



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2024 – 05:08 AM EST

PDB ID : 1WYI  
Title : human triosephosphate isomerase of new crystal form  
Authors : Kinoshita, T.  
Deposited on : 2005-02-14  
Resolution : 2.20 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>.

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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  
Xtrriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

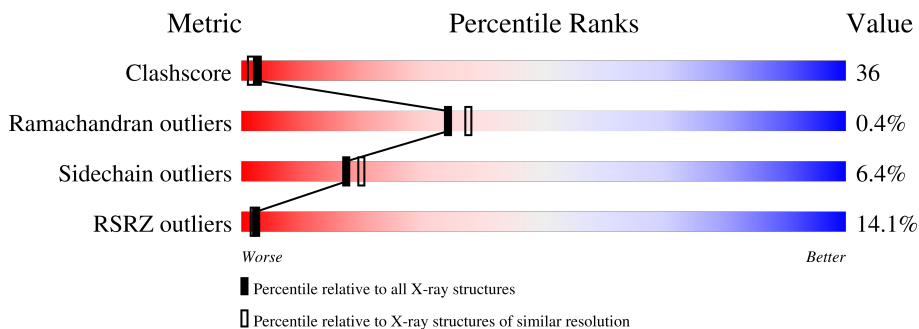
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	250	
1	B	250	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4196 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
			Total	C	N	O	S			
1	A	248	1868	1181	324	356	7	0	0	0
1	B	248	1868	1181	324	356	7	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	cloning artifact	UNP P60174
A	0	SER	-	cloning artifact	UNP P60174
B	-1	GLY	-	cloning artifact	UNP P60174
B	0	SER	-	cloning artifact	UNP P60174

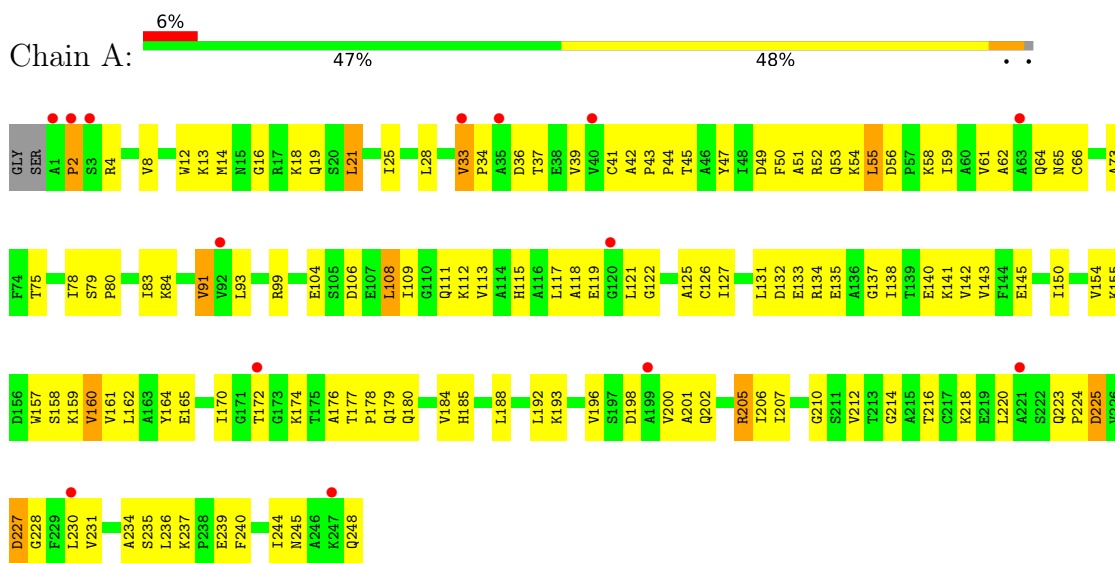
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	240	Total	O	0	0
			240	240		
2	B	220	Total	O	0	0
			220	220		

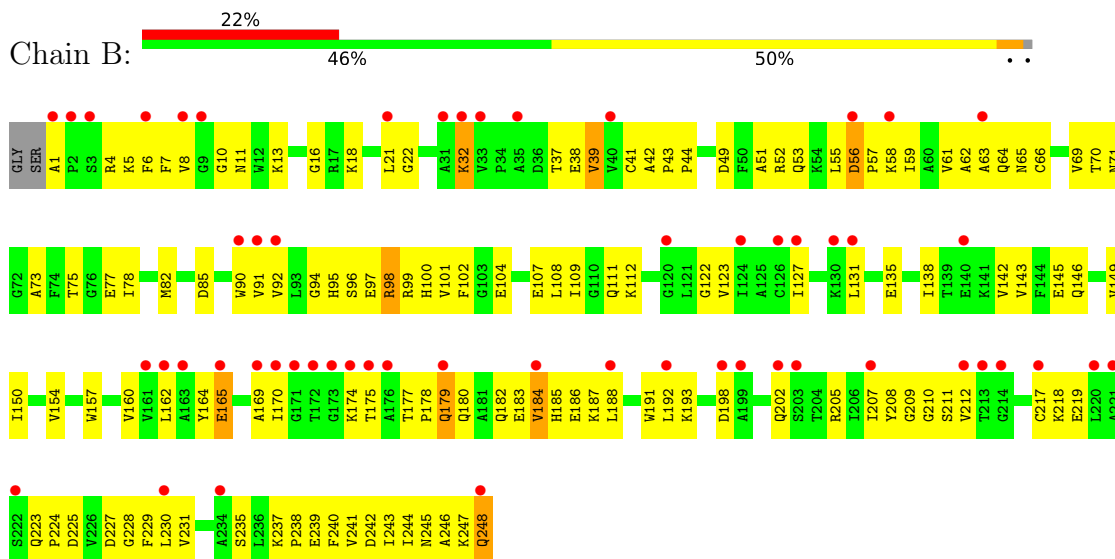
### 3 Residue-property plots

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: Triosephosphate isomerase



- Molecule 1: Triosephosphate isomerase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.56Å 70.88Å 141.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.20 – 2.20 28.33 – 2.20	Depositor EDS
% Data completeness (in resolution range)	64.6 (28.20-2.20) 96.1 (28.33-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.85 (at 2.20Å)	Xtrriage
Refinement program	CNX 2002	Depositor
R, $R_{free}$	0.247 , 0.296 0.267 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.8	Xtrriage
Anisotropy	0.400	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 55.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4196	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.74% 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	0/1903	0.66	1/2577 (0.0%)
