



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:02 PM GMT

PDB ID : 1N8T  
Title : The crystal structure of phosphoglucose isomerase from rabbit muscle  
Authors : Davies, C.; Muirhead, H.  
Deposited on : 2002-11-21  
Resolution : 2.50 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

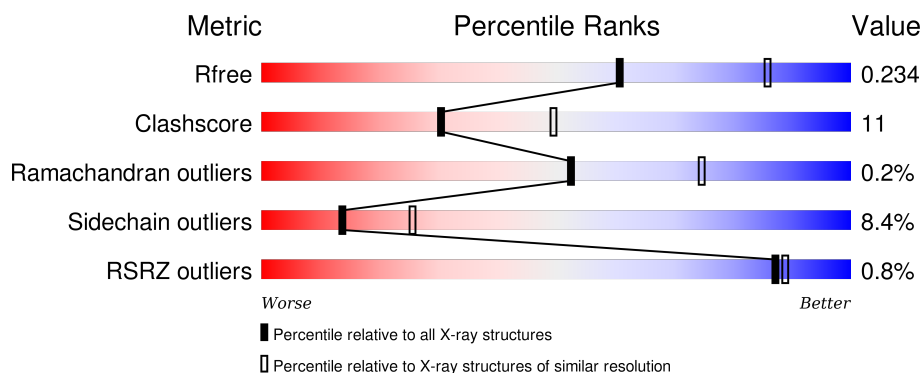
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 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))
$R_{free}$	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (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. 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	557	<div> <div></div> <div>73% 25% .</div> </div>
1	B	557	<div> <div></div> <div>72% 24% .</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9343 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 Glucose-6-phosphate isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	557	Total	C	N	O	S	0	0	0
			4417	2818	768	812	19			
1	B	557	Total	C	N	O	S	0	0	0
			4417	2818	768	812	19			

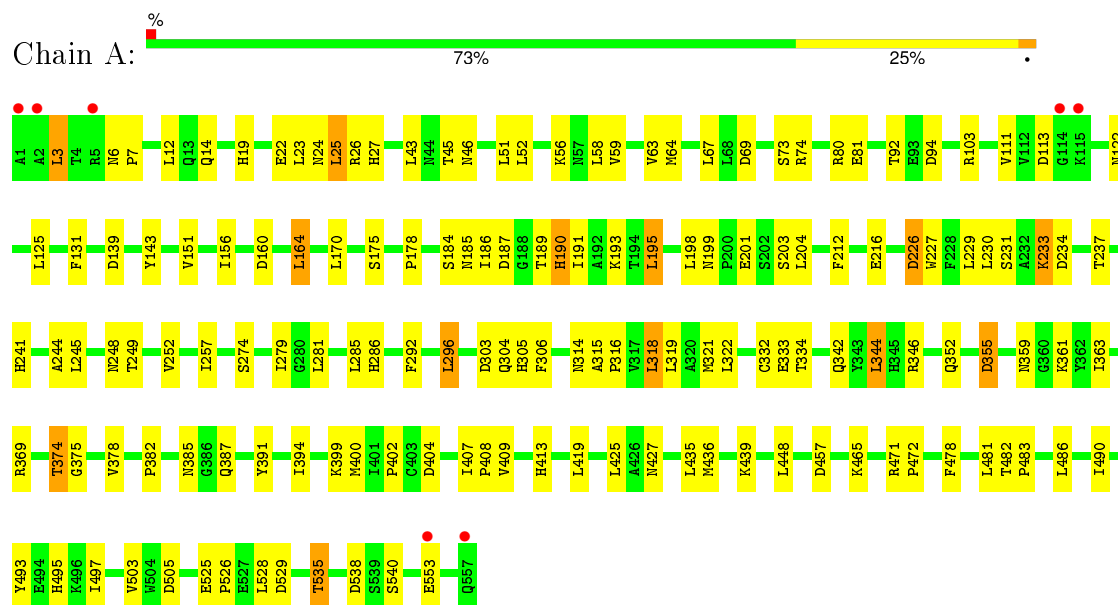
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	249	Total	O	0	0
			249	249		
2	B	260	Total	O	0	0
			260	260		

