Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 11:29 PM GMT

PDB ID	:	1XIQ
Title	:	Plasmodium falciparum Nucleoside diphosphate kinase B
Authors	:	Robien, M.A.; Bosch, J.; Hol, W.G.J.; Structural Genomics of Pathogenic Protozoa Consortium (SGPP)
Deposited on	:	2004-09-21
Resolution	:	3.05 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

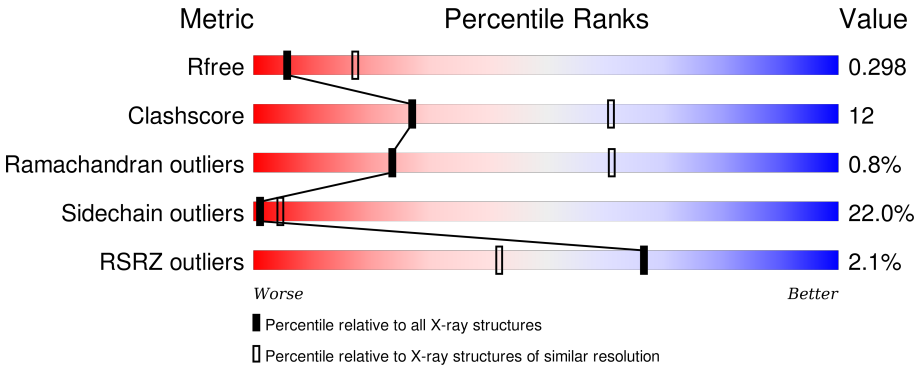
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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	1191 (3.10-3.02)
Clashscore	102246	1303 (3.10-3.02)
Ramachandran outliers	100387	1254 (3.10-3.02)
Sidechain outliers	100360	1254 (3.10-3.02)
RSRZ outliers	91569	1197 (3.10-3.02)

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 <=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	157	2% 60% 31% • 5%
1	B	157	3% 54% 34% 6% • 5%
1	C	157	3% 55% 33% 6% 5%
1	D	157	3% 57% 31% 6% • 5%
1	E	157	% 52% 32% 10% • 5%

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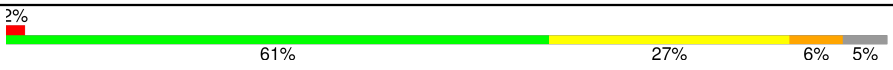
Mol	Chain	Length	Quality of chain
1	F	157	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7164 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 Nucleoside diphosphate kinase B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	0	0
			1194	769	202	215	8			
1	B	149	Total	C	N	O	S	0	0	0
			1194	769	202	215	8			
1	C	149	Total	C	N	O	S	0	0	0
			1194	769	202	215	8			
1	D	149	Total	C	N	O	S	0	0	0
			1194	769	202	215	8			
1	E	149	Total	C	N	O	S	0	0	0
			1194	769	202	215	8			
1	F	149	Total	C	N	O	S	0	0	0
			1194	769	202	215	8			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	EXPRESSION TAG	UNP Q8ID43
A	-6	ALA	-	EXPRESSION TAG	UNP Q8ID43
A	-5	HIS	-	EXPRESSION TAG	UNP Q8ID43
A	-4	HIS	-	EXPRESSION TAG	UNP Q8ID43
A	-3	HIS	-	EXPRESSION TAG	UNP Q8ID43
A	-2	HIS	-	EXPRESSION TAG	UNP Q8ID43
A	-1	HIS	-	EXPRESSION TAG	UNP Q8ID43
A	0	HIS	-	EXPRESSION TAG	UNP Q8ID43
B	-7	MET	-	EXPRESSION TAG	UNP Q8ID43
B	-6	ALA	-	EXPRESSION TAG	UNP Q8ID43
B	-5	HIS	-	EXPRESSION TAG	UNP Q8ID43
B	-4	HIS	-	EXPRESSION TAG	UNP Q8ID43
B	-3	HIS	-	EXPRESSION TAG	UNP Q8ID43
B	-2	HIS	-	EXPRESSION TAG	UNP Q8ID43
B	-1	HIS	-	EXPRESSION TAG	UNP Q8ID43
B	0	HIS	-	EXPRESSION TAG	UNP Q8ID43
C	-7	MET	-	EXPRESSION TAG	UNP Q8ID43

