C:11:ILE:HD11	1:C:44:LYS:HZ3	1.62	0.65
1:C:11:ILE:HG13	1:C:155:ASN:ND2	2.12	0.65
1:B:204:LYS:HE2	1:B:212:GLU:OE2	1.98	0.63
1:B:229:ASN:HD22	1:B:230:PHE:N	1.95	0.63
1:B:176:ILE:CG2	1:B:177:HIS:N	2.61	0.63
1:C:232:CYS:HA	1:C:241:GLU:O	1.99	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:176:ILE:HG21	1:B:185:ILE:HG22	1.81	0.63
1:A:176:ILE:HG21	1:A:185:ILE:HG22	1.80	0.62
1:C:97:GLU:HG3	1:C:103:TYR:CZ	2.34	0.62
1:A:121:ILE:HD13	1:A:165:MSE:HE3	1.81	0.62
1:A:96:ILE:HD13	1:A:165:MSE:HE1	1.82	0.62
1:A:109:ASN:HD21	1:A:111:LEU:HB2	1.65	0.62
1:C:123:ARG:HG3	1:C:125:ILE:HD11	1.82	0.62
1:A:177:HIS:HA	1:A:257:TYR:CE2	2.35	0.61
1:C:176:ILE:HG22	1:C:177:HIS:HB2	1.83	0.61
1:B:32:LYS:O	1:B:36:LEU:HB2	1.99	0.61
1:A:129:GLU:O	1:A:131:ARG:N	2.33	0.61
1:B:38:ILE:HD11	1:B:146:VAL:HG22	1.83	0.60
1:B:98:THR:HG23	1:B:119:LEU:O	2.01	0.60
1:C:233:LEU:HD23	1:C:233:LEU:N	2.17	0.60
1:C:125:ILE:HD12	1:C:125:ILE:N	2.16	0.60
1:B:179:ASP:OD2	1:B:183:ASN:HB2	2.01	0.59
1:C:65:TYR:CD2	1:C:176:ILE:HG21	2.38	0.59
1:A:128:ASP:OD1	1:A:129:GLU:N	2.30	0.59
1:C:179:ASP:HB2	1:C:183:ASN:HB2	1.85	0.58
1:C:98:THR:HB	1:C:102:TYR:H	1.69	0.58
1:C:21:PHE:HE2	1:C:31:MSE:HE1	1.68	0.58
1:B:174:ARG:H	1:B:187:ASN:HD22	1.51	0.58
1:C:98:THR:CG2	1:C:119:LEU:HB3	2.34	0.58
1:B:82:VAL:O	1:B:83:ILE:HG13	2.04	0.58
1:A:232:CYS:HA	1:A:241:GLU:O	2.04	0.57
1:B:186:THR:HG22	1:B:240:LEU:O	2.03	0.57
1:B:174:ARG:HG2	1:B:187:ASN:HD21	1.68	0.57
1:A:113:THR:O	1:A:117:GLU:HG3	2.03	0.57
1:A:131:ARG:NH2	1:A:150:GLU:OE1	2.37	0.57
1:B:173:LYS:HA	1:B:187:ASN:HD22	1.70	0.57
1:C:36:LEU:HD13	1:C:48:ILE:HD11	1.84	0.57
1:C:109:ASN:HD22	1:C:109:ASN:C	2.06	0.57
1:C:176:ILE:HG22	1:C:177:HIS:CB	2.35	0.57
1:B:233:LEU:HD23	1:B:233:LEU:H	1.70	0.57
1:B:34:ARG:NE	1:B:34:ARG:HA	2.20	0.56
1:B:96:ILE:HD13	1:B:165:MSE:HE1	1.86	0.56
1:B:174:ARG:HG2	1:B:187:ASN:ND2	2.20	0.56
1:C:38:ILE:HG22	1:C:153:ILE:HD11	1.86	0.56
1:C:218:LYS:HD3	1:C:226:GLU:CB	2.36	0.56
1:C:98:THR:HG21	1:C:119:LEU:HB3	1.86	0.56
1:B:188:ILE:HD13	1:B:263:ILE:HD12	1.88	0.56
