



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:51 PM GMT

PDB ID : 1VGA  
Title : Structures of unligated and inhibitor complexes of W168F mutant of  
Triosephosphate Isomerase from Plasmodium falciparum  
Authors : Eaazhisai, K.; Balaram, H.; Balaram, P.; Murthy, M.R.N.  
Deposited on : 2004-04-23  
Resolution : 1.80 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

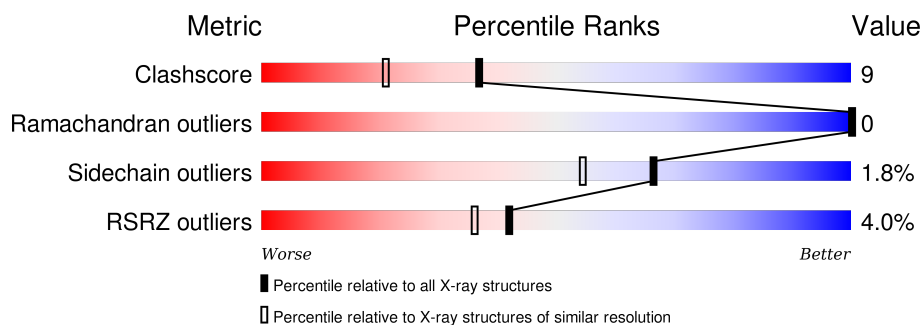
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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. 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	248	<div> <div>2%</div> <div>79%</div> <div>20%</div> <div>.</div> </div>
1	B	248	<div> <div>5%</div> <div>75%</div> <div>24%</div> <div>.</div> </div>
1	C	248	<div> <div>5%</div> <div>84%</div> <div>14%</div> <div>..</div> </div>
1	D	248	<div> <div>4%</div> <div>77%</div> <div>22%</div> <div>.</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8374 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 Triosephosphate isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	246	Total	C	N	O	S	0	0	0
			1954	1237	332	380	5			
1	B	246	Total	C	N	O	S	0	0	0
			1954	1237	332	380	5			
1	C	246	Total	C	N	O	S	0	0	0
			1954	1237	332	380	5			
1	D	246	Total	C	N	O	S	0	0	0
			1954	1237	332	380	5			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	163	VAL	ALA	ENGINEERED	UNP Q07412
A	168	PHE	TRP	ENGINEERED	UNP Q07412
B	163	VAL	ALA	ENGINEERED	UNP Q07412
B	168	PHE	TRP	ENGINEERED	UNP Q07412
C	163	VAL	ALA	ENGINEERED	UNP Q07412
C	168	PHE	TRP	ENGINEERED	UNP Q07412
D	163	VAL	ALA	ENGINEERED	UNP Q07412
D	168	PHE	TRP	ENGINEERED	UNP Q07412

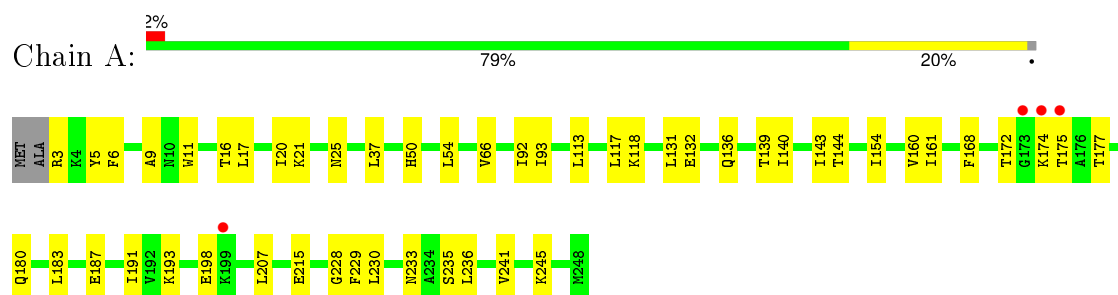
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	172	Total	O	0	0
			172	172		
2	B	144	Total	O	0	0
			144	144		
2	C	111	Total	O	0	0
			111	111		
2	D	131	Total	O	0	0
			131	131		

