



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:09 PM GMT

PDB ID : 1NN4
Title : Structural Genomics, RpiB/AlsB
Authors : Zhang, R.G.; Andersson, C.E.; Mowbray, S.L.; Savchenko, A.; Skarina, T.;
Evdokimova, E.; Beasley, S.L.; Arrowsmith, C.; Edwards, A.M.; Joachimiak,
A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2003-01-12
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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

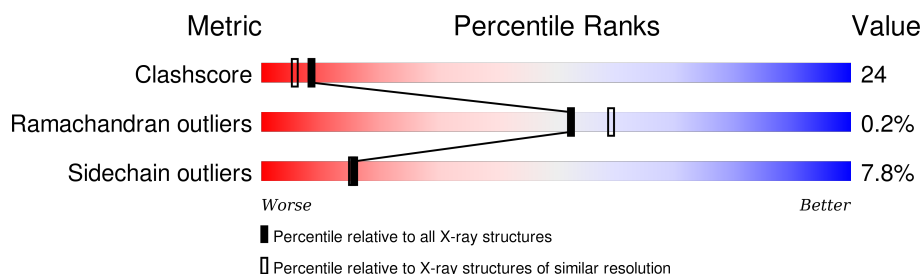
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	162	 63% 29% 6% •
1	B	162	 65% 28% • •
1	C	162	 54% 41% • •
1	D	162	 59% 34% 6% •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5061 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 Ribose 5-phosphate isomerase B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	159	Total	C	N	O	S	0	0	0
			1194	748	219	222	5			
1	B	158	Total	C	N	O	S	0	0	0
			1188	744	217	222	5			
1	C	159	Total	C	N	O	S	0	0	0
			1190	746	219	220	5			
1	D	159	Total	C	N	O	S	0	0	0
			1191	746	219	221	5			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	HIS	-	CLONING ARTIFACT	UNP P37351
A	-11	HIS	-	CLONING ARTIFACT	UNP P37351
A	-10	HIS	-	CLONING ARTIFACT	UNP P37351
A	-9	SER	-	CLONING ARTIFACT	UNP P37351
A	-8	SER	-	CLONING ARTIFACT	UNP P37351
A	-7	GLY	-	CLONING ARTIFACT	UNP P37351
A	-6	LEU	-	CLONING ARTIFACT	UNP P37351
A	-5	THR	-	CLONING ARTIFACT	UNP P37351
A	-4	PRO	-	CLONING ARTIFACT	UNP P37351
A	-3	ARG	-	CLONING ARTIFACT	UNP P37351
A	-2	GLY	-	CLONING ARTIFACT	UNP P37351
A	-1	SER	-	CLONING ARTIFACT	UNP P37351
A	0	GLN	-	CLONING ARTIFACT	UNP P37351
B	-12	HIS	-	CLONING ARTIFACT	UNP P37351
B	-11	HIS	-	CLONING ARTIFACT	UNP P37351
B	-10	HIS	-	CLONING ARTIFACT	UNP P37351
B	-9	SER	-	CLONING ARTIFACT	UNP P37351
B	-8	SER	-	CLONING ARTIFACT	UNP P37351
B	-7	GLY	-	CLONING ARTIFACT	UNP P37351
B	-6	LEU	-	CLONING ARTIFACT	UNP P37351
B	-5	THR	-	CLONING ARTIFACT	UNP P37351