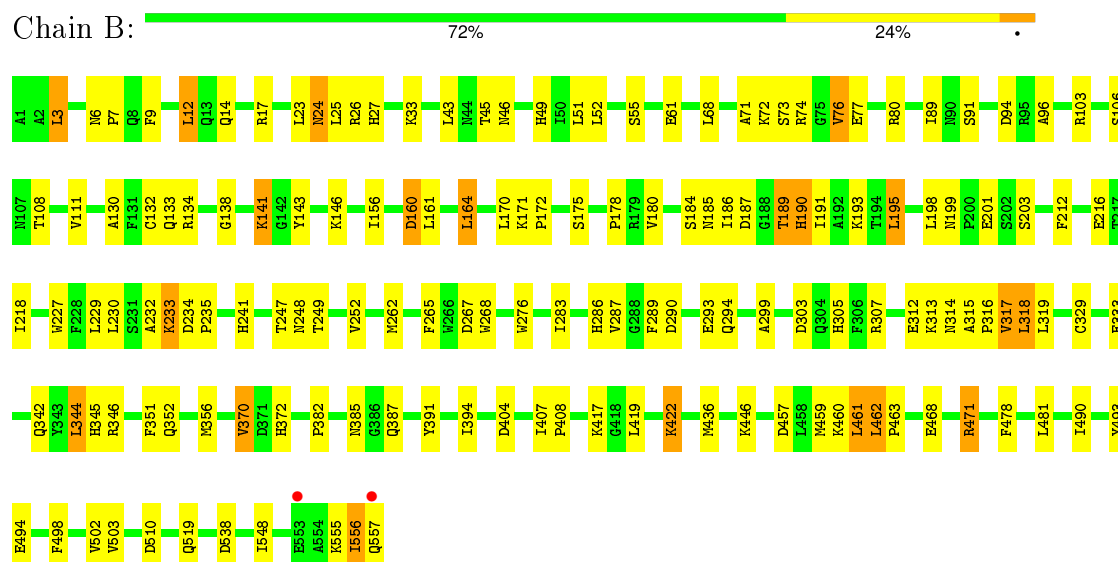
### 3 Residue-property plots

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: Glucose-6-phosphate isomerase



#### • Molecule 1: Glucose-6-phosphate isomerase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.07Å 115.17Å 271.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.50 19.95 – 2.39	Depositor EDS
% Data completeness (in resolution range)	100.0 (15.00-2.50) 91.8 (19.95-2.39)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.33 (at 2.38Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, $R_{free}$	0.178 , 0.234 0.178 , 0.234	Depositor DCC
$R_{free}$ test set	2142 reflections (5.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.0	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 36.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 47432 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9343	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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

### 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	0.51	0/4526	0.81	15/6129 (0.2%)
1	B	0.51	0/4526	0.80	7/6129 (0.1%)
All	All	0.51	0/9052	0.81	22/12258 (0.2%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	164	LEU	CA-CB-CG	6.97	131.33	115.30
1	B	94	ASP	CB-CG-OD2	6.41	124.07	118.30
1	A	69	ASP	CB-CG-OD2	6.21	123.89	118.30
1	B	267	ASP	CB-CG-OD2	5.94	123.65	118.30
1	A	505	ASP	CB-CG-OD2	5.93	123.64	118.30
1	B	404	ASP	CB-CG-OD2	5.83	123.55	118.30
1	A	234	ASP	CB-CG-OD2	5.79	123.51	118.30
1	A	139	ASP	CB-CG-OD2	5.72	123.45	118.30
1	B	510	ASP	CB-CG-OD2	5.67	123.40	118.30
1	B	290	ASP	CB-CG-OD2	5.58	123.32	118.30
1	A	457	ASP	CB-CG-OD2	5.53	123.28	118.30
1	B	160	ASP	CB-CG-OD2	5.49	123.24	118.30
1	A	94	ASP	CB-CG-OD2	5.45	123.20	118.30
1	A	226	ASP	CB-CG-OD1	5.39	123.15	118.30
1	A	160	ASP	CB-CG-OD2	5.28	123.05	118.30
1	A	355	ASP	CB-CG-OD2	5.25	123.03	118.30
1	B	538	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	113	ASP	CB-CG-OD2	5.16	122.95	118.30
1	A	529	ASP	CB-CA-C	-5.14	100.11	110.40
1	A	538	ASP	CB-CG-OD2	5.08	122.87	118.30
1	A	296	LEU	CA-CB-CG	5.05	126.91	115.30
1	A	404	ASP	CB-CG-OD2	5.04	122.83	118.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-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 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	4417	0	4375	106	0
1	B	4417	0	4375	101	0
2	A	249	0	0	6	0
2	B	260	0	0	7	0
All	All	9343	0	8750	193	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 11.

