



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:06 AM UTC

PDB ID : 1TK9 / pdb\_00001tk9  
Title : Crystal Structure of Phosphoheptose isomerase 1  
Authors : Rajashankar, K.R.; Solorzano, V.; Kniewel, R.; Lima, C.D.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)  
Deposited on : 2004-06-08  
Resolution : 2.10 Å(reported)

This is a Full wwPDB X-ray Structure 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

---

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

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

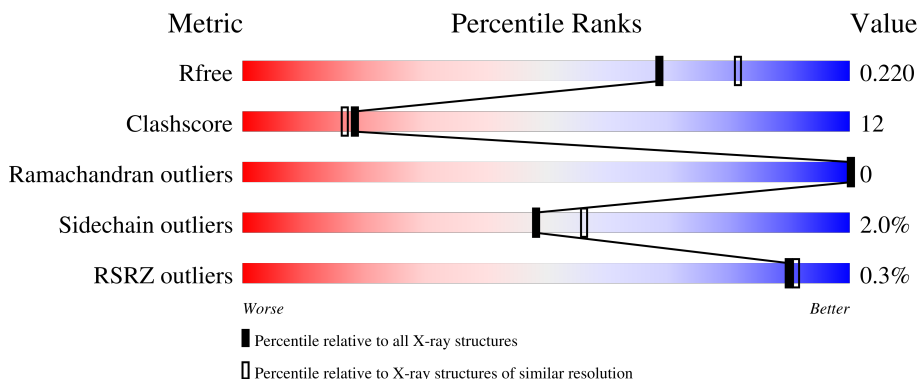
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*





The reported resolution of this entry is 2.10 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))
$R_{free}$	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	188	 79% 20% .
1	B	188	 74% 26%
1	C	188	 74% 25% .
1	D	188	 74% 24% .

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6367 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Phosphoheptose isomerase 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	188	1417	890	244	272	6	5	0	0	0
1	B	188	1417	890	244	272	6	5	0	0	0
1	C	188	1417	890	244	272	6	5	0	0	0
1	D	188	1417	890	244	272	6	5	0	0	0

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	cloning artifact	UNP Q9PNE6
A	2	SER	-	cloning artifact	UNP Q9PNE6
A	3	LEU	-	cloning artifact	UNP Q9PNE6
A	138	MSE	MET	modified residue	UNP Q9PNE6
A	150	MSE	MET	modified residue	UNP Q9PNE6
A	151	MSE	MET	modified residue	UNP Q9PNE6
A	172	MSE	MET	modified residue	UNP Q9PNE6
B	1	MSE	-	cloning artifact	UNP Q9PNE6
B	2	SER	-	cloning artifact	UNP Q9PNE6
B	3	LEU	-	cloning artifact	UNP Q9PNE6
B	138	MSE	MET	modified residue	UNP Q9PNE6
B	150	MSE	MET	modified residue	UNP Q9PNE6
B	151	MSE	MET	modified residue	UNP Q9PNE6
B	172	MSE	MET	modified residue	UNP Q9PNE6
C	1	MSE	-	cloning artifact	UNP Q9PNE6
C	2	SER	-	cloning artifact	UNP Q9PNE6
C	3	LEU	-	cloning artifact	UNP Q9PNE6
C	138	MSE	MET	modified residue	UNP Q9PNE6
C	150	MSE	MET	modified residue	UNP Q9PNE6
C	151	MSE	MET	modified residue	UNP Q9PNE6
C	172	MSE	MET	modified residue	UNP Q9PNE6

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MSE	-	cloning artifact	UNP Q9PNE6
D	2	SER	-	cloning artifact	UNP Q9PNE6
D	3	LEU	-	cloning artifact	UNP Q9PNE6
D	138	MSE	MET	modified residue	UNP Q9PNE6
D	150	MSE	MET	modified residue	UNP Q9PNE6
D	151	MSE	MET	modified residue	UNP Q9PNE6
D	172	MSE	MET	modified residue	UNP Q9PNE6