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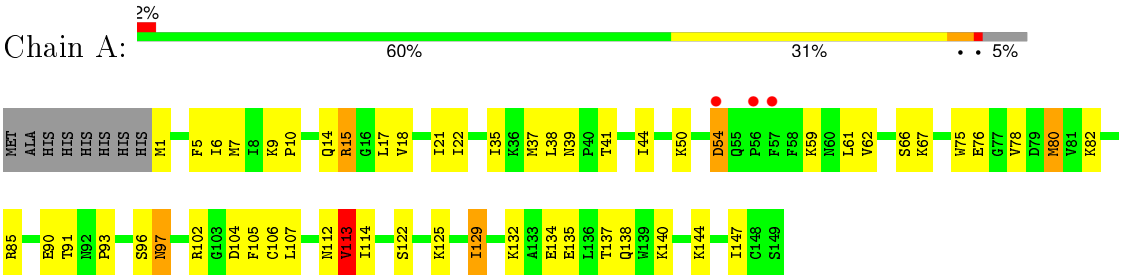
Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	ALA	-	EXPRESSION TAG	UNP Q8ID43
C	-5	HIS	-	EXPRESSION TAG	UNP Q8ID43
C	-4	HIS	-	EXPRESSION TAG	UNP Q8ID43
C	-3	HIS	-	EXPRESSION TAG	UNP Q8ID43
C	-2	HIS	-	EXPRESSION TAG	UNP Q8ID43
C	-1	HIS	-	EXPRESSION TAG	UNP Q8ID43
C	0	HIS	-	EXPRESSION TAG	UNP Q8ID43
D	-7	MET	-	EXPRESSION TAG	UNP Q8ID43
D	-6	ALA	-	EXPRESSION TAG	UNP Q8ID43
D	-5	HIS	-	EXPRESSION TAG	UNP Q8ID43
D	-4	HIS	-	EXPRESSION TAG	UNP Q8ID43
D	-3	HIS	-	EXPRESSION TAG	UNP Q8ID43
D	-2	HIS	-	EXPRESSION TAG	UNP Q8ID43
D	-1	HIS	-	EXPRESSION TAG	UNP Q8ID43
D	0	HIS	-	EXPRESSION TAG	UNP Q8ID43
E	-7	MET	-	EXPRESSION TAG	UNP Q8ID43
E	-6	ALA	-	EXPRESSION TAG	UNP Q8ID43
E	-5	HIS	-	EXPRESSION TAG	UNP Q8ID43
E	-4	HIS	-	EXPRESSION TAG	UNP Q8ID43
E	-3	HIS	-	EXPRESSION TAG	UNP Q8ID43
E	-2	HIS	-	EXPRESSION TAG	UNP Q8ID43
E	-1	HIS	-	EXPRESSION TAG	UNP Q8ID43
E	0	HIS	-	EXPRESSION TAG	UNP Q8ID43
F	-7	MET	-	EXPRESSION TAG	UNP Q8ID43
F	-6	ALA	-	EXPRESSION TAG	UNP Q8ID43
F	-5	HIS	-	EXPRESSION TAG	UNP Q8ID43
F	-4	HIS	-	EXPRESSION TAG	UNP Q8ID43
F	-3	HIS	-	EXPRESSION TAG	UNP Q8ID43
F	-2	HIS	-	EXPRESSION TAG	UNP Q8ID43
F	-1	HIS	-	EXPRESSION TAG	UNP Q8ID43
F	0	HIS	-	EXPRESSION TAG	UNP Q8ID43

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.