All (193) 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:557:GLN:HB2	2:B:788:HOH:O	1.07	1.19
1:B:557:GLN:CB	2:B:788:HOH:O	1.71	1.00
1:A:103:ARG:HD2	1:A:303:ASP:OD2	1.64	0.97
1:B:103:ARG:HD2	1:B:303:ASP:OD2	1.66	0.96
1:B:232:ALA:O	1:B:234:ASP:N	2.04	0.90
1:A:374:THR:CG2	1:A:375:GLY:H	1.86	0.87
1:B:138:GLY:O	1:B:141:LYS:HE2	1.78	0.82
1:A:143:TYR:HB3	1:A:237:THR:HG23	1.61	0.81
1:B:462:LEU:HB3	1:B:463:PRO:HD3	1.64	0.79
1:B:315:ALA:HB3	1:B:316:PRO:HD3	1.66	0.78
1:A:374:THR:HG23	1:A:375:GLY:H	1.47	0.77
1:A:374:THR:CG2	1:A:375:GLY:N	2.48	0.75
1:B:24:ASN:HD22	1:B:27:HIS:H	1.33	0.75
1:B:185:ASN:H	1:B:190:HIS:HD2	1.33	0.73
1:A:369:ARG:HH11	1:A:369:ARG:HG2	1.54	0.72
1:A:25:LEU:HB3	1:A:436:MET:HG2	1.70	0.72
1:A:143:TYR:H	1:A:241:HIS:HE1	1.38	0.71
1:A:24:ASN:HD22	1:A:27:HIS:H	1.37	0.71
1:A:185:ASN:H	1:A:190:HIS:HD2	1.39	0.70
1:B:314:ASN:HB3	1:B:317:VAL:HG22	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:TYR:H	1:B:241:HIS:HE1	1.41	0.69
1:B:345:HIS:HA	1:B:382:PRO:HG3	1.74	0.69
1:B:191:ILE:HG13	1:B:195:LEU:HD22	1.76	0.68
1:B:164:LEU:HD12	1:B:164:LEU:C	2.15	0.67
1:A:342:GLN:HE21	1:B:185:ASN:HD22	1.41	0.66
1:A:185:ASN:HD22	1:B:342:GLN:HE21	1.41	0.66
1:B:557:GLN:OE1	2:B:788:HOH:O	2.10	0.66
1:A:493:TYR:O	1:A:497:ILE:HG13	1.96	0.66
1:B:77:GLU:OE1	2:B:775:HOH:O	2.14	0.65
1:B:305:HIS:HE1	1:B:315:ALA:H	1.46	0.64
1:A:374:THR:HG22	1:A:375:GLY:N	2.12	0.64
1:A:25:LEU:HD13	1:A:58:LEU:HD23	1.80	0.64
1:B:24:ASN:ND2	1:B:27:HIS:H	1.96	0.63
1:A:314:ASN:OD1	1:A:316:PRO:HD2	1.98	0.63
1:B:248:ASN:O	1:B:252:VAL:HG23	1.99	0.63
1:A:19:HIS:O	1:A:22:GLU:HG2	1.99	0.62
1:A:184:SER:OG	1:B:385:ASN:ND2	2.32	0.62
1:A:3:LEU:HG	1:A:503:VAL:O	2.00	0.62
1:A:382:PRO:O	1:A:385:ASN:HB2	1.99	0.62
1:B:45:THR:O	1:B:46:ASN:HB2	1.98	0.62
1:A:346:ARG:CD	2:A:755:HOH:O	2.48	0.60
1:B:12:LEU:HD13	1:B:329:CYS:SG	2.41	0.60
1:A:363:ILE:HG21	1:B:463:PRO:HB2	1.84	0.60
1:A:143:TYR:HB3	1:A:237:THR:CG2	2.30	0.59
1:A:305:HIS:HE1	1:A:315:ALA:H	1.49	0.59
1:B:422:LYS:HB3	1:B:422:LYS:HZ2	1.68	0.59
1:A:346:ARG:HD2	2:A:755:HOH:O	2.03	0.58
1:B:471:ARG:HG2	2:B:631:HOH:O	2.02	0.58
1:A:45:THR:O	1:A:46:ASN:HB2	2.04	0.58