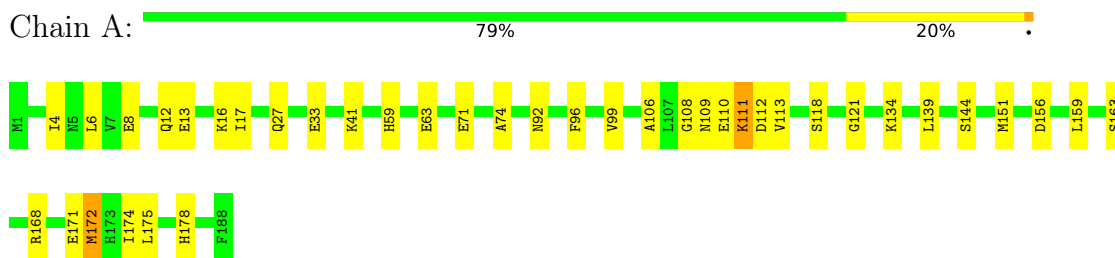
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	186	Total O 186 186	0	0
2	B	183	Total O 183 183	0	0
2	C	169	Total O 169 169	0	0
2	D	161	Total O 161 161	0	0

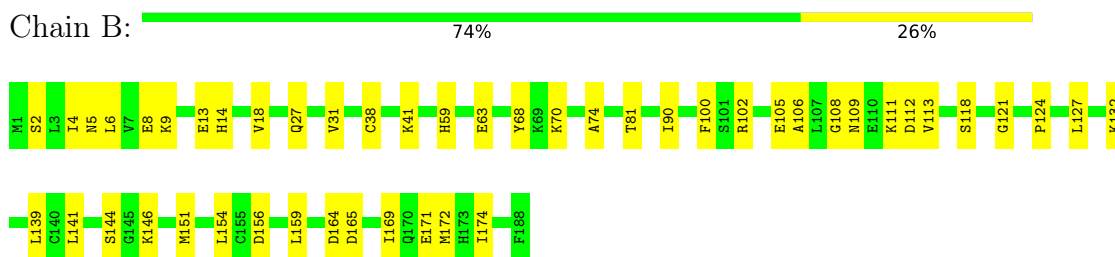
### 3 Residue-property plots [i](#)

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

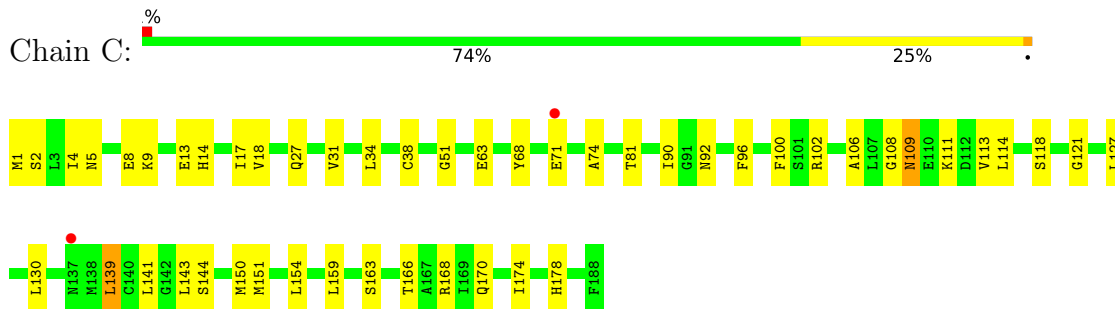
- Molecule 1: Phosphoheptose isomerase 1



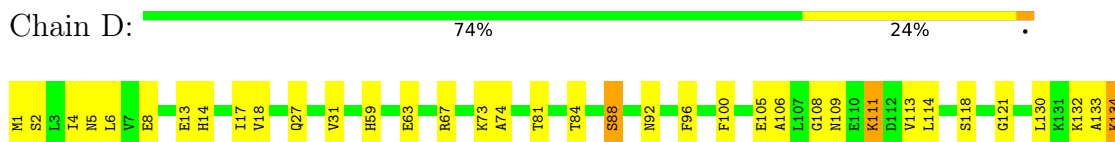
- Molecule 1: Phosphoheptose isomerase 1