1:A:75:PRO:HB3	1:B:37:ASN:ND2	2.21	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:50:ASP:OD1	1:B:53:HIS:NE2	2.36	0.55
1:C:142:ASP:O	1:C:146:VAL:HG23	2.06	0.55
1:C:147:VAL:O	1:C:151:ILE:HG13	2.07	0.55
1:B:12:LEU:H	1:B:155:ASN:ND2	2.05	0.55
1:C:176:ILE:HG23	2:C:276:HOH:O	2.07	0.55
1:C:189:LYS:C	1:C:191:ASP:H	2.10	0.55
1:A:176:ILE:HD12	1:A:187:ASN:HA	1.87	0.55
1:C:205:ILE:HG12	1:C:261:ILE:CD1	2.37	0.55
1:C:201:ILE:HD13	1:C:219:PHE:HB2	1.89	0.55
1:B:233:LEU:HD23	1:B:233:LEU:N	2.22	0.54
1:A:126:LYS:HB2	1:A:162:LEU:HD23	1.90	0.54
1:B:174:ARG:H	1:B:187:ASN:ND2	2.05	0.54
1:A:125:ILE:N	1:A:125:ILE:HD12	2.22	0.54
1:C:186:THR:HG23	1:C:188:ILE:N	2.22	0.54
1:C:172:LYS:O	1:C:174:ARG:HG2	2.08	0.54
1:C:185:ILE:N	1:C:185:ILE:HD12	2.24	0.53
1:A:177:HIS:HD2	1:A:185:ILE:HD12	1.74	0.53
1:B:176:ILE:HG23	1:B:177:HIS:N	2.23	0.53
1:C:222:SER:OG	1:C:225:GLU:HG3	2.09	0.52
1:B:207:HIS:HE1	1:B:213:LYS:HG3	1.74	0.52
1:C:109:ASN:ND2	1:C:111:LEU:H	2.07	0.52
1:A:246:MSE:HA	1:A:246:MSE:HE2	1.91	0.52
1:C:37:ASN:C	1:C:37:ASN:HD22	2.13	0.52
1:A:176:ILE:CG2	1:A:185:ILE:HG22	2.40	0.51
1:C:234:ILE:N	1:C:234:ILE:HD12	2.25	0.51
1:B:172:LYS:HE2	1:B:260:GLU:CG	2.39	0.51
1:B:174:ARG:N	1:B:187:ASN:HD22	2.08	0.51
1:B:232:CYS:HA	1:B:241:GLU:O	2.09	0.51
1:A:194:THR:HG22	1:A:195:PHE:N	2.20	0.51
1:C:36:LEU:HD13	1:C:48:ILE:CD1	2.41	0.50
1:B:263:ILE:HG22	1:B:264:GLU:N	2.25	0.50
1:C:41:LYS:HG2	1:C:41:LYS:O	2.11	0.50
1:A:176:ILE:HG22	1:A:185:ILE:C	2.32	0.50
1:B:82:VAL:C	1:B:83:ILE:HG13	2.32	0.50
1:C:128:ASP:O	1:C:130:GLU:N	2.44	0.49
1:C:96:ILE:CD1	1:C:165:MSE:HE1	2.42	0.49
1:C:169:ASP:OD2	1:C:172:LYS:HB2	2.12	0.49
1:C:109:ASN:HD22	1:C:111:LEU:H	1.60	0.49
1:C:21:PHE:HE2	1:C:31:MSE:CE	2.26	0.49
1:B:56:LYS:HG3	1:B:62:HIS:ND1	2.28	0.49
1:B:147:VAL:O	1:B:151:ILE:HG13	2.13	0.49
1:C:96:ILE:HD13	1:C:165:MSE:HE1	1.95	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:97:GLU:OE1	1:A:123:ARG:HD3	2.13	0.49
1:B:176:ILE:HG22	1:B:185:ILE:C	2.33	0.49
1:C:162:LEU:HD23	1:C:163:ASP:N	2.25	0.48
1:B:97:GLU:CB	1:B:123:ARG:HG3	2.43	0.48