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	PRO	-	CLONING ARTIFACT	UNP P37351
B	-3	ARG	-	CLONING ARTIFACT	UNP P37351
B	-2	GLY	-	CLONING ARTIFACT	UNP P37351
B	-1	SER	-	CLONING ARTIFACT	UNP P37351
B	0	GLN	-	CLONING ARTIFACT	UNP P37351
C	-12	HIS	-	CLONING ARTIFACT	UNP P37351
C	-11	HIS	-	CLONING ARTIFACT	UNP P37351
C	-10	HIS	-	CLONING ARTIFACT	UNP P37351
C	-9	SER	-	CLONING ARTIFACT	UNP P37351
C	-8	SER	-	CLONING ARTIFACT	UNP P37351
C	-7	GLY	-	CLONING ARTIFACT	UNP P37351
C	-6	LEU	-	CLONING ARTIFACT	UNP P37351
C	-5	THR	-	CLONING ARTIFACT	UNP P37351
C	-4	PRO	-	CLONING ARTIFACT	UNP P37351
C	-3	ARG	-	CLONING ARTIFACT	UNP P37351
C	-2	GLY	-	CLONING ARTIFACT	UNP P37351
C	-1	SER	-	CLONING ARTIFACT	UNP P37351
C	0	GLN	-	CLONING ARTIFACT	UNP P37351
D	-12	HIS	-	CLONING ARTIFACT	UNP P37351
D	-11	HIS	-	CLONING ARTIFACT	UNP P37351
D	-10	HIS	-	CLONING ARTIFACT	UNP P37351
D	-9	SER	-	CLONING ARTIFACT	UNP P37351
D	-8	SER	-	CLONING ARTIFACT	UNP P37351
D	-7	GLY	-	CLONING ARTIFACT	UNP P37351
D	-6	LEU	-	CLONING ARTIFACT	UNP P37351
D	-5	THR	-	CLONING ARTIFACT	UNP P37351
D	-4	PRO	-	CLONING ARTIFACT	UNP P37351
D	-3	ARG	-	CLONING ARTIFACT	UNP P37351
D	-2	GLY	-	CLONING ARTIFACT	UNP P37351
D	-1	SER	-	CLONING ARTIFACT	UNP P37351
D	0	GLN	-	CLONING ARTIFACT	UNP P37351

- Molecule 2 is water.

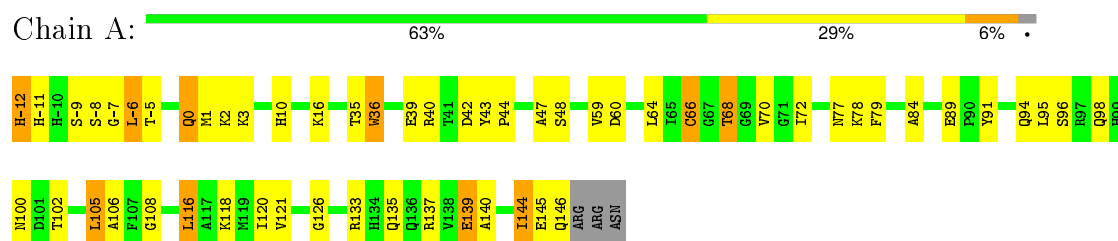
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	92	Total O 92 92	0	0
2	B	66	Total O 66 66	0	0
2	C	66	Total O 66 66	0	0
2	D	74	Total O 74 74	0	0

3 Residue-property plots [i](#)

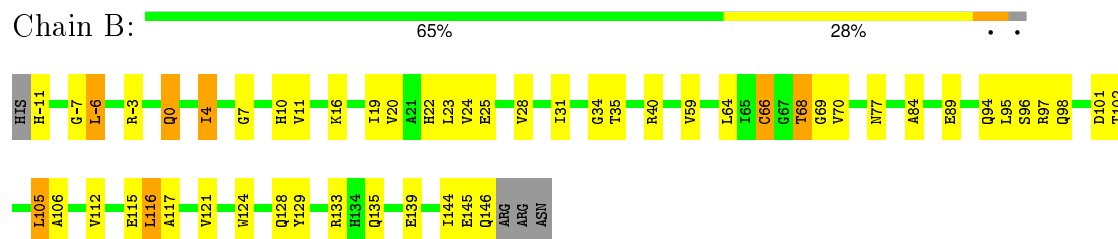
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

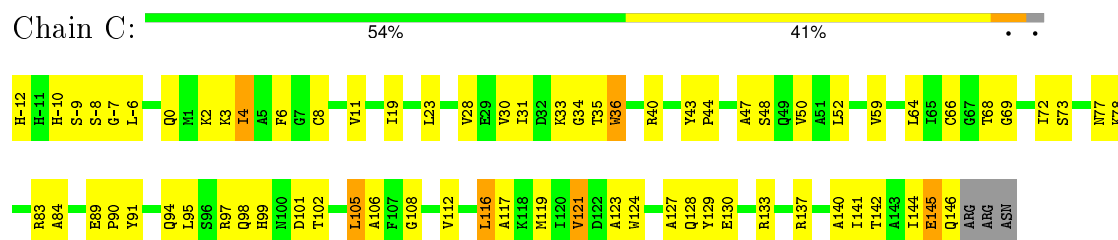
• Molecule 1: Ribose 5-phosphate isomerase B