- Molecule 1: Phosphoheptose isomerase 1



- Molecule 1: Phosphoheptose isomerase 1





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.66Å 123.29Å 127.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.78 – 2.10 19.78 – 2.10	Depositor EDS
% Data completeness (in resolution range)	94.8 (19.78-2.10) 94.8 (19.78-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.94 (at 2.10Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.182 , 0.211 0.194 , 0.220	Depositor DCC
$R_{free}$ test set	4553 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.1	Xtrriage
Anisotropy	0.083	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 53.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.008 for -h,l,k	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6367	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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	0.42	0/1428	0.85	3/1906 (0.2%)
1	B	0.40	0/1428	0.86	3/1906 (0.2%)
1	C	0.38	0/1428	0.78	1/1906 (0.1%)
1	D	0.40	0/1428	0.82	3/1906 (0.2%)
All	All	0.40	0/5712	0.83	10/7624 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	108	GLY	N-CA-C	7.01	122.18	112.37
1	C	108	GLY	N-CA-C	6.52	120.55	112.14
1	B	108	GLY	N-CA-C	6.50	120.52	112.14
1	B	112	ASP	N-CA-C	5.93	119.19	110.59
1	D	156	ASP	N-CA-C	-5.85	104.86	112.23
1	A	156	ASP	N-CA-C	-5.67	105.24	111.82
1	D	108	GLY	N-CA-C	5.66	120.63	112.82
1	B	156	ASP	N-CA-C	-5.17	105.83	111.82
1	A	112	ASP	N-CA-C	5.15	118.08	110.46
1	D	143	LEU	N-CA-C	-5.06	100.17	108.41

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1417	0	1460	34	0
1	B	1417	0	1460	39	0
1	C	1417	0	1460	39	0
1	D	1417	0	1460	34	0
2	A	186	0	0	8	0
2	B	183	0	0	3	0
2	C	169	0	0	5	0
2	D	161	0	0	1	0
All	All	6367	0	5840	138	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (138) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:367:HOH:O	1:C:90:ILE:HG12	1.66	0.93
1:B:4:ILE:H	1:B:4:ILE:HD12	1.42	0.84
1:C:90:ILE:HD12	2:C:254:HOH:O	1.79	0.81
1:A:172:MSE:HE3	1:A:172:MSE:HA	1.69	0.75
1:B:90:ILE:HD12	2:B:264:HOH:O	1.87	0.74
1:D:134:LYS:HE3	1:D:134:LYS:HA	1.69	0.74
1:A:63:GLU:HG3	2:A:374:HOH:O	1.89	0.73
1:A:109:ASN:HD22	1:A:111:LYS:H	1.36	0.72
1:C:4:ILE:O	1:C:8:GLU:HG3	1.93	0.69
1:C:51:GLY:H	1:D:88:SER:HB3	1.58	0.67
1:B:90:ILE:HD11	1:B:102:ARG:HG2	1.78	0.66
1:D:113:VAL:HG22	1:D:139:LEU:HB3	1.77	0.66
1:D:109:ASN:HD22	1:D:111:LYS:H	1.44	0.65
1:A:172:MSE:HA	1:A:172:MSE:CE	2.26	0.65
1:A:171:GLU:O	1:A:174:ILE:HG22	1.97	0.63
1:A:41:LYS:HE3	1:A:111:LYS:HG3	1.81	0.63
1:B:90:ILE:CD1	1:B:102:ARG:HG2	2.29	0.62
1:C:114:LEU:HD12	1:C:130:LEU:HD23	1.79	0.62
1:A:74:ALA:HB1	1:C:106:ALA:HB2	1.80	0.61
1:B:109:ASN:HD22	1:B:111:LYS:H	1.47	0.61
1:C:2:SER:H	1:C:5:ASN:HD22	1.49	0.60
1:A:59:HIS:HE1	1:A:178:HIS:NE2	2.00	0.60
1:D:84:THR:O	1:D:88:SER:HB2	2.01	0.60
2:A:239:HOH:O	1:C:90:ILE:HD11	2.02	0.60
1:A:13:GLU:HG3	2:A:373:HOH:O	2.02	0.59
1:B:113:VAL:HG22	1:B:139:LEU:HB3	1.85	0.59