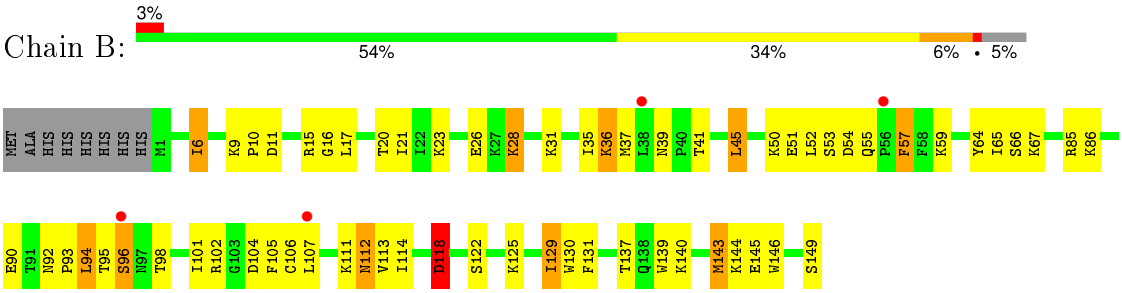
- Molecule 1: Nucleoside diphosphate kinase B

Chain A:



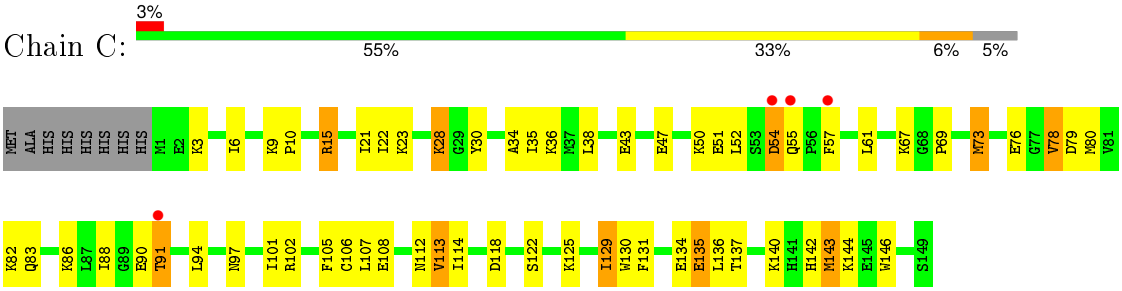
- Molecule 1: Nucleoside diphosphate kinase B

Chain B:



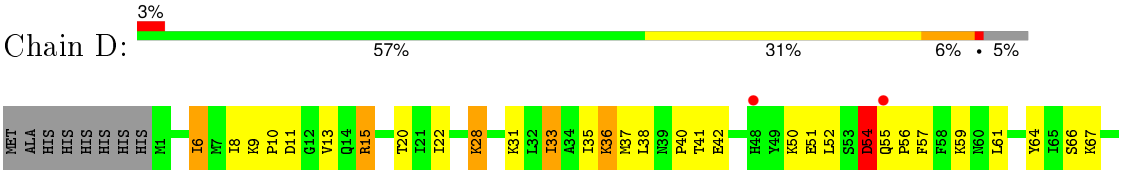
- Molecule 1: Nucleoside diphosphate kinase B

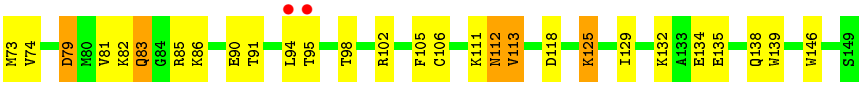
Chain C:



- Molecule 1: Nucleoside diphosphate kinase B

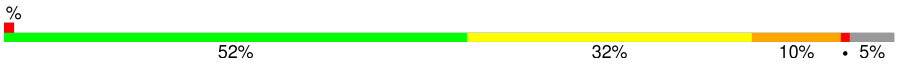
Chain D:

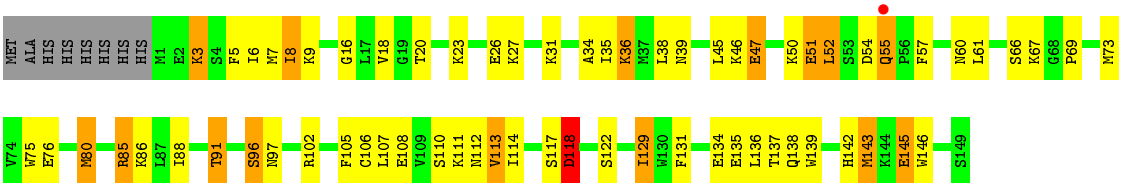




• Molecule 1: Nucleoside diphosphate kinase B

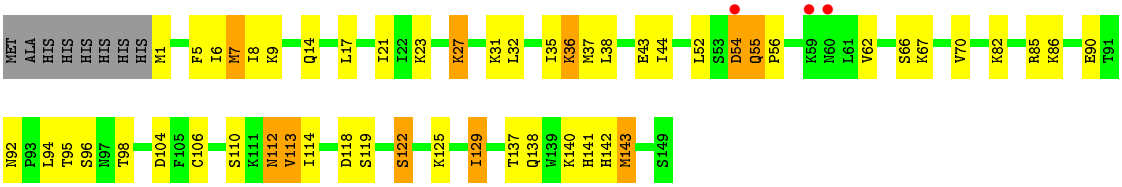
Chain E:





• Molecule 1: Nucleoside diphosphate kinase B

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	68.59Å 112.42Å 68.59Å 90.00° 119.05° 90.00°	Depositor
Resolution (Å)	52.70 – 3.05 41.01 – 3.05	Depositor EDS
% Data completeness (in resolution range)	99.5 (52.70-3.05) 99.4 (41.01-3.05)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$< I/\sigma(I) >$ [1]	1.96 (at 3.06Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.229 , 0.286 0.239 , 0.298	Depositor DCC
R_{free} test set	876 reflections (5.33%)	DCC
Wilson B-factor (Å^2)	69.2	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 8.6	EDS
Estimated twinning fraction	0.011 for -h-l,k,h 0.011 for l,k,-h-l 0.026 for h,-k,-h-l 0.023 for -h-l,-k,l 0.448 for l,-k,h	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 17324 reflections	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	7164	wwPDB-VP
Average B, all atoms (Å^2)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.46% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.47	0/1219	0.71	2/1638 (0.1%)
1	B	0.47	0/1219	0.67	2/1638 (0.1%)
1	C	0.48	0/1219	0.66	2/1638 (0.1%)
1	D	0.47	0/1219	0.70	3/1638 (0.2%)
1	E	0.45	0/1219	0.66	2/1638 (0.1%)
1	F	0.48	0/1219	0.70	2/1638 (0.1%)
All	All	0.47	0/7314	0.68	13/9828 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	104	ASP	CB-CG-OD2	6.67	124.30	118.30
1	C	118	ASP	CB-CG-OD2	6.10	123.79	118.30
1	D	79	ASP	CB-CG-OD2	5.82	123.54	118.30
1	F	104	ASP	CB-CG-OD2	5.52	123.27	118.30
1	B	118	ASP	CB-CG-OD2	5.49	123.24	118.30
1	D	118	ASP	CB-CG-OD2	5.48	123.23	118.30
1	C	54	ASP	CB-CG-OD2	5.42	123.18	118.30
1	F	54	ASP	CB-CG-OD2	5.39	123.15	118.30
1	E	54	ASP	CB-CG-OD2	5.38	123.14	118.30
1	E	118	ASP	CB-CG-OD2	5.30	123.07	118.30
1	A	54	ASP	CB-CG-OD2	5.30	123.07	118.30
1	B	54	ASP	CB-CG-OD2	5.25	123.02	118.30
1	D	54	ASP	CB-CG-OD2	5.03	122.82	118.30

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	1194	0	1228	34	0
1	B	1194	0	1228	46	0
1	C	1194	0	1228	37	0
1	D	1194	0	1228	30	0
1	E	1194	0	1228	41	0
1	F	1194	0	1228	28	0
All	All	7164	0	7368	174	0