1	B	0.33	0/1903	0.57	0/2577
All	All	0.37	0/3806	0.62	1/5154 (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	A	230	LEU	N-CA-C	-5.37	96.50	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	1868	0	1879	138	0
1	B	1868	0	1879	145	0
2	A	240	0	0	46	0
2	B	220	0	0	37	0
All	All	4196	0	3758	270	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 36.

All (270) 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:B:78:ILE:HB	2:B:421:HOH:O	1.62	1.00
1:A:108:LEU:HD11	1:A:112:LYS:HE3	1.48	0.95
1:A:138:ILE:HG22	1:A:142:VAL:HG23	1.50	0.93
1:A:25:ILE:HG23	1:A:55:LEU:HD13	1.56	0.88
1:B:98:ARG:HA	1:B:98:ARG:HH11	1.37	0.87
1:B:92:VAL:HG21	2:B:450:HOH:O	1.75	0.85
1:A:25:ILE:HA	2:A:380:HOH:O	1.76	0.85
1:B:127:ILE:HD12	1:B:143:VAL:HG13	1.58	0.83
1:A:93:LEU:HB2	2:A:382:HOH:O	1.79	0.83
1:A:143:VAL:HG11	1:A:188:LEU:HD11	1.60	0.83
1:B:142:VAL:O	1:B:146:GLN:HG3	1.78	0.82
1:B:13:LYS:HA	2:B:373:HOH:O	1.79	0.81
1:B:98:ARG:HH12	1:B:102:PHE:HB2	1.46	0.80
1:B:44:PRO:HD3	2:B:368:HOH:O	1.82	0.79
1:A:125:ALA:HB1	1:A:150:ILE:HD13	1.64	0.78
1:B:92:VAL:HG12	2:B:394:HOH:O	1.82	0.77
1:A:225:ASP:HA	2:A:459:HOH:O	1.83	0.77
1:A:207:ILE:HG22	1:A:228:GLY:HA3	1.69	0.75
1:B:177:THR:HB	1:B:179:GLN:HE21	1.53	0.74
1:A:75:THR:HG23	1:B:64:GLN:HB3	1.71	0.73
1:B:108:LEU:HD11	1:B:112:LYS:HE3	1.69	0.72
1:A:188:LEU:HD22	2:A:413:HOH:O	1.88	0.72
1:B:175:THR:HA	1:B:211:SER:HB2	1.72	0.71
1:B:183:GLU:HG3	2:B:454:HOH:O	1.89	0.71
1:A:212:VAL:HG13	2:A:479:HOH:O	1.90	0.70
1:A:8:VAL:HG23	2:A:444:HOH:O	1.91	0.70
1:A:108:LEU:CD1	1:A:112:LYS:HE3	2.21	0.70
1:A:12:TRP:HD1	1:A:235:SER:HG	1.40	0.69
1:A:218:LYS:HB2	2:A:383:HOH:O	1.91	0.69
1:A:51:ALA:O	1:A:55:LEU:HB2	1.92	0.69
1:A:170:ILE:O	1:A:172:THR:HG23	1.93	0.69
1:A:138:ILE:HG22	1:A:138:ILE:O	1.93	0.69
1:B:98:ARG:NH1	1:B:102:PHE:HB2	2.08	0.68
1:A:118:ALA:HB2	2:A:260:HOH:O	1.93	0.68
1:B:164:TYR:CE2	1:B:184:VAL:HG11	2.29	0.68
1:B:127:ILE:HG22	2:B:444:HOH:O	1.94	0.67
1:B:100:HIS:HB2	2:B:361:HOH:O	1.94	0.67
1:B:212:VAL:HG12	1:B:243:ILE:HD13	1.76	0.67
1:A:28:LEU:HB2	2:A:380:HOH:O	1.94	0.67
1:B:98:ARG:HG3	1:B:104:GLU:OE2	1.94	0.67