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:150:MSE:HG3	2:C:252:HOH:O	2.04	0.58
1:B:14:HIS:O	1:B:18:VAL:HG23	2.04	0.58
1:B:121:GLY:C	1:B:151:MSE:HE3	2.29	0.58
1:A:59:HIS:CE1	1:A:174:ILE:HD11	2.39	0.57
1:B:31:VAL:HG13	1:B:141:LEU:HD13	1.86	0.57
1:A:4:ILE:O	1:A:8:GLU:HG3	2.04	0.56
1:A:106:ALA:HB2	1:C:74:ALA:HB1	1.88	0.55
1:B:109:ASN:ND2	1:B:111:LYS:H	2.04	0.55
1:B:70:LYS:HE3	2:B:367:HOH:O	2.06	0.55
1:D:109:ASN:ND2	1:D:111:LYS:H	2.04	0.55
1:D:105:GLU:HG3	1:D:132:LYS:HG2	1.88	0.55
1:B:4:ILE:HD12	1:B:4:ILE:N	2.18	0.54
1:A:113:VAL:HG22	1:A:139:LEU:HB3	1.89	0.54
1:C:166:THR:O	1:C:170:GLN:HG3	2.08	0.54
1:C:31:VAL:HG13	1:C:141:LEU:HD13	1.89	0.54
1:C:127:LEU:HD22	1:C:154:LEU:HD12	1.89	0.53
1:D:1:MSE:HG3	1:D:5:ASN:HB2	1.90	0.53
1:A:6:LEU:C	1:A:6:LEU:HD23	2.34	0.53
1:B:146:LYS:HG2	1:B:164:ASP:HA	1.90	0.53
1:C:118:SER:O	1:C:144:SER:HA	2.09	0.53
1:B:4:ILE:H	1:B:4:ILE:CD1	2.16	0.52
1:C:34:LEU:HD22	1:C:139:LEU:HD11	1.90	0.52
1:C:90:ILE:HD13	1:C:102:ARG:CB	2.39	0.52
1:C:17:ILE:HD12	1:C:163:SER:HB2	1.92	0.52
1:D:118:SER:O	1:D:144:SER:HA	2.09	0.52
1:A:59:HIS:HD2	2:A:217:HOH:O	1.93	0.52
1:C:17:ILE:HG13	1:C:168:ARG:HB3	1.90	0.52
1:B:4:ILE:O	1:B:8:GLU:HG3	2.10	0.51
1:C:14:HIS:O	1:C:18:VAL:HG23	2.09	0.51
1:A:118:SER:O	1:A:144:SER:HA	2.11	0.51
1:D:109:ASN:HD21	1:D:111:LYS:HB2	1.75	0.51
1:A:109:ASN:ND2	1:A:111:LYS:H	2.07	0.51
1:A:59:HIS:O	1:A:63:GLU:HG2	2.10	0.51
1:B:165:ASP:O	1:B:169:ILE:HG13	2.11	0.51
1:D:27:GLN:HB3	1:D:159:LEU:CD2	2.40	0.51
1:C:90:ILE:HD13	1:C:102:ARG:HB3	1.93	0.51
1:B:146:LYS:HA	1:B:164:ASP:OD1	2.11	0.50
1:B:127:LEU:HD22	1:B:154:LEU:HD12	1.93	0.50
1:B:74:ALA:HB1	1:D:106:ALA:HB2	1.93	0.50
1:C:27:GLN:HB3	1:C:159:LEU:CD2	2.40	0.50
1:B:59:HIS:O	1:B:63:GLU:HG2	2.12	0.49