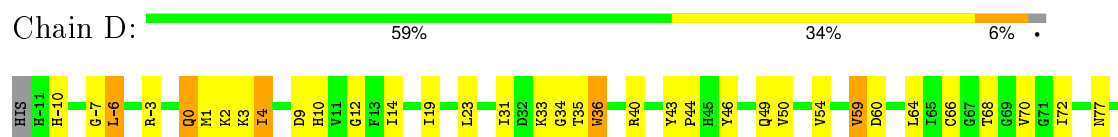
• Molecule 1: Ribose 5-phosphate isomerase B



• Molecule 1: Ribose 5-phosphate isomerase B



• Molecule 1: Ribose 5-phosphate isomerase B



R83	A84	E89	P90	Y91	Q94	L95	S96	R97	Q98	D101	V104	L105	A106	R110	V111	V112	E115	L116	A117	K118	V121	A127	Q128	Y129	R137	A140	L141	T142	A143	I144	E145	Q146	R147	ARG	ASN
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4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	145.59 Å 145.59 Å 74.78 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.40 – 2.20	Depositor
% Data completeness (in resolution range)	89.6 (36.40-2.20)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.216 , 0.262	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5061	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

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	0.48	1/1217 (0.1%)	0.68	1/1646 (0.1%)
1	B	0.37	0/1210	0.67	1/1636 (0.1%)
1	C	0.48	1/1213 (0.1%)	0.65	0/1641
1	D	0.47	1/1213 (0.1%)	0.64	0/1640
All	All	0.45	3/4853 (0.1%)	0.66	2/6563 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	36	TRP	NE1-CE2	8.81	1.49	1.37
1	D	36	TRP	NE1-CE2	8.75	1.49	1.37
1	A	36	TRP	NE1-CE2	8.69	1.48	1.37

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	CYS	CA-CB-SG	5.78	124.41	114.00
1	B	66	CYS	CA-CB-SG	5.27	123.49	114.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	133	ARG	Sidechain
1	B	133	ARG	Sidechain
1	C	133	ARG	Sidechain

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	1194	0	1178	61	0
1	B	1188	0	1177	57	0
1	C	1190	0	1174	75	0
1	D	1191	0	1178	71	0
2	A	92	0	0	6	0
2	B	66	0	0	1	0
2	C	66	0	0	1	0
2	D	74	0	0	5	0
All	All	5061	0	4707	227	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 24.