1:B:101:VAL:HG23	2:B:361:HOH:O	1.95	0.67
1:B:240:PHE:O	1:B:244:ILE:HG13	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:LEU:HB3	2:A:391:HOH:O	1.95	0.66
1:B:98:ARG:HH11	1:B:98:ARG:CA	2.08	0.66
1:B:43:PRO:HD3	2:B:422:HOH:O	1.94	0.66
1:B:170:ILE:HG23	2:B:342:HOH:O	1.96	0.65
1:A:127:ILE:HD12	1:A:164:TYR:HA	1.77	0.65
1:B:193:LYS:HG3	1:B:198:ASP:HA	1.79	0.65
1:A:138:ILE:HG23	1:A:141:LYS:HB3	1.77	0.65
1:A:41:CYS:O	1:A:43:PRO:HD3	1.97	0.64
1:A:138:ILE:CG2	1:A:142:VAL:HG23	2.25	0.64
1:A:231:VAL:HG23	2:A:444:HOH:O	1.97	0.63
1:A:174:LYS:HE2	2:A:469:HOH:O	1.98	0.63
1:B:237:LYS:HB2	1:B:239:GLU:HG2	1.81	0.63
1:B:16:GLY:HA3	1:B:21:LEU:CD1	2.29	0.62
1:A:45:THR:N	1:B:78:ILE:HD11	2.15	0.62
1:B:62:ALA:HB2	1:B:90:TRP:HB2	1.80	0.62
1:B:99:ARG:HG3	1:B:100:HIS:CD2	2.35	0.62
1:B:143:VAL:HG11	1:B:188:LEU:HD11	1.81	0.62
1:A:179:GLN:H	1:A:179:GLN:NE2	1.98	0.62
1:A:184:VAL:HA	2:A:424:HOH:O	1.98	0.61
1:A:73:ALA:HA	1:B:13:LYS:HD3	1.81	0.61
1:A:161:VAL:HG23	2:A:315:HOH:O	1.99	0.61
1:A:75:THR:CG2	1:B:64:GLN:HB3	2.29	0.61
1:A:108:LEU:HA	1:A:111:GLN:OE1	2.01	0.61
1:B:99:ARG:HD3	2:B:252:HOH:O	2.01	0.61
1:B:13:LYS:N	1:B:64:GLN:HE22	1.99	0.60
1:B:247:LYS:HD3	1:B:248:GLN:N	2.16	0.60
1:B:98:ARG:HH12	1:B:102:PHE:CB	2.13	0.60
1:A:138:ILE:CG2	1:A:141:LYS:HB3	2.32	0.60
1:A:185:HIS:HA	2:A:413:HOH:O	2.00	0.60
1:A:216:THR:HB	2:A:300:HOH:O	2.02	0.60
1:A:19:GLN:HB3	2:A:250:HOH:O	2.00	0.59
1:B:52:ARG:HB2	2:B:365:HOH:O	2.02	0.59
1:A:176:ALA:HA	1:A:180:GLN:OE1	2.03	0.58
1:A:64:GLN:HB3	1:B:75:THR:HG23	1.85	0.58
1:B:56:ASP:OD2	1:B:58:LYS:HB2	2.03	0.58
1:A:133:GLU:HG3	2:A:435:HOH:O	2.03	0.58
1:B:184:VAL:O	1:B:188:LEU:HD13	2.04	0.58
1:B:64:GLN:HG3	2:B:294:HOH:O	2.02	0.58
1:B:98:ARG:HA	1:B:98:ARG:NH1	2.14	0.58
1:B:138:ILE:HG22	2:B:298:HOH:O	2.03	0.58
1:A:54:LYS:HE3	2:A:465:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:TRP:HZ3	1:B:162:LEU:HD21	1.68	0.57
1:A:134:ARG:HB3	2:A:440:HOH:O	2.04	0.57
1:B:91:VAL:HG12	1:B:123:VAL:HG12	1.86	0.57
1:A:104:GLU:OE1	1:A:112:LYS:NZ	2.36	0.57
1:B:94:GLY:HA2	1:B:109:ILE:HD13	1.85	0.57
1:A:108:LEU:HD11	1:A:112:LYS:CE	2.31	0.57