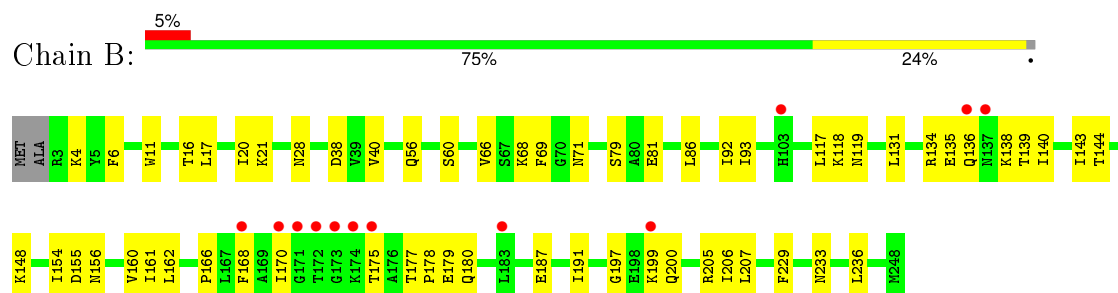
### 3 Residue-property plots [i](#)

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.

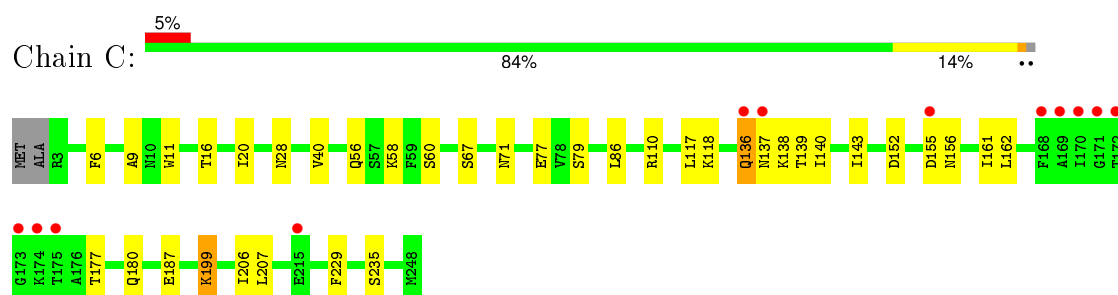
- Molecule 1: Triosephosphate isomerase



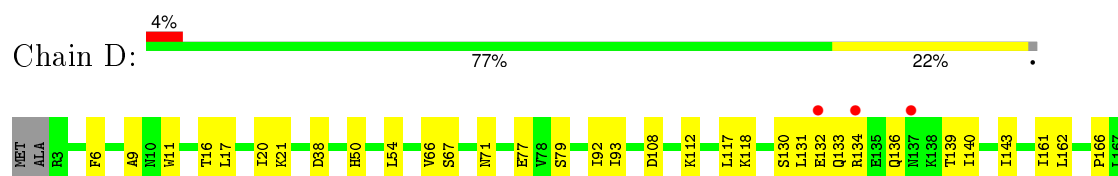
- Molecule 1: Triosephosphate isomerase

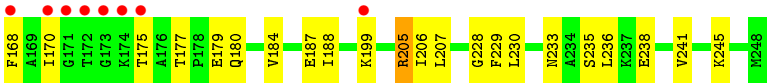


- Molecule 1: Triosephosphate isomerase



- Molecule 1: Triosephosphate isomerase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.73Å 106.16Å 89.53Å 90.00° 92.42° 90.00°	Depositor
Resolution (Å)	17.20 – 1.80 17.21 – 1.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (17.20-1.80) 93.6 (17.21-1.80)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.48 (at 1.80Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.207 , 0.229 0.210 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	19.9	Xtriage
Anisotropy	0.299	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 50.0	EDS
Estimated twinning fraction	0.028 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 87947 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8374	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.68% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.31	0/1985	0.60	1/2675 (0.0%)
1	B	0.32	0/1985	0.59	0/2675
1	C	0.31	0/1985	0.59	0/2675
1	D	0.32	0/1985	0.61	1/2675 (0.0%)
All	All	0.31	0/7940	0.60	2/10700 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	230	LEU	N-CA-C	-5.40	96.42	111.00
1	A	230	LEU	N-CA-C	-5.30	96.69	111.00

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	1954	0	1952	42	0
1	B	1954	0	1952	40	0
1	C	1954	0	1952	37	0
1	D	1954	0	1952	36	0
2	A	172	0	0	5	0
2	B	144	0	0	0	0
2	C	111	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	131	0	0	2	0
All	All	8374	0	7808	147	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 9.