1:B:218:LYS:HD2	1:B:226:GLU:CD	2.34	0.48
1:C:12:LEU:HD12	1:C:155:ASN:HA	1.96	0.48
1:A:109:ASN:HD22	1:A:111:LEU:H	1.60	0.48
1:B:58:PHE:HB3	1:B:85:PRO:HB3	1.96	0.48
1:B:98:THR:CG2	1:B:119:LEU:HB3	2.41	0.48
1:A:129:GLU:O	1:A:130:GLU:C	2.52	0.48
1:B:203:ILE:HD12	1:B:217:CYS:SG	2.54	0.48
1:B:38:ILE:HG22	1:B:153:ILE:HD11	1.96	0.48
1:C:124:ILE:C	1:C:125:ILE:HD12	2.35	0.48
1:B:112:PHE:HD1	1:B:115:VAL:HG11	1.80	0.47
1:B:13:ASP:HB2	1:C:40:LYS:NZ	2.29	0.47
1:A:36:LEU:HD13	1:A:48:ILE:HD11	1.97	0.47
1:B:11:ILE:CA	1:B:155:ASN:HD21	2.13	0.47
1:A:194:THR:O	1:A:195:PHE:HB3	2.14	0.47
1:B:246:MSE:HA	1:B:246:MSE:HE2	1.95	0.47
1:B:18:THR:OG1	1:B:51:ILE:HB	2.14	0.47
1:C:143:VAL:O	1:C:147:VAL:HG23	2.14	0.47
1:C:121:ILE:HD13	1:C:165:MSE:HE3	1.97	0.47
1:A:179:ASP:CB	1:A:183:ASN:HB2	2.44	0.46
1:B:179:ASP:HB2	1:B:183:ASN:H	1.80	0.46
1:B:224:PHE:O	1:B:227:LYS:HE3	2.15	0.46
1:A:195:PHE:CD1	1:A:195:PHE:O	2.69	0.46
1:A:147:VAL:O	1:A:151:ILE:HG13	2.16	0.46
1:C:55:ILE:HD12	1:C:63:GLY:HA2	1.96	0.46
1:A:234:ILE:N	1:A:234:ILE:HD12	2.30	0.46
1:C:186:THR:HG21	1:C:188:ILE:HD12	1.97	0.46
1:A:224:PHE:O	1:A:227:LYS:HG3	2.15	0.46
1:B:204:LYS:HE3	1:B:214:ILE:HD11	1.98	0.46
1:B:151:ILE:HG12	1:B:157:TYR:HB2	1.97	0.45
1:B:173:LYS:HD2	1:B:188:ILE:HG13	1.97	0.45
1:B:219:PHE:HA	1:B:232:CYS:O	2.17	0.45
1:A:173:LYS:HD3	1:A:173:LYS:HA	1.74	0.45
1:A:126:LYS:HD3	1:A:160:GLU:CD	2.37	0.45
1:B:177:HIS:HD2	1:B:185:ILE:HD12	1.82	0.45
1:C:98:THR:HG22	1:C:99:LYS:N	2.30	0.45
1:A:20:ASP:O	1:C:180:ARG:NH1	2.50	0.45
1:B:42:TYR:CD1	1:B:42:TYR:N	2.85	0.45
1:C:98:THR:HG22	1:C:100:SER:H	1.82	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:21:PHE:CE2	1:C:31:MSE:HE1	2.51	0.45
1:A:36:LEU:HD13	1:A:48:ILE:CD1	2.47	0.45
1:A:109:ASN:ND2	1:A:111:LEU:H	2.15	0.44
1:A:109:ASN:ND2	1:A:109:ASN:C	2.66	0.44
1:C:205:ILE:HG12	1:C:261:ILE:HD13	1.99	0.44
1:B:71:ILE:HD12	1:B:115:VAL:HG21	2.00	0.44
1:C:109:ASN:ND2	1:C:109:ASN:C	2.71	0.44
1:A:40:LYS:C	1:A:42:TYR:H	2.20	0.44
1:B:44:LYS:HE2	1:B:152:LEU:O	2.18	0.44
1:C:162:LEU:CD2	1:C:163:ASP:N	2.81	0.44