1:A:374:THR:HG22	1:A:375:GLY:H	1.64	0.57
1:A:315:ALA:HB3	1:A:316:PRO:HD3	1.87	0.57
1:B:156:ILE:HG21	1:B:216:GLU:CG	2.34	0.57
1:B:164:LEU:CD1	1:B:164:LEU:C	2.73	0.57
1:A:274:SER:HB2	1:A:279:ILE:HG13	1.86	0.57
1:A:24:ASN:HD21	1:A:26:ARG:HB3	1.70	0.57
1:B:233:LYS:O	1:B:235:PRO:HD3	2.05	0.57
1:A:59:VAL:HG21	1:A:64:MET:SD	2.45	0.57
1:A:359:ASN:O	1:A:374:THR:CG2	2.52	0.57
1:A:359:ASN:O	1:A:374:THR:HG23	2.05	0.56
1:B:14:GLN:OE1	1:B:17:ARG:NH1	2.36	0.56
1:B:314:ASN:OD1	1:B:316:PRO:HD2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:ILE:HG21	1:B:216:GLU:HG3	1.88	0.55
1:A:425:LEU:HD22	1:B:548:ILE:HD13	1.88	0.55
1:A:203:SER:O	1:A:241:HIS:HD2	1.90	0.55
1:B:394:ILE:O	1:B:471:ARG:HD3	2.06	0.55
1:B:146:LYS:HE3	1:B:201:GLU:OE2	2.08	0.55
1:B:12:LEU:CD1	1:B:329:CYS:SG	2.95	0.54
1:A:361:LYS:H	1:A:374:THR:HG21	1.73	0.54
1:B:186:ILE:HD13	1:B:216:GLU:HA	1.90	0.54
1:A:361:LYS:O	1:A:374:THR:HG21	2.09	0.53
1:A:3:LEU:HD23	1:A:74:ARG:HD3	1.89	0.53
1:B:103:ARG:HD3	1:B:299:ALA:CB	2.39	0.53
1:B:143:TYR:N	1:B:241:HIS:HE1	2.07	0.53
1:A:394:ILE:O	1:A:471:ARG:HD3	2.09	0.53
1:A:103:ARG:CD	1:A:303:ASP:OD2	2.49	0.52
1:B:14:GLN:OE1	1:B:17:ARG:NH2	2.42	0.52
1:A:92:THR:O	1:B:460:LYS:HE3	2.10	0.52
1:B:103:ARG:HD3	1:B:299:ALA:HB1	1.91	0.52
1:A:342:GLN:NE2	1:B:185:ASN:HD22	2.05	0.51
1:A:394:ILE:O	1:A:471:ARG:NH2	2.39	0.51
1:A:387:GLN:HA	1:A:391:TYR:CG	2.45	0.51
1:A:346:ARG:HD3	2:A:755:HOH:O	2.09	0.51
1:B:80:ARG:HD3	1:B:307:ARG:HG3	1.92	0.51
1:B:203:SER:O	1:B:241:HIS:HD2	1.93	0.51
1:A:387:GLN:HA	1:A:391:TYR:CD1	2.46	0.51
1:A:525:GLU:HB2	1:A:526:PRO:HD3	1.91	0.50
1:A:59:VAL:HG23	1:A:63:VAL:CG1	2.41	0.50
1:B:498:PHE:O	1:B:502:VAL:HG23	2.12	0.50
1:A:189:THR:HG21	2:A:734:HOH:O	2.12	0.50
1:B:370:VAL:HG13	1:B:372:HIS:O	2.12	0.50
1:A:24:ASN:HD22	1:A:27:HIS:N	2.09	0.50
1:A:439:LYS:NZ	1:B:519:GLN:HE22	2.08	0.50
1:B:184:SER:H	1:B:190:HIS:CD2	2.30	0.49
1:B:407:ILE:HG13	1:B:408:PRO:CD	2.43	0.49
1:B:96:ALA:HB3	1:B:268:TRP:HB3	1.93	0.49
1:A:185:ASN:HD22	1:B:342:GLN:NE2	2.09	0.49
1:A:274:SER:O	1:A:279:ILE:HB	2.13	0.49
1:B:370:VAL:HG22	1:B:372:HIS:CE1	2.48	0.48
1:A:333:GLU:HB2	1:B:333:GLU:CD	2.33	0.48