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:GLU:HG2	2:C:338:HOH:O	2.12	0.49
1:A:172:MSE:HE2	1:A:175:LEU:HD23	1.95	0.49
1:A:17:ILE:HG13	1:A:168:ARG:HB3	1.95	0.49
1:B:118:SER:CB	1:B:151:MSE:HE1	2.43	0.49
1:D:146:LYS:HE2	1:D:164:ASP:O	2.13	0.49
1:A:27:GLN:HB3	1:A:159:LEU:HD21	1.95	0.48
1:B:174:ILE:HD13	2:C:220:HOH:O	2.13	0.48
1:A:59:HIS:ND1	1:A:174:ILE:HD11	2.28	0.48
1:C:90:ILE:CD1	1:C:102:ARG:HB3	2.44	0.48
1:D:105:GLU:OE2	1:D:132:LYS:HE3	2.14	0.47
1:D:146:LYS:HG2	1:D:164:ASP:HA	1.96	0.47
1:A:17:ILE:HD12	1:A:163:SER:HB2	1.95	0.47
1:D:165:ASP:O	1:D:169:ILE:HG13	2.14	0.47
1:C:1:MSE:HB2	1:C:5:ASN:HB2	1.97	0.47
1:B:118:SER:HB3	1:B:151:MSE:HE1	1.95	0.47
1:D:14:HIS:O	1:D:18:VAL:HG23	2.14	0.47
1:D:27:GLN:HB3	1:D:159:LEU:HD21	1.97	0.47
1:D:31:VAL:HG13	1:D:141:LEU:HD13	1.96	0.47
1:B:2:SER:H	1:B:5:ASN:HD22	1.63	0.47
1:D:27:GLN:O	1:D:31:VAL:HG23	2.15	0.47
1:A:17:ILE:HD11	2:A:259:HOH:O	2.15	0.47
1:B:63:GLU:OE1	1:B:68:TYR:HA	2.15	0.46
1:A:172:MSE:HE2	1:A:175:LEU:CD2	2.45	0.46
1:A:134:LYS:HD3	2:A:238:HOH:O	2.14	0.46
1:B:171:GLU:O	1:B:174:ILE:HG22	2.16	0.46
1:B:106:ALA:HB2	1:D:74:ALA:HB1	1.98	0.46
1:B:6:LEU:C	1:B:6:LEU:HD23	2.40	0.46
1:C:109:ASN:ND2	1:C:111:LYS:H	2.13	0.46
1:B:81:THR:HG22	1:B:100:PHE:CE2	2.51	0.45
1:C:143:LEU:CD2	1:C:159:LEU:HD12	2.47	0.45
1:B:27:GLN:HB3	1:B:159:LEU:CD2	2.47	0.45
1:C:81:THR:HG22	1:C:100:PHE:CE2	2.51	0.45
1:D:4:ILE:O	1:D:8:GLU:HG3	2.15	0.45
1:A:92:ASN:HD21	1:B:124:PRO:HD2	1.82	0.45
1:B:9:LYS:HE3	2:B:252:HOH:O	2.16	0.45
1:D:114:LEU:HD23	1:D:130:LEU:HA	1.98	0.44
1:A:13:GLU:O	1:A:17:ILE:HG12	2.17	0.44
1:A:96:PHE:O	1:A:99:VAL:HG22	2.17	0.44
1:A:121:GLY:HA2	1:A:151:MSE:SE	2.67	0.44
1:C:96:PHE:HE2	1:D:92:ASN:HB2	1.82	0.44
1:B:118:SER:O	1:B:144:SER:HA	2.17	0.44