All (147) 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)
1:D:140:ILE:HD12	1:D:187:GLU:HG2	1.45	0.98
1:A:154:ILE:HD12	1:A:160:VAL:HG21	1.46	0.94
1:B:177:THR:HB	1:C:136:GLN:OE1	1.74	0.88
1:C:177:THR:HG22	1:C:180:GLN:HG3	1.60	0.82
1:A:233:ASN:HA	1:A:236:LEU:HD23	1.65	0.77
1:B:154:ILE:HD12	1:B:160:VAL:HG21	1.66	0.75
1:C:136:GLN:CB	1:C:138:LYS:HE2	2.17	0.74
1:B:233:ASN:HA	1:B:236:LEU:HD23	1.75	0.69
1:C:140:ILE:HD12	1:C:187:GLU:HG2	1.75	0.68
1:A:154:ILE:CD1	1:A:160:VAL:HG21	2.24	0.67
1:C:86:LEU:HD21	1:D:17:LEU:HD21	1.77	0.66
1:D:11:TRP:CZ3	1:D:20:ILE:HD12	2.31	0.66
1:B:177:THR:OG1	1:B:180:GLN:HG3	1.96	0.65
1:C:16:THR:O	1:C:20:ILE:HG12	1.96	0.65
1:D:132:GLU:O	1:D:136:GLN:HG3	1.95	0.65
1:A:154:ILE:HD12	1:A:160:VAL:CG2	2.24	0.64
1:C:177:THR:HG22	1:C:180:GLN:CG	2.28	0.63
1:B:11:TRP:CZ3	1:B:20:ILE:HD12	2.33	0.63
1:A:236:LEU:HD21	2:A:360:HOH:O	1.98	0.63
1:B:178:PRO:HD2	1:C:136:GLN:OE1	2.01	0.61
1:C:136:GLN:HB3	1:C:138:LYS:HE2	1.83	0.61
1:B:38:ASP:OD2	1:B:205:ARG:NH2	2.34	0.61
1:D:38:ASP:OD2	1:D:205:ARG:NH2	2.30	0.61
1:B:177:THR:HB	1:C:136:GLN:CD	2.21	0.60
1:B:144:THR:O	1:B:148:LYS:HG3	2.01	0.60
1:D:162:LEU:HD12	1:D:206:ILE:HD12	1.84	0.59
1:A:11:TRP:CZ3	1:A:20:ILE:HD12	2.39	0.58
1:B:162:LEU:HD12	1:B:206:ILE:HD12	1.86	0.58
1:D:16:THR:O	1:D:20:ILE:HG12	2.04	0.57
1:B:140:ILE:HD12	1:B:187:GLU:HG2	1.85	0.57
1:C:136:GLN:HA	1:C:136:GLN:HE21	1.69	0.57
1:D:92:ILE:O	1:D:93:ILE:HD13	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:LEU:CD1	1:A:117:LEU:HD22	2.35	0.57
1:D:162:LEU:HD12	1:D:206:ILE:CD1	2.35	0.56
1:A:66:VAL:HG12	1:A:93:ILE:HD11	1.87	0.56
1:A:50:HIS:HD2	2:A:380:HOH:O	1.88	0.56
1:B:131:LEU:O	1:B:135:GLU:HG3	2.06	0.56
1:C:162:LEU:HD12	1:C:206:ILE:HD12	1.86	0.56
1:A:236:LEU:N	1:A:236:LEU:HD22	2.22	0.55
1:C:199:LYS:HE2	2:C:296:HOH:O	2.05	0.55
1:C:58:LYS:HD2	2:C:328:HOH:O	2.06	0.55
1:C:136:GLN:HB2	1:C:138:LYS:HE2	1.89	0.55
1:C:110:ARG:NH1	1:C:152:ASP:OD1	2.40	0.55
1:C:177:THR:CG2	1:C:180:GLN:HG3	2.35	0.55
1:B:205:ARG:O	1:B:206:ILE:HD13	2.07	0.55
1:C:6:PHE:HB3	1:C:207:LEU:HD21	1.89	0.55
1:B:236:LEU:N	1:B:236:LEU:HD22	2.22	0.54
1:A:172:THR:OG1	1:A:174:LYS:HB2	2.08	0.54
1:B:16:THR:O	1:B:20:ILE:HG12	2.07	0.54
1:B:136:GLN:CB	1:B:138:LYS:HE2	2.37	0.54
1:C:177:THR:HG23	1:C:180:GLN:H	1.72	0.54
1:B:66:VAL:HG12	1:B:93:ILE:CD1	2.38	0.54