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

All (174) 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:A:15:ARG:HH11	1:A:15:ARG:HG2	1.17	1.04
1:E:75:TRP:CB	1:E:80:MET:HE3	1.88	1.04
1:A:75:TRP:HB3	1:A:80:MET:HE3	1.38	1.02
1:D:15:ARG:HG2	1:D:15:ARG:HH11	1.26	0.99
1:E:75:TRP:HB3	1:E:80:MET:HE3	1.42	0.99
1:A:75:TRP:CB	1:A:80:MET:HE3	1.95	0.95
1:B:102:ARG:HD3	1:B:112:ASN:HB2	1.47	0.93
1:C:15:ARG:HH11	1:C:15:ARG:HG2	1.33	0.93
1:A:75:TRP:HB3	1:A:80:MET:CE	1.98	0.93
1:E:75:TRP:HB3	1:E:80:MET:CE	2.00	0.90
1:D:102:ARG:HD3	1:D:112:ASN:HB2	1.57	0.85
1:E:75:TRP:HB2	1:E:80:MET:HE3	1.64	0.77
1:E:3:LYS:HD2	1:E:76:GLU:HG3	1.65	0.76
1:E:102:ARG:HD3	1:E:112:ASN:HB2	1.67	0.76
1:C:102:ARG:HD3	1:C:112:ASN:HB2	1.69	0.74
1:A:106:CYS:SG	1:A:113:VAL:HG13	2.28	0.74
1:A:15:ARG:HG2	1:A:15:ARG:NH1	1.95	0.73
1:B:125:LYS:O	1:B:129:ILE:HG23	1.89	0.73
1:C:22:ILE:HG12	1:C:73:MET:CE	2.21	0.70
1:C:15:ARG:NH1	1:C:15:ARG:HG2	2.04	0.69
1:F:44:ILE:HA	1:F:129:ILE:HD11	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:ASP:HB2	1:C:146:TRP:O	1.95	0.67
1:C:3:LYS:HG3	1:C:76:GLU:HG3	1.76	0.66
1:F:125:LYS:O	1:F:129:ILE:HG22	1.95	0.66
1:A:44:ILE:HA	1:A:129:ILE:HD11	1.79	0.65
1:B:31:LYS:HG3	1:B:139:TRP:CZ2	2.32	0.65
1:B:6:ILE:HG23	1:B:114:ILE:HD11	1.78	0.64
1:D:15:ARG:NH2	1:D:105:PHE:O	2.30	0.64
1:A:125:LYS:O	1:A:129:ILE:HG22	1.97	0.63
1:E:88:ILE:HG12	1:E:114:ILE:HG21	1.81	0.63
1:A:15:ARG:NH2	1:A:105:PHE:O	2.31	0.63
1:D:9:LYS:HB3	1:D:10:PRO:HD2	1.80	0.62
1:C:15:ARG:NH2	1:C:105:PHE:O	2.32	0.62
1:B:21:ILE:HG12	1:B:101:ILE:HD13	1.82	0.61
1:D:15:ARG:NH1	1:D:15:ARG:HG2	2.01	0.61
1:B:94:LEU:HD13	1:C:79:ASP:OD2	2.00	0.60
1:B:145:GLU:OE2	1:F:142:HIS:CE1	2.54	0.60
1:B:15:ARG:NH2	1:B:105:PHE:O	2.30	0.59
1:A:38:LEU:HA	1:F:137:THR:HG21	1.85	0.59
1:C:22:ILE:HG12	1:C:73:MET:HE1	1.84	0.58
1:D:11:ASP:HB2	1:E:146:TRP:O	2.03	0.58
1:B:9:LYS:HB3	1:B:10:PRO:HD2	1.85	0.57
1:C:35:ILE:HD12	1:C:73:MET:HG2	1.84	0.57
1:D:40:PRO:HG2	1:D:66:SER:HA	1.85	0.57
1:A:93:PRO:HA	1:A:96:SER:OG	2.05	0.57
1:A:137:THR:HG21	1:F:38:LEU:HA	1.86	0.56
1:C:36:LYS:HE3	1:C:38:LEU:HD13	1.86	0.56
1:E:31:LYS:HG3	1:E:139:TRP:CZ2	2.40	0.56