All (227) 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:C:35:THR:HG22	1:C:36:TRP:H	1.20	1.01
1:A:144:ILE:O	1:A:146:GLN:N	1.95	0.99
1:C:83:ARG:HH12	1:C:101:ASP:HB3	1.32	0.93
1:C:77:ASN:HD21	1:C:84:ALA:H	1.13	0.90
1:D:77:ASN:HD21	1:D:84:ALA:H	1.16	0.90
1:B:68:THR:HG22	1:B:70:VAL:H	1.37	0.89
1:A:68:THR:HG23	1:A:70:VAL:H	1.39	0.88
1:A:144:ILE:C	1:A:146:GLN:H	1.77	0.88
1:C:35:THR:HG22	1:C:36:TRP:N	1.86	0.87
1:D:68:THR:HG23	1:D:70:VAL:H	1.38	0.86
1:C:102:THR:HG23	1:C:124:TRP:HE1	1.42	0.84
1:B:77:ASN:HD21	1:B:84:ALA:H	1.23	0.83
1:C:43:TYR:HB2	1:C:72:ILE:HG23	1.59	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:CYS:HB3	1:B:69:GLY:H	1.44	0.82
1:C:-7:GLY:H	1:D:94:GLN:HE21	1.26	0.81
1:C:35:THR:CG2	1:C:36:TRP:H	1.95	0.79
1:A:-7:GLY:H	1:B:94:GLN:HE21	1.29	0.79
1:B:66:CYS:HB3	1:B:69:GLY:N	1.98	0.78
1:A:77:ASN:HD21	1:A:84:ALA:H	1.30	0.78
1:B:102:THR:HG23	1:B:124:TRP:HE1	1.48	0.76
1:D:77:ASN:ND2	1:D:84:ALA:H	1.84	0.76
1:C:77:ASN:ND2	1:C:84:ALA:H	1.85	0.75
1:A:68:THR:CG2	1:A:70:VAL:H	2.01	0.73
1:B:4:ILE:HD13	1:B:4:ILE:H	1.54	0.73
1:A:94:GLN:NE2	1:B:-6:LEU:H	1.87	0.72
1:B:77:ASN:ND2	1:B:84:ALA:H	1.88	0.71
1:C:94:GLN:HE21	1:D:-7:GLY:H	1.37	0.70
2:A:435:HOH:O	1:D:68:THR:HG21	1.91	0.70
1:C:83:ARG:NH1	1:C:101:ASP:HB3	2.05	0.70
1:A:77:ASN:ND2	1:A:84:ALA:H	1.88	0.70
1:D:142:THR:HG23	2:D:342:HOH:O	1.92	0.70
1:D:112:VAL:HG22	1:D:116:LEU:HB3	1.73	0.69
1:C:102:THR:CG2	1:C:124:TRP:HE1	2.07	0.67
1:D:142:THR:O	1:D:145:GLU:HG2	1.94	0.67
1:B:68:THR:CG2	1:B:70:VAL:H	2.07	0.66
1:B:23:LEU:HB3	1:B:28:VAL:CG2	2.25	0.66
1:A:135:GLN:O	1:A:139:GLU:HB2	1.96	0.66
1:B:102:THR:CG2	1:B:124:TRP:HE1	2.08	0.65
1:A:-7:GLY:H	1:B:94:GLN:NE2	1.94	0.65
1:C:94:GLN:NE2	1:D:-6:LEU:H	1.95	0.65
1:C:117:ALA:O	1:C:121:VAL:HG13	1.97	0.65
1:A:-6:LEU:H	1:B:94:GLN:NE2	1.94	0.65
1:C:94:GLN:O	1:C:98:GLN:HG2	1.97	0.65
1:C:72:ILE:HD12	1:C:73:SER:N	2.12	0.64
1:D:43:TYR:HB2	1:D:72:ILE:HG23	1.78	0.64
1:C:142:THR:O	1:C:145:GLU:HG3	1.97	0.63
1:D:137:ARG:HG3	1:D:137:ARG:HH11	1.63	0.63
1:D:140:ALA:O	1:D:144:ILE:HG23	1.99	0.63
1:C:123:ALA:HB2	1:D:-6:LEU:HD22	1.82	0.62
1:C:3:LYS:CD	1:C:59:VAL:HG12	2.29	0.62
1:C:48:SER:O	1:C:52:LEU:HG	1.99	0.62
1:A:0:GLN:HB2	2:A:315:HOH:O	1.99	0.61
1:A:35:THR:OG1	1:A:36:TRP:N	2.34	0.60
1:C:72:ILE:HD12	1:C:73:SER:H	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:0:GLN:OE1	1:C:128:GLN:HG2	2.02	0.60
1:C:119:MET:HE1	1:D:91:TYR:N	2.15	0.60
1:C:119:MET:HE1	1:D:91:TYR:H	1.66	0.60