1:D:130:SER:OG	1:D:133:GLN:HG3	2.08	0.53
1:A:140:ILE:HD12	1:A:187:GLU:HG2	1.90	0.53
1:A:66:VAL:HG12	1:A:93:ILE:CD1	2.39	0.53
1:B:68:LYS:HE3	1:B:69:PHE:CZ	2.44	0.52
1:A:241:VAL:O	1:A:245:LYS:HG3	2.10	0.52
1:C:139:THR:O	1:C:143:ILE:HG12	2.09	0.52
1:C:118:LYS:NZ	1:C:118:LYS:HB2	2.24	0.52
1:C:86:LEU:HD23	1:D:17:LEU:HD11	1.92	0.52
1:A:174:LYS:HD2	2:A:350:HOH:O	2.09	0.52
1:B:92:ILE:O	1:B:93:ILE:HD13	2.10	0.52
1:A:92:ILE:O	1:A:93:ILE:HD13	2.10	0.51
1:C:9:ALA:HB1	1:C:235:SER:HB2	1.91	0.51
1:A:131:LEU:HD13	1:A:168:PHE:HB2	1.91	0.51
1:C:136:GLN:CA	1:C:136:GLN:HE21	2.22	0.51
1:D:238:GLU:O	1:D:241:VAL:HG22	2.11	0.51
1:A:16:THR:O	1:A:20:ILE:HG12	2.09	0.51
1:B:155:ASP:OD1	1:B:156:ASN:N	2.43	0.51
1:B:161:ILE:N	1:B:161:ILE:HD12	2.26	0.50
1:C:162:LEU:HD12	1:C:206:ILE:CD1	2.41	0.50
1:D:179:GLU:HB2	2:D:350:HOH:O	2.11	0.50
1:D:118:LYS:NZ	1:D:118:LYS:HB2	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:6:PHE:HB3	1:D:207:LEU:HD21	1.94	0.50
1:B:6:PHE:HB3	1:B:207:LEU:HD21	1.93	0.49
1:D:177:THR:OG1	1:D:180:GLN:HG3	2.12	0.49
1:D:139:THR:O	1:D:143:ILE:HG12	2.11	0.49
1:A:118:LYS:HB2	1:A:118:LYS:NZ	2.26	0.49
1:B:71:ASN:ND2	1:B:79:SER:OG	2.46	0.49
1:D:175:THR:HG23	2:D:348:HOH:O	2.13	0.49
1:C:86:LEU:CD2	1:D:17:LEU:HD21	2.43	0.48
1:A:17:LEU:HD21	1:B:86:LEU:CD2	2.43	0.48
1:D:131:LEU:HD13	1:D:168:PHE:HB2	1.95	0.48
1:B:81:GLU:HG3	1:B:119:ASN:OD1	2.13	0.48
1:B:134:ARG:HD2	1:B:168:PHE:CD1	2.48	0.48
1:D:241:VAL:O	1:D:245:LYS:HG3	2.13	0.48
1:D:71:ASN:ND2	1:D:79:SER:OG	2.47	0.48
1:A:9:ALA:HB1	1:A:235:SER:HB2	1.95	0.47
1:B:118:LYS:HB2	1:B:118:LYS:NZ	2.29	0.47
1:A:161:ILE:N	1:A:161:ILE:HD12	2.30	0.47
1:A:113:LEU:HD11	1:A:117:LEU:HD22	1.95	0.47
1:B:200:GLN:OE1	1:B:200:GLN:N	2.46	0.47
1:D:166:PRO:O	1:D:170:ILE:HG13	2.14	0.47
1:B:179:GLU:OE1	1:C:136:GLN:OE1	2.33	0.47
1:A:6:PHE:HB3	1:A:207:LEU:HD21	1.95	0.47
1:B:187:GLU:O	1:B:191:ILE:HG13	2.15	0.47
1:B:166:PRO:HB2	1:B:168:PHE:CE1	2.50	0.47
1:B:17:LEU:O	1:B:21:LYS:HG3	2.14	0.46
1:C:177:THR:CG2	1:C:180:GLN:H	2.28	0.46
1:D:17:LEU:O	1:D:50:HIS:HE1	1.98	0.46
1:D:21:LYS:HA	1:D:54:LEU:HD13	1.96	0.46
1:C:71:ASN:ND2	1:C:79:SER:OG	2.49	0.46
1:A:21:LYS:HA	1:A:54:LEU:HD13	1.97	0.46
1:D:66:VAL:HG12	1:D:93:ILE:CD1	2.45	0.46
1:B:136:GLN:HB3	1:B:138:LYS:HE2	1.98	0.46
1:B:40:VAL:HG22	1:B:60:SER:HB2	1.99	0.45