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:LYS:HG2	1:A:111:LYS:HD2	2.00	0.44
1:B:27:GLN:HB3	1:B:159:LEU:HD21	1.99	0.43
1:C:9:LYS:HD3	2:C:228:HOH:O	2.18	0.43
1:D:114:LEU:HD22	1:D:133:ALA:HB2	2.00	0.43
1:C:174:ILE:HD11	1:C:178:HIS:NE2	2.33	0.43
1:C:109:ASN:HD21	1:C:111:LYS:HD3	1.83	0.43
1:C:121:GLY:HA2	1:C:151:MSE:SE	2.69	0.43
1:D:81:THR:HG22	1:D:100:PHE:CZ	2.54	0.43
1:D:2:SER:H	1:D:5:ASN:HD22	1.64	0.43
1:D:6:LEU:C	1:D:6:LEU:HD23	2.43	0.43
1:B:90:ILE:HD11	2:D:283:HOH:O	2.18	0.43
1:C:27:GLN:HB3	1:C:159:LEU:HD22	2.01	0.43
1:B:9:LYS:O	1:B:13:GLU:HG2	2.19	0.42
1:B:38:CYS:HA	1:B:41:LYS:HE2	2.02	0.42
1:B:105:GLU:HG3	1:B:132:LYS:HG2	2.01	0.42
1:C:13:GLU:O	1:C:17:ILE:HG12	2.19	0.42
1:D:121:GLY:HA2	1:D:151:MSE:SE	2.69	0.42
1:D:59:HIS:O	1:D:63:GLU:HG2	2.20	0.42
1:A:33:GLU:HG2	2:A:314:HOH:O	2.19	0.41
1:C:63:GLU:OE1	1:C:68:TYR:HA	2.20	0.41
1:C:71:GLU:HA	1:C:71:GLU:OE1	2.19	0.41
1:A:12:GLN:O	1:A:16:LYS:HG3	2.20	0.41
1:C:92:ASN:HB2	1:D:96:PHE:HE2	1.85	0.41
1:D:13:GLU:O	1:D:17:ILE:HG12	2.21	0.41
1:C:114:LEU:C	1:C:114:LEU:HD13	2.46	0.41
1:A:27:GLN:HB3	1:A:159:LEU:CD2	2.51	0.41
1:D:18:VAL:HG22	1:D:172:MSE:HE1	2.03	0.41
1:D:81:THR:HG22	1:D:100:PHE:CE2	2.56	0.40
1:C:38:CYS:SG	1:C:113:VAL:HG23	2.62	0.40
1:B:90:ILE:HD13	1:B:102:ARG:CB	2.52	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [\(i\)](#)

### 5.3.1 Protein backbone [\(i\)](#)

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	186/188 (99%)	180 (97%)	6 (3%)	0	100	100
1	B	186/188 (99%)	181 (97%)	5 (3%)	0	100	100
1	C	186/188 (99%)	180 (97%)	6 (3%)	0	100	100
1	D	186/188 (99%)	181 (97%)	5 (3%)	0	100	100
All	All	744/752 (99%)	722 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	153/148 (103%)	149 (97%)	4 (3%)	40	46
1	B	153/148 (103%)	152 (99%)	1 (1%)	76	83
1	C	153/148 (103%)	151 (99%)	2 (1%)	61	69
1	D	153/148 (103%)	148 (97%)	5 (3%)	33	37
All	All	612/592 (103%)	600 (98%)	12 (2%)	48	56

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	GLU
1	A	110	GLU
1	A	111	LYS
1	A	172	MSE
1	B	172	MSE
1	C	109	ASN
1	C	139	LEU
1	D	67	ARG
1	D	73	LYS
1	D	88	SER

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	D	111	LYS
1	D	134	LYS

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

Mol	Chain	Res	Type
1	A	19	GLN
1	A	27	GLN
1	A	59	HIS
1	A	92	ASN
1	A	109	ASN
1	B	5	ASN
1	B	27	GLN
1	B	103	GLN
1	B	109	ASN
1	C	5	ASN
1	C	103	GLN
1	C	109	ASN
1	C	157	HIS
1	D	5	ASN
1	D	12	GLN
1	D	58	GLN
1	D	109	ASN
1	D	158	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	183/188 (97%)	-0.36	0 <a href="#">100</a> <a href="#">100</a>	16, 21, 33, 43	0
1	B	183/188 (97%)	-0.30	0 <a href="#">100</a> <a href="#">100</a>	16, 23, 35, 44	0
1	C	183/188 (97%)	-0.15	2 (1%) <a href="#">78</a> <a href="#">80</a>	18, 27, 37, 46	0
1	D	183/188 (97%)	-0.26	0 <a href="#">100</a> <a href="#">100</a>	17, 25, 36, 41	0
All	All	732/752 (97%)	-0.27	2 (0%) <a href="#">90</a> <a href="#">91</a>	16, 23, 36, 46	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	71	GLU	2.2
1	C	137	ASN	2.2

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.

### 6.5 Other polymers [i](#)

There are no such residues in this entry.