1:D:4:ILE:HD11	1:D:23:LEU:HD13	1.82	0.60
1:B:145:GLU:HG2	1:C:78:LYS:HD2	1.83	0.60
1:A:-8:SER:O	1:B:-3:ARG:HD2	2.01	0.60
1:C:97:ARG:HG2	1:C:127:ALA:HB3	1.84	0.60
1:A:116:LEU:O	1:A:120:ILE:HG13	2.02	0.60
1:A:47:ALA:HB2	1:A:72:ILE:HG13	1.84	0.59
1:C:90:PRO:HG2	1:C:119:MET:HE3	1.83	0.59
1:A:118:LYS:HA	1:A:121:VAL:HG22	1.84	0.59
1:B:96:SER:OG	1:B:102:THR:HG21	2.03	0.59
1:D:64:LEU:HB3	1:D:72:ILE:HD12	1.84	0.58
1:B:11:VAL:HG21	2:B:464:HOH:O	2.02	0.58
1:B:117:ALA:O	1:B:121:VAL:HG23	2.03	0.58
1:D:112:VAL:HG23	1:D:116:LEU:HD13	1.85	0.58
1:C:112:VAL:HG22	1:C:116:LEU:HB3	1.86	0.58
1:A:-7:GLY:N	1:B:94:GLN:HE21	2.01	0.58
1:C:90:PRO:HG2	1:C:119:MET:CE	2.34	0.57
1:A:68:THR:CG2	1:A:70:VAL:HB	2.34	0.57
1:D:97:ARG:HG2	1:D:127:ALA:HB3	1.86	0.57
1:D:31:ILE:HG22	1:D:33:LYS:HE2	1.85	0.57
1:C:140:ALA:O	1:C:144:ILE:HG12	2.04	0.57
1:C:98:GLN:HB2	1:C:130:GLU:HG3	1.86	0.57
1:A:105:LEU:HD22	1:A:106:ALA:N	2.19	0.57
1:A:16:LYS:NZ	1:A:35:THR:O	2.27	0.57
1:C:94:GLN:NE2	1:D:-7:GLY:H	2.02	0.57
1:B:7:GLY:O	1:B:64:LEU:HA	2.04	0.57
1:C:40:ARG:HG3	1:C:40:ARG:HH11	1.69	0.57
1:D:101:ASP:HB2	1:D:129:TYR:CD1	2.40	0.57
1:D:35:THR:OG1	1:D:36:TRP:N	2.38	0.56
1:A:144:ILE:C	1:A:146:GLN:N	2.44	0.56
1:A:94:GLN:HE21	1:B:-7:GLY:H	1.52	0.56
1:C:-7:GLY:N	1:D:94:GLN:HE21	2.01	0.55
1:B:68:THR:HG23	1:B:70:VAL:HG23	1.89	0.55
1:C:105:LEU:HD22	1:C:106:ALA:N	2.23	0.54
1:C:3:LYS:HD2	1:C:59:VAL:HG12	1.89	0.54
1:D:14:ILE:HD11	1:D:110:ARG:NH2	2.23	0.54
1:D:31:ILE:CG2	1:D:33:LYS:HE2	2.38	0.54
1:D:112:VAL:CG2	1:D:116:LEU:HB3	2.37	0.54
1:B:31:ILE:HD12	1:B:59:VAL:CG1	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:112:VAL:CG2	1:C:116:LEU:HD13	2.38	0.53
1:A:48:SER:HG	1:A:79:PHE:HZ	1.56	0.53
1:A:10:HIS:CG	1:A:40:ARG:HA	2.43	0.53
1:C:0:GLN:HG2	1:C:0:GLN:O	2.07	0.53
1:D:66:CYS:HB3	1:D:72:ILE:HD13	1.90	0.53
1:B:112:VAL:HG13	1:B:116:LEU:HB3	1.89	0.53
1:A:3:LYS:O	1:A:59:VAL:HG22	2.09	0.53
1:C:69:GLY:HA2	1:C:72:ILE:HD11	1.91	0.52
1:C:-10:HIS:HB3	1:D:-3:ARG:NH1	2.23	0.52
1:A:3:LYS:HG3	1:A:59:VAL:HG23	1.90	0.52
1:C:43:TYR:N	1:C:44:PRO:CD	2.73	0.52
1:C:31:ILE:N	1:C:31:ILE:HD12	2.24	0.52
1:C:47:ALA:HB2	1:C:72:ILE:HG22	1.91	0.52
1:A:2:LYS:NZ	2:A:465:HOH:O	2.43	0.52
1:D:50:VAL:HG21	1:D:64:LEU:HD21	1.92	0.51
1:A:68:THR:HG21	1:A:70:VAL:HB	1.91	0.51
1:C:102:THR:HG23	1:C:124:TRP:NE1	2.21	0.51
1:B:4:ILE:HD13	1:B:4:ILE:N	2.25	0.51
1:C:40:ARG:NH1	2:C:345:HOH:O	2.42	0.51
1:B:102:THR:HG23	1:B:124:TRP:NE1	2.22	0.51
1:C:112:VAL:HG23	1:C:116:LEU:HD13	1.93	0.51