1:C:11:TRP:CZ3	1:C:20:ILE:HD12	2.52	0.45
1:A:21:LYS:O	1:A:25:ASN:ND2	2.50	0.45
1:D:233:ASN:HA	1:D:236:LEU:HD23	1.97	0.45
1:A:113:LEU:HD13	1:A:117:LEU:HD22	1.98	0.45
1:A:233:ASN:CA	1:A:236:LEU:HD23	2.40	0.45
1:A:132:GLU:HG2	1:A:136:GLN:NE2	2.33	0.44
1:D:134:ARG:HD2	1:D:168:PHE:CD1	2.53	0.44
1:B:139:THR:O	1:B:143:ILE:HG12	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:THR:OG1	1:A:180:GLN:HG3	2.18	0.44
1:A:233:ASN:O	1:A:236:LEU:HD23	2.18	0.44
1:B:28:ASN:HB3	1:B:56:GLN:NE2	2.33	0.43
1:A:6:PHE:O	1:A:228:GLY:HA3	2.17	0.43
1:C:136:GLN:NE2	1:C:136:GLN:HA	2.33	0.43
1:C:161:ILE:N	1:C:161:ILE:HD12	2.34	0.43
1:B:4:LYS:HB3	1:B:205:ARG:HH21	1.83	0.43
1:D:161:ILE:HD12	1:D:161:ILE:N	2.34	0.43
1:A:154:ILE:CD1	1:A:160:VAL:HG11	2.49	0.43
1:A:139:THR:O	1:A:143:ILE:HG12	2.18	0.43
1:A:233:ASN:HA	1:A:236:LEU:CD2	2.42	0.42
1:C:28:ASN:HB3	1:C:56:GLN:NE2	2.34	0.42
1:C:67:SER:HB2	1:C:77:GLU:HB3	2.00	0.42
1:B:166:PRO:O	1:B:170:ILE:HG12	2.20	0.42
1:D:67:SER:HB2	1:D:77:GLU:HB3	2.00	0.42
1:D:9:ALA:HB1	1:D:235:SER:HB2	2.01	0.42
1:A:5:TYR:O	1:A:37:LEU:HD12	2.20	0.42
1:D:184:VAL:O	1:D:188:ILE:HG13	2.19	0.42
1:A:193:LYS:HG3	1:A:198:GLU:HA	2.02	0.42
1:A:113:LEU:C	1:A:113:LEU:HD13	2.41	0.41
1:D:236:LEU:N	1:D:236:LEU:HD22	2.35	0.41
1:A:236:LEU:CD2	1:A:236:LEU:N	2.83	0.41
1:D:6:PHE:O	1:D:228:GLY:HA3	2.20	0.41
1:D:108:ASP:O	1:D:112:LYS:HG3	2.20	0.41
1:A:183:LEU:HG	2:A:414:HOH:O	2.21	0.41
1:A:3:ARG:NH2	2:A:307:HOH:O	2.52	0.41
1:B:236:LEU:N	1:B:236:LEU:CD2	2.83	0.41
1:C:40:VAL:HG22	1:C:60:SER:HB2	2.02	0.41
1:A:144:THR:HA	1:A:191:ILE:HD13	2.01	0.41
1:D:131:LEU:CD1	1:D:168:PHE:HB2	2.51	0.41
1:C:155:ASP:OD1	1:C:156:ASN:N	2.54	0.41
1:B:197:GLY:HA3	1:B:200:GLN:NE2	2.36	0.40
1:C:28:ASN:HB3	1:C:56:GLN:HE21	1.86	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	244/248 (98%)	239 (98%)	5 (2%)	0	100	100
1	B	244/248 (98%)	236 (97%)	8 (3%)	0	100	100
1	C	244/248 (98%)	234 (96%)	10 (4%)	0	100	100
1	D	244/248 (98%)	237 (97%)	7 (3%)	0	100	100
All	All	976/992 (98%)	946 (97%)	30 (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	219/220 (100%)	216 (99%)	3 (1%)	74	65
1	B	219/220 (100%)	215 (98%)	4 (2%)	66	54
1	C	219/220 (100%)	214 (98%)	5 (2%)	58	42
1	D	219/220 (100%)	215 (98%)	4 (2%)	66	54
All	All	876/880 (100%)	860 (98%)	16 (2%)	66	54