1:A:78:ILE:HG13	2:A:375:HOH:O	2.03	0.57
1:B:223:GLN:HE21	1:B:224:PRO:HD2	1.70	0.57
1:A:193:LYS:HG3	1:A:198:ASP:HA	1.87	0.57
1:B:127:ILE:O	1:B:165:GLU:HG3	2.05	0.57
1:A:161:VAL:HG12	1:A:205:ARG:NH1	2.20	0.56
1:A:84:LYS:HG2	1:A:121:LEU:HD13	1.86	0.56
1:B:162:LEU:O	1:B:207:ILE:HG12	2.05	0.56
1:B:10:GLY:HA3	2:B:405:HOH:O	2.05	0.56
1:A:193:LYS:N	1:A:201:ALA:HB2	2.20	0.56
1:A:212:VAL:HA	2:A:479:HOH:O	2.05	0.56
1:A:117:LEU:HD11	1:A:154:VAL:HG22	1.88	0.56
1:A:126:CYS:HB3	1:A:165:GLU:OE1	2.05	0.56
1:A:132:ASP:HA	2:A:400:HOH:O	2.05	0.55
1:A:109:ILE:O	1:A:113:VAL:HG23	2.05	0.55
1:B:157:TRP:O	1:B:160:VAL:HG12	2.05	0.55
1:A:59:ILE:HB	2:A:391:HOH:O	2.06	0.55
1:B:77:GLU:O	1:B:78:ILE:HD13	2.05	0.55
1:A:178:PRO:HB3	1:A:220:LEU:HD23	1.88	0.55
1:B:96:SER:HB2	1:B:165:GLU:OE1	2.07	0.55
1:A:84:LYS:HD2	1:A:119:GLU:O	2.06	0.55
1:A:240:PHE:O	1:A:244:ILE:HG13	2.06	0.55
1:B:51:ALA:O	1:B:55:LEU:HB2	2.07	0.54
1:B:13:LYS:CA	1:B:64:GLN:HE22	2.21	0.54
1:B:94:GLY:CA	1:B:99:ARG:HH21	2.21	0.54
1:A:207:ILE:HG22	1:A:228:GLY:CA	2.38	0.54
1:B:94:GLY:HA3	1:B:99:ARG:HH21	1.73	0.53
1:B:217:CYS:SG	1:B:243:ILE:HA	2.48	0.53
1:A:2:PRO:HB3	1:A:4:ARG:HH12	1.73	0.53
1:A:34:PRO:HB2	1:A:37:THR:HG23	1.90	0.53
1:A:135:GLU:HG3	2:A:400:HOH:O	2.08	0.53
1:B:91:VAL:CG1	1:B:123:VAL:HG12	2.39	0.53
1:B:182:GLN:OE1	1:B:223:GLN:HB3	2.09	0.53
1:B:90:TRP:HA	1:B:122:GLY:O	2.09	0.53
1:A:184:VAL:HG12	2:A:413:HOH:O	2.08	0.53
1:A:196:VAL:HG11	1:A:200:VAL:HG11	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:ASP:HB2	2:B:366:HOH:O	2.09	0.53
1:A:14:MET:HE1	1:B:69:VAL:HG23	1.90	0.52
1:A:214:GLY:H	1:A:239:GLU:CD	2.13	0.52
1:A:18:LYS:HG2	1:B:85:ASP:OD2	2.10	0.52
1:A:157:TRP:HA	1:A:160:VAL:HG13	1.92	0.52
1:A:127:ILE:CD1	1:A:164:TYR:HA	2.38	0.52
1:A:210:GLY:O	1:A:212:VAL:HG23	2.10	0.52
1:B:198:ASP:HB2	2:B:300:HOH:O	2.10	0.52
1:A:115:HIS:HE1	2:A:270:HOH:O	1.92	0.52
1:B:127:ILE:HD11	1:B:188:LEU:HD21	1.91	0.52
1:B:5:LYS:HB2	2:B:351:HOH:O	2.08	0.51
1:B:16:GLY:HA3	1:B:21:LEU:HD13	1.92	0.51
1:A:174:LYS:NZ	2:A:353:HOH:O	2.43	0.51
1:A:47:TYR:HE2	1:B:82:MET:HG2	1.76	0.51