1:A:91:TYR:OH	1:B:115:GLU:HG3	2.12	0.50
2:A:380:HOH:O	1:B:-11:HIS:HE1	1.94	0.50
1:A:94:GLN:NE2	1:B:-7:GLY:H	2.09	0.50
1:B:66:CYS:HB3	1:B:69:GLY:CA	2.41	0.50
1:C:145:GLU:OE1	1:C:146:GLN:N	2.45	0.50
1:A:78:LYS:HB2	1:D:144:ILE:HD11	1.94	0.50
1:C:94:GLN:HE21	1:D:-7:GLY:N	2.07	0.49
1:A:-6:LEU:HD23	1:B:-6:LEU:HD23	1.94	0.49
1:A:-11:HIS:O	2:A:322:HOH:O	2.20	0.49
1:B:68:THR:HG22	1:B:70:VAL:N	2.18	0.49
1:D:68:THR:HG23	1:D:70:VAL:N	2.17	0.49
1:B:20:VAL:O	1:B:24:VAL:HG23	2.12	0.49
1:A:79:PHE:CE1	1:D:144:ILE:HD12	2.47	0.49
1:B:97:ARG:HA	1:B:102:THR:HG22	1.94	0.48
1:D:101:ASP:HB2	1:D:129:TYR:CG	2.48	0.48
1:A:68:THR:HG21	2:D:390:HOH:O	2.13	0.48
1:B:10:HIS:CG	1:B:40:ARG:HA	2.48	0.48
1:A:-9:SER:HB2	1:B:128:GLN:O	2.14	0.48
1:A:-12:HIS:HA	2:A:516:HOH:O	2.13	0.47
1:A:43:TYR:CG	1:A:44:PRO:HD3	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:84:ALA:HB2	1:D:104:VAL:HB	1.95	0.47
1:D:3:LYS:HB3	1:D:3:LYS:NZ	2.29	0.47
1:C:8:CYS:N	1:C:34:GLY:O	2.47	0.47
1:A:-6:LEU:H	1:B:94:GLN:HE22	1.62	0.47
1:C:4:ILE:HD11	1:C:30:VAL:HG22	1.96	0.47
1:C:4:ILE:HD13	1:C:4:ILE:H	1.79	0.47
1:B:101:ASP:HB2	1:B:129:TYR:CD1	2.49	0.47
1:D:9:ASP:HB2	1:D:66:CYS:HA	1.97	0.46
1:A:1:MET:HE2	1:A:60:ASP:O	2.16	0.46
1:D:50:VAL:O	1:D:54:VAL:HG23	2.16	0.46
1:A:126:GLY:C	1:B:-7:GLY:HA2	2.36	0.46
1:A:-7:GLY:N	1:B:94:GLN:NE2	2.62	0.46
1:C:69:GLY:O	1:C:72:ILE:HD12	2.15	0.46
1:B:135:GLN:O	1:B:139:GLU:HB2	2.15	0.46
1:A:91:TYR:CZ	1:B:115:GLU:HB3	2.50	0.46
1:D:33:LYS:HD2	1:D:49:GLN:HB3	1.98	0.46
1:C:68:THR:HA	1:C:108:GLY:HA3	1.98	0.46
1:D:9:ASP:HB3	1:D:12:GLY:H	1.81	0.46
1:C:-8:SER:HB2	1:D:94:GLN:NE2	2.31	0.46
1:B:101:ASP:HB2	1:B:129:TYR:CG	2.50	0.45
1:A:94:GLN:HE21	1:B:-6:LEU:H	1.63	0.45
1:D:142:THR:HG22	1:D:145:GLU:OE1	2.16	0.45
1:D:10:HIS:CG	1:D:40:ARG:HA	2.51	0.45
1:A:0:GLN:HB2	1:A:0:GLN:HE21	1.58	0.45
1:D:-3:ARG:NH1	2:D:426:HOH:O	2.49	0.45
1:D:95:LEU:HA	1:D:98:GLN:HE21	1.81	0.45
1:C:137:ARG:O	1:C:141:ILE:HG12	2.15	0.45
1:D:83:ARG:HH11	1:D:83:ARG:HG3	1.81	0.45
1:D:3:LYS:HG2	1:D:59:VAL:HG23	1.98	0.45
1:A:144:ILE:O	1:A:146:GLN:HG2	2.17	0.44
1:C:31:ILE:CD1	1:C:31:ILE:N	2.81	0.44
1:C:-6:LEU:H	1:D:94:GLN:NE2	2.16	0.44
1:A:47:ALA:HB2	1:A:72:ILE:CG1	2.47	0.44
1:D:118:LYS:O	1:D:121:VAL:HG22	2.17	0.44
1:C:6:PHE:O	1:C:33:LYS:HB2	2.17	0.44
1:C:3:LYS:HD2	1:C:59:VAL:HA	1.99	0.44
1:C:43:TYR:HB2	1:C:72:ILE:CG2	2.38	0.44
1:C:66:CYS:SG	1:C:69:GLY:N	2.86	0.44
1:D:49:GLN:NE2	2:D:343:HOH:O	2.50	0.44
1:C:94:GLN:HE22	1:D:-6:LEU:H	1.65	0.44