1:A:15:ILE:HA	1:A:78:VAL:O	2.16	0.44
1:B:263:ILE:CG2	1:B:264:GLU:N	2.80	0.44
1:A:176:ILE:HG23	1:A:177:HIS:HB3	1.99	0.44
1:B:129:GLU:CD	1:B:129:GLU:O	2.56	0.44
1:B:176:ILE:CG2	1:B:185:ILE:HG22	2.46	0.43
1:A:182:GLY:HA3	1:A:248:ASN:OD1	2.19	0.43
1:B:187:ASN:O	1:B:239:PHE:HD2	2.02	0.43
1:B:245:PHE:CE2	1:B:246:MSE:HG2	2.53	0.43
1:C:125:ILE:HG22	1:C:126:LYS:O	2.19	0.43
1:A:47:LYS:HE3	2:B:279:HOH:O	2.18	0.43
1:A:150:GLU:OE1	1:A:157:TYR:HE1	2.01	0.43
1:B:36:LEU:HD13	1:B:48:ILE:CD1	2.49	0.43
1:B:229:ASN:ND2	1:B:230:PHE:N	2.51	0.43
1:A:38:ILE:HG22	1:A:153:ILE:HD11	2.01	0.43
1:C:109:ASN:ND2	1:C:111:LEU:HB2	2.25	0.42
1:A:31:MSE:HE2	1:A:141:ARG:O	2.19	0.42
1:C:201:ILE:HG23	1:C:263:ILE:HG23	2.01	0.42
1:A:16:THR:HG22	1:A:49:ILE:HB	2.00	0.42
1:C:189:LYS:O	1:C:191:ASP:N	2.50	0.42
1:A:176:ILE:CG2	1:A:177:HIS:N	2.82	0.42
1:B:125:ILE:CD1	1:B:125:ILE:N	2.77	0.42
1:B:186:THR:HG23	1:B:188:ILE:N	2.35	0.42
1:B:95:VAL:HG23	1:B:127:ILE:CG1	2.49	0.42
1:B:15:ILE:HA	1:B:78:VAL:O	2.20	0.42
1:B:98:THR:HG22	1:B:100:SER:H	1.84	0.42
1:B:176:ILE:HG22	1:B:177:HIS:H	1.85	0.42
1:C:209:ASN:OD1	1:C:211:ILE:N	2.41	0.42
1:A:186:THR:O	1:A:239:PHE:HB3	2.20	0.42
1:A:34:ARG:HD3	1:A:146:VAL:HG22	2.02	0.42
1:A:179:ASP:HB3	1:A:183:ASN:H	1.85	0.41
1:C:184:ILE:HB	1:C:242:ILE:HB	2.00	0.41
1:B:34:ARG:HH21	1:B:37:ASN:CB	2.33	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:186:THR:HG21	1:B:240:LEU:HB3	2.03	0.41
1:C:106:GLY:HA2	1:C:144:TYR:CE1	2.55	0.41
1:B:173:LYS:HA	1:B:173:LYS:HD3	1.80	0.41
1:A:131:ARG:N	1:A:131:ARG:HD2	2.36	0.41
1:B:240:LEU:CD2	1:B:263:ILE:HD11	2.51	0.41
1:C:176:ILE:HG22	1:C:177:HIS:N	2.34	0.41
1:B:67:LEU:HD22	1:B:71:ILE:HD11	2.02	0.41
1:A:12:LEU:HG	1:A:155:ASN:ND2	2.35	0.41
1:C:125:ILE:CD1	1:C:125:ILE:N	2.83	0.41
1:A:182:GLY:HA3	1:A:248:ASN:CG	2.42	0.41
1:B:16:THR:HA	1:B:49:ILE:O	2.21	0.41
1:A:129:GLU:C	1:A:131:ARG:N	2.74	0.40
1:A:256:ASP:CG	1:A:257:TYR:H	2.24	0.40
1:B:58:PHE:CD1	1:B:85:PRO:HA	2.57	0.40
1:B:108:ASP:OD2	1:B:166:VAL:HG23	2.21	0.40
1:B:71:ILE:HB	1:B:72:PRO:HD3	2.03	0.40