1:B:63:ALA:O	1:B:92:VAL:HG23	2.11	0.51
1:A:207:ILE:HG12	2:A:315:HOH:O	2.09	0.51
1:A:145:GLU:HG3	2:A:385:HOH:O	2.10	0.51
1:B:235:SER:HB2	2:B:405:HOH:O	2.10	0.51
1:B:71:ASN:HB2	2:B:308:HOH:O	2.11	0.51
1:A:91:VAL:HG23	1:A:122:GLY:O	2.11	0.51
1:B:248:GLN:HE21	1:B:248:GLN:C	2.14	0.50
1:A:231:VAL:CG1	1:A:234:ALA:HB3	2.41	0.50
1:A:58:LYS:HG3	1:A:59:ILE:HG13	1.93	0.50
1:A:185:HIS:ND1	1:A:206:ILE:HG22	2.27	0.50
1:B:6:PHE:O	1:B:38:GLU:N	2.45	0.50
1:B:7:PHE:HD2	1:B:207:ILE:HG21	1.77	0.50
1:B:90:TRP:CE3	1:B:122:GLY:HA3	2.47	0.50
1:A:16:GLY:HA3	1:A:21:LEU:HD13	1.93	0.50
1:A:235:SER:HA	1:A:240:PHE:CD1	2.47	0.49
1:A:56:ASP:OD2	1:A:58:LYS:HG2	2.13	0.49
1:B:164:TYR:HB3	1:B:185:HIS:NE2	2.28	0.49
1:B:49:ASP:O	1:B:53:GLN:HG3	2.12	0.49
1:B:66:CYS:HA	2:B:421:HOH:O	2.13	0.49
1:B:207:ILE:HG22	1:B:228:GLY:HA3	1.95	0.49
1:A:14:MET:CE	1:B:69:VAL:HG23	2.43	0.48
1:A:49:ASP:O	1:A:53:GLN:HG3	2.12	0.48
1:A:42:ALA:HA	1:A:62:ALA:O	2.13	0.48
1:A:155:LYS:HD2	2:A:395:HOH:O	2.13	0.48
1:A:117:LEU:CD1	1:A:154:VAL:HG22	2.42	0.48
1:A:237:LYS:HA	2:A:481:HOH:O	2.13	0.48
1:A:159:LYS:HA	2:A:310:HOH:O	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:LEU:HB3	2:B:460:HOH:O	2.14	0.48
1:B:180:GLN:O	1:B:184:VAL:HG12	2.14	0.48
1:B:95:HIS:HB2	2:B:394:HOH:O	2.14	0.48
1:B:231:VAL:HB	1:B:235:SER:HB3	1.95	0.48
1:A:66:CYS:HB3	1:A:83:ILE:HD11	1.96	0.48
1:A:133:GLU:HB3	2:A:402:HOH:O	2.14	0.48
1:A:84:LYS:CG	1:A:121:LEU:HD13	2.44	0.47
1:B:109:ILE:HG12	2:B:437:HOH:O	2.13	0.47
1:B:42:ALA:HA	1:B:62:ALA:O	2.13	0.47
1:B:145:GLU:O	1:B:149:VAL:HG23	2.14	0.47
1:B:146:GLN:HB3	2:B:444:HOH:O	2.14	0.47
1:B:241:VAL:HG12	1:B:241:VAL:O	2.14	0.47
1:A:4:ARG:NH2	2:A:459:HOH:O	2.48	0.47
1:B:32:LYS:HE2	2:B:333:HOH:O	2.13	0.47
1:A:79:SER:C	2:A:375:HOH:O	2.53	0.47
1:B:202:GLN:HG3	2:B:374:HOH:O	2.13	0.47
1:A:184:VAL:O	1:A:188:LEU:HD13	2.14	0.47
1:B:217:CYS:O	1:B:246:ALA:HB1	2.14	0.47
1:A:141:LYS:O	1:A:145:GLU:HG3	2.15	0.47
1:A:99:ARG:HG2	2:A:253:HOH:O	2.15	0.47
1:A:80:PRO:HB2	1:A:119:GLU:HG3	1.96	0.47
1:B:6:PHE:HB3	1:B:37:THR:HG23	1.96	0.47
1:B:191:TRP:CE3	1:B:192:LEU:HD23	2.50	0.47