1:C:119:MET:HE3	1:D:90:PRO:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:137:ARG:NH1	1:D:137:ARG:HG3	2.29	0.44
1:B:16:LYS:O	1:B:20:VAL:HG23	2.17	0.44
1:C:95:LEU:HD11	1:C:99:HIS:CE1	2.53	0.43
1:C:-7:GLY:H	1:D:94:GLN:NE2	2.05	0.43
1:C:97:ARG:HG3	1:C:97:ARG:NH1	2.33	0.43
1:D:95:LEU:HD12	1:D:98:GLN:HE21	1.83	0.43
1:A:66:CYS:O	1:A:108:GLY:HA2	2.18	0.43
1:B:77:ASN:HD21	1:B:84:ALA:N	2.04	0.43
1:D:43:TYR:N	1:D:44:PRO:CD	2.81	0.43
1:C:50:VAL:HB	1:C:64:LEU:HD21	1.99	0.43
1:B:64:LEU:O	1:B:106:ALA:HA	2.19	0.43
1:A:43:TYR:N	1:A:44:PRO:CD	2.80	0.43
1:D:64:LEU:O	1:D:106:ALA:HA	2.19	0.43
1:D:33:LYS:O	1:D:46:TYR:HD2	2.02	0.43
1:D:0:GLN:HB3	1:D:0:GLN:HE21	1.61	0.43
1:A:42:ASP:HB2	1:A:44:PRO:HD2	2.01	0.42
1:D:1:MET:HG2	1:D:60:ASP:HB3	2.01	0.42
1:C:66:CYS:O	1:C:108:GLY:HA2	2.18	0.42
1:C:98:GLN:HA	1:C:129:TYR:HA	2.00	0.42
1:A:96:SER:HB2	1:A:102:THR:HG21	2.00	0.42
1:B:95:LEU:HA	1:B:98:GLN:HE21	1.84	0.42
1:D:118:LYS:HA	1:D:121:VAL:HG22	2.01	0.42
1:A:68:THR:HG23	1:A:70:VAL:HG23	2.01	0.42
1:D:2:LYS:HD3	1:D:2:LYS:HA	1.90	0.42
1:C:91:TYR:CZ	1:D:115:GLU:HB3	2.54	0.42
1:B:66:CYS:HB3	1:B:69:GLY:HA2	2.02	0.42
1:B:105:LEU:CD2	1:B:106:ALA:N	2.83	0.42
1:B:22:HIS:O	1:B:25:GLU:HB3	2.20	0.42
1:A:10:HIS:HB3	1:A:39:GLU:O	2.20	0.42
1:D:146:GLN:OE1	1:D:147:ARG:HB2	2.20	0.42
1:A:95:LEU:HD12	1:A:98:GLN:HE21	1.84	0.41
1:A:-5:THR:OG1	1:B:-3:ARG:NH1	2.53	0.41
1:D:34:GLY:HA3	1:D:36:TRP:CH2	2.55	0.41
1:D:9:ASP:CB	1:D:66:CYS:HA	2.51	0.41
1:A:137:ARG:O	1:A:140:ALA:HB3	2.19	0.41
1:C:89:GLU:OE1	1:C:91:TYR:HB3	2.21	0.41
1:C:105:LEU:CD2	1:C:106:ALA:N	2.84	0.41
1:C:-9:SER:O	1:D:127:ALA:HA	2.19	0.41
1:C:40:ARG:HG3	1:C:40:ARG:NH1	2.34	0.41
1:C:23:LEU:HB3	1:C:28:VAL:CG2	2.50	0.41
1:D:2:LYS:HE2	2:D:514:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:ILE:C	1:B:146:GLN:H	2.24	0.41
1:A:68:THR:CG2	1:A:70:VAL:N	2.77	0.40
1:B:105:LEU:HD23	1:B:106:ALA:H	1.86	0.40
1:A:64:LEU:O	1:A:106:ALA:HA	2.21	0.40
1:D:128:GLN:HB2	1:D:128:GLN:HE21	1.65	0.40
1:B:0:GLN:HG3	1:B:128:GLN:CG	2.51	0.40
1:A:96:SER:O	1:A:100:ASN:HB2	2.21	0.40
1:A:95:LEU:HD12	1:A:98:GLN:NE2	2.35	0.40
1:B:34:GLY:HA3	1:B:35:THR:HA	1.91	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	157/162 (97%)	150 (96%)	6 (4%)	1 (1%)	30	29
1	B	156/162 (96%)	154 (99%)	2 (1%)	0	100	100
1	C	157/162 (97%)	153 (98%)	4 (2%)	0	100	100
1	D	157/162 (97%)	154 (98%)	3 (2%)	0	100	100
All	All	627/648 (97%)	611 (97%)	15 (2%)	1 (0%)	52	59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	145	GLU