1:A:131:PHE:HZ	1:A:204:LEU:HD21	1.77	0.48
1:A:189:THR:O	1:A:193:LYS:HG2	2.13	0.48
1:B:283:ILE:O	1:B:287:VAL:HG22	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:PRO:HA	1:A:478:PHE:O	2.13	0.48
1:A:6:ASN:ND2	1:A:73:SER:OG	2.33	0.48
1:A:59:VAL:HG23	1:A:63:VAL:HG11	1.96	0.48
1:B:143:TYR:H	1:B:241:HIS:CE1	2.26	0.47
1:B:199:ASN:OD1	1:B:201:GLU:HG2	2.13	0.47
1:A:231:SER:O	1:A:233:LYS:HG2	2.13	0.47
1:A:178:PRO:O	1:A:286:HIS:HD2	1.98	0.47
1:A:413:HIS:CD2	1:B:189:THR:HB	2.50	0.47
1:B:407:ILE:HG13	1:B:408:PRO:HD2	1.97	0.47
1:A:191:ILE:HG13	1:A:195:LEU:HD22	1.97	0.46
1:B:556:ILE:HG13	1:B:557:GLN:N	2.30	0.46
1:B:556:ILE:CG1	1:B:557:GLN:N	2.78	0.46
1:A:387:GLN:HE22	1:A:427:ASN:HB3	1.81	0.46
1:B:24:ASN:HD21	1:B:26:ARG:HB3	1.81	0.46
1:A:435:LEU:HB3	1:A:471:ARG:HB2	1.97	0.46
1:B:407:ILE:CG1	1:B:408:PRO:CD	2.94	0.46
1:B:351:PHE:CZ	1:B:493:TYR:HB2	2.50	0.46
1:A:244:ALA:HB2	1:A:257:ILE:HD13	1.98	0.46
1:B:171:LYS:HB3	1:B:172:PRO:HD3	1.98	0.46
1:B:289:PHE:O	1:B:293:GLU:HG3	2.16	0.46
1:A:199:ASN:OD1	1:A:201:GLU:HB2	2.16	0.46
1:B:252:VAL:HG21	1:B:262:MET:SD	2.57	0.45
1:B:178:PRO:O	1:B:286:HIS:HD2	1.98	0.45
1:A:482:THR:HB	1:A:483:PRO:HD2	1.98	0.45
1:B:3:LEU:HD22	1:B:9:PHE:CG	2.50	0.45
1:B:344:LEU:HD22	1:B:481:LEU:HD22	1.97	0.45
1:B:460:LYS:HG3	1:B:461:LEU:N	2.29	0.45
1:A:355:ASP:HA	1:A:497:ILE:HD13	1.98	0.45
1:A:407:ILE:HD13	1:A:425:LEU:HD23	1.97	0.45
1:B:312:GLU:O	1:B:318:LEU:HD21	2.17	0.45
1:B:352:GLN:O	1:B:356:MET:HB2	2.17	0.45
1:B:49:HIS:O	1:B:478:PHE:HA	2.17	0.45
1:A:3:LEU:HD23	1:A:3:LEU:HA	1.74	0.45
1:B:74:ARG:HD3	1:B:503:VAL:O	2.17	0.44
1:B:156:ILE:HG21	1:B:216:GLU:HG2	1.99	0.44
1:A:67:LEU:HB3	1:A:321:MET:HB3	1.99	0.44
1:A:175:SER:HB2	2:A:758:HOH:O	2.17	0.44
1:A:369:ARG:HH11	1:A:369:ARG:CG	2.26	0.44
1:B:387:GLN:HA	1:B:391:TYR:CG	2.53	0.44
1:A:80:ARG:HD2	1:A:81:GLU:OE2	2.18	0.44
1:A:306:PHE:CD2	1:A:495:HIS:CD2	3.06	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:LEU:HA	1:A:230:LEU:HD12	1.77	0.44
1:A:369:ARG:NH1	1:A:369:ARG:HG2	2.28	0.44
1:A:439:LYS:HZ3	1:B:519:GLN:HE22	1.66	0.44
1:A:540:SER:HB2	1:B:55:SER:OG	2.18	0.44
1:A:407:ILE:HG12	1:A:408:PRO:HD2	2.00	0.43