1:B:170:ILE:HG12	1:B:210:GLY:HA2	1.97	0.46
1:B:218:LYS:HB2	1:B:218:LYS:NZ	2.30	0.46
1:A:131:LEU:HD12	2:A:440:HOH:O	2.15	0.46
1:B:98:ARG:NH1	1:B:98:ARG:HG3	2.30	0.46
1:B:183:GLU:HG2	1:B:187:LYS:NZ	2.31	0.46
1:A:52:ARG:NH2	1:A:55:LEU:HB3	2.31	0.46
1:A:106:ASP:CG	2:A:356:HOH:O	2.53	0.46
1:B:107:GLU:O	1:B:111:GLN:HG3	2.15	0.46
1:B:64:GLN:O	1:B:65:ASN:HB2	2.16	0.46
1:B:184:VAL:HG13	1:B:185:HIS:H	1.81	0.46
1:A:52:ARG:HA	1:A:61:VAL:HG21	1.98	0.46
1:A:34:PRO:HB3	1:A:248:GLN:NE2	2.31	0.46
1:B:241:VAL:HA	1:B:244:ILE:HD12	1.98	0.46
1:A:161:VAL:HG12	1:A:205:ARG:HH12	1.79	0.45
1:B:182:GLN:HE21	1:B:186:GLU:HB2	1.81	0.45
1:A:52:ARG:NH2	2:A:391:HOH:O	2.48	0.45
1:B:164:TYR:CD1	1:B:184:VAL:HG21	2.52	0.45
1:B:131:LEU:HG	1:B:135:GLU:OE1	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:238:PRO:HD3	2:B:389:HOH:O	2.16	0.44
1:A:138:ILE:HG23	1:A:141:LYS:HD3	1.99	0.44
1:A:12:TRP:HD1	1:A:235:SER:OG	1.97	0.44
1:A:170:ILE:HD11	2:A:401:HOH:O	2.18	0.44
1:A:150:ILE:O	1:A:154:VAL:HG23	2.17	0.44
1:A:177:THR:OG1	1:A:180:GLN:HG2	2.18	0.44
1:A:162:LEU:C	2:A:315:HOH:O	2.56	0.43
1:B:66:CYS:O	1:B:112:LYS:HE2	2.18	0.43
1:B:237:LYS:HD2	1:B:239:GLU:OE1	2.17	0.43
1:B:1:ALA:HB1	2:B:425:HOH:O	2.19	0.43
1:B:57:PRO:HB3	2:B:423:HOH:O	2.17	0.43
1:A:73:ALA:O	1:B:97:GLU:HG2	2.18	0.43
1:B:98:ARG:HH11	1:B:98:ARG:CG	2.30	0.43
1:B:108:LEU:HD11	1:B:112:LYS:CE	2.46	0.43
1:B:39:VAL:HG13	1:B:59:ILE:HG12	1.99	0.43
1:B:182:GLN:HE22	1:B:225:ASP:HB2	1.83	0.43
1:A:13:LYS:HD2	1:B:73:ALA:HA	2.01	0.43
1:A:45:THR:H	1:B:78:ILE:HD11	1.83	0.43
1:A:227:ASP:HB3	2:A:422:HOH:O	2.19	0.43
1:A:137:GLY:HA2	2:A:387:HOH:O	2.19	0.42
1:A:223:GLN:HA	1:A:224:PRO:HD3	1.87	0.42
1:B:18:LYS:O	1:B:22:GLY:N	2.48	0.42
1:B:97:GLU:HA	2:B:361:HOH:O	2.18	0.42
1:A:51:ALA:O	1:A:55:LEU:N	2.52	0.42
1:A:127:ILE:HD11	1:A:188:LEU:HD21	2.01	0.42
1:B:8:VAL:HG13	1:B:39:VAL:HA	2.00	0.42
1:B:99:ARG:HA	1:B:104:GLU:H	1.84	0.42
1:B:11:ASN:OD1	2:B:373:HOH:O	2.21	0.42
1:A:33:VAL:HG21	1:A:39:VAL:HG23	2.02	0.42
1:A:188:LEU:O	1:A:192:LEU:HG	2.20	0.42
1:B:142:VAL:HG23	1:B:143:VAL:N	2.35	0.42