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	122/126 (97%)	113 (93%)	9 (7%)	17	17
1	B	122/126 (97%)	114 (93%)	8 (7%)	21	22
1	C	121/126 (96%)	112 (93%)	9 (7%)	17	17
1	D	121/126 (96%)	109 (90%)	12 (10%)	10	9
All	All	486/504 (96%)	448 (92%)	38 (8%)	16	15

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-12	HIS
1	A	-6	LEU
1	A	0	GLN
1	A	68	THR
1	A	89	GLU
1	A	105	LEU
1	A	116	LEU
1	A	139	GLU
1	A	144	ILE
1	B	-6	LEU
1	B	0	GLN
1	B	4	ILE
1	B	19	ILE
1	B	68	THR
1	B	89	GLU
1	B	105	LEU
1	B	116	LEU
1	C	-12	HIS
1	C	2	LYS
1	C	4	ILE
1	C	11	VAL
1	C	19	ILE
1	C	105	LEU
1	C	116	LEU
1	C	121	VAL
1	C	145	GLU
1	D	-10	HIS
1	D	-6	LEU

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Mol	Chain	Res	Type
1	D	0	GLN
1	D	4	ILE
1	D	19	ILE
1	D	59	VAL
1	D	89	GLU
1	D	105	LEU
1	D	116	LEU
1	D	128	GLN
1	D	142	THR
1	D	146	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	0	GLN
1	A	77	ASN
1	A	94	GLN
1	A	98	GLN
1	A	146	GLN
1	B	0	GLN
1	B	77	ASN
1	B	94	GLN
1	B	98	GLN
1	B	100	ASN
1	B	135	GLN
1	C	17	HIS
1	C	77	ASN
1	C	94	GLN
1	C	99	HIS
1	C	100	ASN
1	C	136	GLN
1	D	0	GLN
1	D	77	ASN
1	D	94	GLN
1	D	98	GLN
1	D	100	ASN
1	D	128	GLN
1	D	134	HIS

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 ⓘ

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.