1:A:186:ILE:O	1:A:187:ASP:C	2.56	0.43
1:B:71:ALA:O	1:B:76:VAL:HG13	2.18	0.43
1:A:6:ASN:HA	1:A:7:PRO:HD3	1.84	0.43
1:B:68:LEU:HD13	1:B:312:GLU:HG3	2.00	0.43
1:A:292:PHE:CZ	1:A:296:LEU:HD21	2.53	0.43
1:B:345:HIS:O	1:B:346:ARG:HD2	2.19	0.43
1:B:106:SER:HB2	1:B:108:THR:HG23	2.00	0.43
1:A:245:LEU:HD13	1:A:279:ILE:HA	2.00	0.43
1:A:151:VAL:HG21	1:A:285:LEU:HD13	2.01	0.43
1:B:130:ALA:O	1:B:134:ARG:HG3	2.19	0.43
1:B:193:LYS:NZ	2:B:746:HOH:O	2.51	0.43
1:A:344:LEU:HD13	1:A:481:LEU:HD21	2.00	0.43
1:B:407:ILE:CG1	1:B:408:PRO:HD2	2.49	0.42
1:A:184:SER:H	1:A:190:HIS:CD2	2.38	0.42
1:B:6:ASN:HA	1:B:7:PRO:HD3	1.90	0.42
1:B:187:ASP:OD1	1:B:189:THR:HG23	2.19	0.42
1:A:156:ILE:HG21	1:A:216:GLU:CG	2.49	0.42
1:A:486:LEU:O	1:A:490:ILE:HG13	2.20	0.42
1:A:56:LYS:O	1:A:472:PRO:HA	2.20	0.42
1:A:334:THR:O	1:A:402:PRO:HD2	2.19	0.42
1:B:387:GLN:HA	1:B:391:TYR:CD1	2.54	0.41
1:B:218:ILE:HD13	1:B:218:ILE:HA	1.90	0.41
1:A:409:VAL:HG13	1:B:548:ILE:HG23	2.02	0.41
1:B:490:ILE:O	1:B:494:GLU:HG3	2.21	0.41
1:A:535:THR:CG2	1:A:535:THR:O	2.69	0.41
1:A:332:CYS:SG	1:A:400:MET:HG2	2.61	0.41
1:A:156:ILE:HG21	1:A:216:GLU:HG3	2.03	0.41
1:B:89:ILE:O	1:B:91:SER:N	2.53	0.41
1:B:276:TRP:CZ3	1:B:299:ALA:HB2	2.55	0.41
1:A:25:LEU:HD13	1:A:58:LEU:CD2	2.48	0.41
1:B:195:LEU:HG	1:B:227:TRP:CD2	2.56	0.41
1:A:186:ILE:HD13	1:A:216:GLU:HA	2.02	0.41
1:A:304:GLN:HG2	2:A:700:HOH:O	2.20	0.41
1:A:122:ASN:HD22	1:A:125:LEU:HD12	1.85	0.41
1:A:352:GLN:HG3	1:A:378:VAL:O	2.21	0.41
1:B:195:LEU:HG	1:B:227:TRP:CG	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:557:GLN:HB3	2:B:788:HOH:O	1.70	0.40
1:A:143:TYR:CB	1:A:237:THR:HG23	2.42	0.40
1:B:468:GLU:HB2	1:B:471:ARG:NH1	2.36	0.40
1:B:407:ILE:HA	1:B:408:PRO:HD3	1.87	0.40
1:A:195:LEU:HG	1:A:227:TRP:CD2	2.57	0.40
1:B:230:LEU:HD23	1:B:230:LEU:HA	1.91	0.40
1:A:248:ASN:O	1:A:252:VAL:HG23	2.21	0.40
1:A:448:LEU:HD21	1:A:465:LYS:HE2	2.03	0.40
1:A:318:LEU:HD12	1:A:318:LEU:HA	1.90	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	555/557 (100%)	531 (96%)	24 (4%)	0	100	100
1	B	555/557 (100%)	533 (96%)	20 (4%)	2 (0%)	39	61
All	All	1110/1114 (100%)	1064 (96%)	44 (4%)	2 (0%)	52	75