1:B:237:LYS:HB3	2:B:281:HOH:O	2.20	0.42
1:B:242:ASP:O	1:B:245:ASN:N	2.52	0.42
1:A:18:LYS:HD2	1:A:50:PHE:CD1	2.55	0.41
1:A:13:LYS:N	1:A:64:GLN:NE2	2.68	0.41
1:A:13:LYS:N	1:A:64:GLN:HE22	2.17	0.41
1:A:231:VAL:HG11	1:A:234:ALA:HB3	2.03	0.41
1:B:5:LYS:HB2	1:B:5:LYS:HE3	1.86	0.41
1:B:174:LYS:HE3	2:B:407:HOH:O	2.19	0.41
1:B:178:PRO:HD2	1:B:179:GLN:NE2	2.34	0.41
1:B:150:ILE:O	1:B:154:VAL:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:TYR:O	1:B:229:PHE:HA	2.21	0.41
1:A:12:TRP:HB3	1:A:236:LEU:HD21	2.02	0.41
1:B:219:GLU:HB3	2:B:455:HOH:O	2.19	0.41
1:A:8:VAL:CG2	2:A:444:HOH:O	2.59	0.41
1:A:245:ASN:HA	2:A:297:HOH:O	2.20	0.41
1:B:97:GLU:O	1:B:101:VAL:HB	2.20	0.41
1:A:73:ALA:HA	1:B:13:LYS:CD	2.47	0.41
1:A:143:VAL:CG1	1:A:188:LEU:HD11	2.40	0.41
1:B:218:LYS:HG2	2:B:403:HOH:O	2.21	0.41
1:A:13:LYS:CA	1:A:64:GLN:HE22	2.34	0.41
1:B:99:ARG:NH2	1:B:109:ILE:HD13	2.36	0.41
1:B:209:GLY:HA2	1:B:230:LEU:HB3	2.01	0.41
1:B:191:TRP:HE3	1:B:192:LEU:HD23	1.85	0.41
1:A:138:ILE:O	1:A:138:ILE:CG2	2.64	0.40
1:A:52:ARG:CA	1:A:61:VAL:HG21	2.51	0.40
1:A:162:LEU:HD23	1:A:162:LEU:HA	1.86	0.40
1:B:13:LYS:H	1:B:64:GLN:NE2	2.19	0.40
1:A:34:PRO:C	1:A:36:ASP:H	2.24	0.40
1:B:52:ARG:HA	1:B:61:VAL:HG21	2.04	0.40
1:B:78:ILE:CG2	1:B:82:MET:HB2	2.51	0.40
1:B:157:TRP:CE3	1:B:160:VAL:HG11	2.57	0.40
1:B:41:CYS:O	1:B:43:PRO:HD3	2.22	0.40
1:A:64:GLN:O	1:A:65:ASN:HB2	2.21	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	246/250 (98%)	230 (94%)	15 (6%)	1 (0%)	34 37
1	B	246/250 (98%)	206 (84%)	39 (16%)	1 (0%)	34 37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	492/500 (98%)	436 (89%)	54 (11%)	2 (0%)	34	37

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	169	ALA
1	A	2	PRO

### 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	195/196 (100%)	182 (93%)	13 (7%)	16	18
1	B	195/196 (100%)	183 (94%)	12 (6%)	18	21
All	All	390/392 (100%)	365 (94%)	25 (6%)	17	20

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	LEU
1	A	33	VAL
1	A	44	PRO
1	A	55	LEU
1	A	91	VAL
1	A	108	LEU
1	A	140	GLU
1	A	158	SER
1	A	160	VAL
1	A	202	GLN
1	A	205	ARG
1	A	225	ASP