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	175	THR
1	A	215	GLU

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Mol	Chain	Res	Type
1	A	229	PHE
1	B	117	LEU
1	B	175	THR
1	B	199	LYS
1	B	229	PHE
1	C	117	LEU
1	C	136	GLN
1	C	137	ASN
1	C	199	LYS
1	C	229	PHE
1	D	117	LEU
1	D	199	LYS
1	D	205	ARG
1	D	229	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	50	HIS
1	A	71	ASN
1	A	87	ASN
1	A	159	ASN
1	A	203	GLN
1	A	222	GLN
1	B	29	ASN
1	B	56	GLN
1	B	71	ASN
1	B	87	ASN
1	B	159	ASN
1	B	222	GLN
1	C	29	ASN
1	C	50	HIS
1	C	56	GLN
1	C	71	ASN
1	C	87	ASN
1	C	159	ASN
1	C	203	GLN
1	C	222	GLN
1	D	29	ASN
1	D	50	HIS
1	D	71	ASN

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Mol	Chain	Res	Type
1	D	87	ASN
1	D	159	ASN
1	D	203	GLN

### 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 carbohydrates 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

### 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	246/248 (99%)	0.08	4 (1%) 74 71	13, 19, 38, 47	0
1	B	246/248 (99%)	0.21	12 (4%) 33 27	13, 20, 42, 61	0
1	C	246/248 (99%)	0.47	12 (4%) 33 27	14, 21, 47, 84	0
1	D	246/248 (99%)	0.19	11 (4%) 37 31	14, 20, 42, 64	0
All	All	984/992 (99%)	0.24	39 (3%) 42 36	13, 20, 43, 84	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	173	GLY	21.8
1	C	171	GLY	10.6
1	C	172	THR	8.7
1	C	170	ILE	7.5
1	C	168	PHE	6.7
1	D	171	GLY	6.4
1	C	174	LYS	6.0
1	B	172	THR	5.0
1	B	173	GLY	4.4
1	B	170	ILE	4.3
1	D	173	GLY	4.2
1	C	136	GLN	4.1
1	C	215	GLU	3.9
1	B	183	LEU	3.8
1	D	174	LYS	3.8
1	C	175	THR	3.6
1	B	136	GLN	3.5
1	D	172	THR	3.4
1	D	137	ASN	3.3
1	C	155	ASP	3.3
1	B	174	LYS	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	168	PHE	3.2
1	B	175	THR	3.2
1	D	170	ILE	2.8
1	A	174	LYS	2.7
1	C	137	ASN	2.6
1	B	137	ASN	2.6
1	A	173	GLY	2.5
1	D	132	GLU	2.4
1	B	171	GLY	2.3
1	C	169	ALA	2.3
1	D	134	ARG	2.2
1	D	168	PHE	2.2
1	D	175	THR	2.1
1	B	103	HIS	2.1
1	A	199	LYS	2.1
1	A	175	THR	2.1
1	D	199	LYS	2.0
1	B	199	LYS	2.0

## 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 carbohydrates 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.