1:B:42:TYR:HD1	1:B:42:TYR:N	2.20	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	227/273 (83%)	208 (92%)	16 (7%)	3 (1%)	18	29
1	B	231/273 (85%)	209 (90%)	19 (8%)	3 (1%)	18	29
1	C	231/273 (85%)	214 (93%)	14 (6%)	3 (1%)	18	29
All	All	689/819 (84%)	631 (92%)	49 (7%)	9 (1%)	18	29

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	130	GLU
1	C	129	GLU

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Mol	Chain	Res	Type
1	C	190	LYS
1	A	195	PHE
1	C	227	LYS
1	A	120	GLY
1	B	236	SER
1	B	188	ILE
1	B	57	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	194/238 (82%)	184 (95%)	10 (5%)	32	55
1	B	197/238 (83%)	181 (92%)	16 (8%)	17	30
1	C	192/238 (81%)	184 (96%)	8 (4%)	40	66
All	All	583/714 (82%)	549 (94%)	34 (6%)	28	49

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ASP
1	A	23	THR
1	A	36	LEU
1	A	109	ASN
1	A	111	LEU
1	A	155	ASN
1	A	158	ASP
1	A	162	LEU
1	A	229	ASN
1	A	233	LEU
1	B	13	ASP
1	B	23	THR
1	B	34	ARG
1	B	37	ASN
1	B	42	TYR
1	B	109	ASN

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Mol	Chain	Res	Type
1	B	111	LEU
1	B	123	ARG
1	B	129	GLU
1	B	141	ARG
1	B	186	THR
1	B	197	TYR
1	B	199	ASP
1	B	208	LYS
1	B	229	ASN
1	B	247	ASP
1	C	36	LEU
1	C	37	ASN
1	C	39	LEU
1	C	109	ASN
1	C	111	LEU
1	C	162	LEU
1	C	179	ASP
1	C	186	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	37	ASN
1	A	109	ASN
1	A	154	ASN
1	A	155	ASN
1	A	229	ASN
1	B	43	ASN
1	B	109	ASN
1	B	154	ASN
1	B	155	ASN
1	B	187	ASN
1	B	229	ASN
1	B	254	ASN
1	C	37	ASN
1	C	109	ASN
1	C	154	ASN
1	C	155	ASN
1	C	254	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	237/273 (86%)	0.09	7 (2%) 48 50	27, 49, 71, 86	0
1	B	239/273 (87%)	0.06	3 (1%) 74 76	29, 52, 71, 82	0
1	C	239/273 (87%)	-0.09	0 100 100	19, 45, 66, 76	0
All	All	715/819 (87%)	0.02	10 (1%) 72 74	19, 49, 69, 86	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	9	ASP	4.7
1	A	224	PHE	4.1
1	A	10	ASP	2.7
1	B	193	VAL	2.6
1	B	86	THR	2.4
1	A	246	MSE	2.3
1	A	176	ILE	2.3
1	B	264	GLU	2.3
1	A	228	ASN	2.3
1	A	163	ASP	2.3

### 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.