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	233	LYS
1	B	462	LEU

### 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	478/478 (100%)	448 (94%)	30 (6%)	22	40
1	B	478/478 (100%)	428 (90%)	50 (10%)	8	16
All	All	956/956 (100%)	876 (92%)	80 (8%)	14	25

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	12	LEU
1	A	14	GLN
1	A	23	LEU
1	A	25	LEU
1	A	43	LEU
1	A	51	LEU
1	A	52	LEU
1	A	111	VAL
1	A	164	LEU
1	A	170	LEU
1	A	190	HIS
1	A	195	LEU
1	A	198	LEU
1	A	212	PHE
1	A	226	ASP
1	A	229	LEU
1	A	233	LYS
1	A	249	THR
1	A	281	LEU
1	A	318	LEU
1	A	319	LEU
1	A	322	LEU
1	A	344	LEU
1	A	374	THR
1	A	399	LYS
1	A	419	LEU
1	A	528	LEU
1	A	535	THR
1	A	553	GLU
1	B	3	LEU
1	B	12	LEU
1	B	23	LEU

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Mol	Chain	Res	Type
1	B	24	ASN
1	B	25	LEU
1	B	33	LYS
1	B	43	LEU
1	B	51	LEU
1	B	52	LEU
1	B	61	GLU
1	B	72	LYS
1	B	73	SER
1	B	76	VAL
1	B	111	VAL
1	B	132	CYS
1	B	133	GLN
1	B	141	LYS
1	B	160	ASP
1	B	161	LEU
1	B	164	LEU
1	B	170	LEU
1	B	175	SER
1	B	180	VAL
1	B	189	THR
1	B	190	HIS
1	B	195	LEU
1	B	198	LEU
1	B	212	PHE
1	B	229	LEU
1	B	247	THR
1	B	249	THR
1	B	265	PHE
1	B	294	GLN
1	B	313	LYS
1	B	317	VAL
1	B	318	LEU
1	B	319	LEU
1	B	344	LEU
1	B	370	VAL
1	B	417	LYS
1	B	419	LEU
1	B	422	LYS
1	B	436	MET
1	B	446	LYS
1	B	457	ASP

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Mol	Chain	Res	Type
1	B	459	MET
1	B	461	LEU
1	B	471	ARG
1	B	555	LYS
1	B	556	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	24	ASN
1	A	57	ASN
1	A	107	ASN
1	A	122	ASN
1	A	133	GLN
1	A	190	HIS
1	A	241	HIS
1	A	286	HIS
1	A	291	ASN
1	A	305	HIS
1	A	342	GLN
1	A	359	ASN
1	A	385	ASN
1	A	387	GLN
1	A	474	ASN
1	A	500	GLN
1	A	507	ASN
1	A	551	GLN
1	B	19	HIS
1	B	24	ASN
1	B	57	ASN
1	B	107	ASN
1	B	122	ASN
1	B	190	HIS
1	B	241	HIS
1	B	286	HIS
1	B	291	ASN
1	B	305	HIS
1	B	342	GLN
1	B	352	GLN
1	B	353	GLN
1	B	359	ASN

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Mol	Chain	Res	Type
1	B	385	ASN
1	B	387	GLN
1	B	474	ASN
1	B	500	GLN
1	B	507	ASN
1	B	519	GLN
1	B	551	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 [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	557/557 (100%)	-0.49	7 (1%) 79 82	15, 23, 35, 55	0
1	B	557/557 (100%)	-0.55	2 (0%) 93 93	14, 22, 34, 62	0
All	All	1114/1114 (100%)	-0.52	9 (0%) 87 89	14, 23, 35, 62	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	557	GLN	8.2
1	A	2	ALA	3.1
1	A	1	ALA	2.9
1	A	557	GLN	2.6
1	A	553	GLU	2.6
1	A	114	GLY	2.3
1	A	5	ARG	2.2
1	A	115	LYS	2.2
1	B	553	GLU	2.1

### 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.