1	A	227	ASP
1	B	4	ARG
1	B	32	LYS
1	B	39	VAL

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Mol	Chain	Res	Type
1	B	56	ASP
1	B	70	THR
1	B	98	ARG
1	B	165	GLU
1	B	179	GLN
1	B	184	VAL
1	B	205	ARG
1	B	227	ASP
1	B	248	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	GLN
1	A	115	HIS
1	A	179	GLN
1	A	248	GLN
1	B	11	ASN
1	B	100	HIS
1	B	115	HIS
1	B	179	GLN
1	B	180	GLN
1	B	223	GLN
1	B	245	ASN
1	B	248	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 monosaccharides 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	248/250 (99%)	0.44	14 (5%) 24 23	21, 33, 51, 84	0
1	B	248/250 (99%)	1.43	56 (22%) 0 0	21, 48, 85, 102	0
All	All	496/500 (99%)	0.94	70 (14%) 2 2	21, 37, 78, 102	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	ALA	13.8
1	B	175	THR	13.3
1	B	173	GLY	10.2
1	A	1	ALA	9.8
1	A	2	PRO	9.8
1	B	172	THR	9.0
1	B	176	ALA	8.9
1	B	170	ILE	7.6
1	B	174	LYS	7.5
1	B	171	GLY	6.6
1	B	199	ALA	5.9
1	B	214	GLY	5.7
1	B	2	PRO	5.2
1	B	222	SER	5.2
1	B	31	ALA	5.0
1	B	169	ALA	4.4
1	B	217	CYS	4.2
1	B	124	ILE	4.1
1	B	163	ALA	4.1
1	B	184	VAL	3.8
1	A	247	LYS	3.8
1	B	220	LEU	3.7
1	B	33	VAL	3.7
1	B	207	ILE	3.5

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	33	VAL	3.4
1	B	230	LEU	3.3
1	B	248	GLN	3.3
1	B	56	ASP	3.3
1	B	3	SER	3.3
1	B	35	ALA	3.2
1	B	131	LEU	3.2
1	B	91	VAL	3.0
1	B	198	ASP	3.0
1	B	234	ALA	3.0
1	B	63	ALA	2.9
1	B	130	LYS	2.8
1	B	221	ALA	2.8
1	A	92	VAL	2.6
1	A	3	SER	2.6
1	B	8	VAL	2.6
1	A	35	ALA	2.6
1	A	230	LEU	2.5
1	B	21	LEU	2.5
1	B	213	THR	2.4
1	B	126	CYS	2.4
1	B	212	VAL	2.4
1	B	140	GLU	2.4
1	B	162	LEU	2.4
1	B	58	LYS	2.3
1	A	40	VAL	2.3
1	B	203	SER	2.3
1	B	161	VAL	2.3
1	B	127	ILE	2.3
1	A	221	ALA	2.3
1	B	40	VAL	2.2
1	B	92	VAL	2.2
1	B	6	PHE	2.2
1	B	202	GLN	2.2
1	B	165	GLU	2.2
1	A	63	ALA	2.1
1	B	192	LEU	2.1
1	B	90	TRP	2.1
1	B	179	GLN	2.1
1	A	172	THR	2.1
1	B	120	GLY	2.1
1	A	199	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	9	GLY	2.1
1	B	188	LEU	2.0
1	B	32	LYS	2.0
1	A	120	GLY	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 monosaccharides 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.