1:F:55:GLN:HG2	1:F:56:PRO:HD2	1.86	0.56
1:C:30:TYR:CD1	1:C:80:MET:SD	2.98	0.56
1:C:6:ILE:HG23	1:C:114:ILE:HD11	1.87	0.56
1:E:6:ILE:HG23	1:E:114:ILE:HD11	1.88	0.56
1:D:31:LYS:HG3	1:D:139:TRP:CZ2	2.41	0.56
1:F:62:VAL:O	1:F:66:SER:HB2	2.07	0.55
1:A:75:TRP:HB2	1:A:80:MET:HE3	1.84	0.55
1:E:105:PHE:HE1	1:F:27:LYS:HD3	1.71	0.55
1:F:17:LEU:O	1:F:21:ILE:HD12	2.07	0.55
1:B:145:GLU:OE2	1:F:142:HIS:HE1	1.89	0.54
1:D:102:ARG:CD	1:D:112:ASN:HB2	2.32	0.54
1:B:111:LYS:HD2	1:C:146:TRP:HA	1.89	0.54
1:A:132:LYS:HB2	1:A:135:GLU:HG3	1.89	0.53
1:C:91:THR:HA	1:C:102:ARG:NH1	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:ILE:CG2	1:C:107:LEU:HD12	2.39	0.53
1:A:14:GLN:HG3	1:B:146:TRP:CE2	2.43	0.53
1:D:55:GLN:HG3	1:D:56:PRO:HD2	1.91	0.53
1:F:6:ILE:HG23	1:F:114:ILE:HD11	1.91	0.53
1:B:17:LEU:O	1:B:21:ILE:HD12	2.08	0.53
1:C:34:ALA:HB1	1:C:131:PHE:CE2	2.44	0.53
1:E:91:THR:HA	1:E:102:ARG:NH1	2.24	0.52
1:B:45:LEU:HD11	1:B:65:ILE:HG22	1.92	0.52
1:B:102:ARG:CD	1:B:112:ASN:HB2	2.30	0.52
1:B:36:LYS:HA	1:E:35:ILE:O	2.11	0.51
1:C:106:CYS:SG	1:C:113:VAL:HG13	2.51	0.51
1:E:34:ALA:HB1	1:E:131:PHE:CE1	2.46	0.51
1:B:37:MET:HB2	1:E:35:ILE:HG22	1.94	0.50
1:D:146:TRP:CE2	1:F:14:GLN:HG3	2.47	0.50
1:B:52:LEU:HD23	1:B:55:GLN:HG3	1.94	0.50
1:C:35:ILE:O	1:D:36:LYS:HA	2.10	0.50
1:A:102:ARG:HD3	1:A:112:ASN:HB2	1.94	0.49
1:C:143:MET:SD	1:E:143:MET:CE	3.00	0.49
1:A:18:VAL:O	1:A:22:ILE:HG13	2.12	0.49
1:E:85:ARG:HD2	1:E:118:ASP:HA	1.92	0.49
1:B:16:GLY:HA2	1:E:26:GLU:OE2	2.13	0.49
1:C:35:ILE:HG22	1:D:37:MET:HB2	1.95	0.49
1:B:36:LYS:HD2	1:B:131:PHE:CE1	2.48	0.48
1:A:9:LYS:HB3	1:A:10:PRO:CD	2.43	0.48
1:B:35:ILE:O	1:E:36:LYS:HA	2.14	0.48
1:A:62:VAL:O	1:A:66:SER:HB2	2.12	0.48
1:D:102:ARG:HD3	1:D:112:ASN:CB	2.38	0.48
1:B:143:MET:HE1	1:F:143:MET:HG2	1.96	0.48
1:F:32:LEU:HD11	1:F:35:ILE:HB	1.95	0.48
1:A:17:LEU:O	1:A:21:ILE:HD12	2.14	0.47
1:D:15:ARG:HA	1:E:143:MET:HE3	1.94	0.47
1:C:88:ILE:HG12	1:C:114:ILE:HG21	1.96	0.47
1:B:137:THR:HG21	1:E:38:LEU:HA	1.95	0.47
1:F:9:LYS:NZ	1:F:112:ASN:HD21	2.13	0.47
1:D:6:ILE:O	1:D:73:MET:HB2	2.15	0.47
1:D:8:ILE:N	1:D:8:ILE:HD12	2.30	0.47
1:B:28:LYS:O	1:B:28:LYS:HG3	2.14	0.47
1:B:26:GLU:OE1	1:E:16:GLY:HA2	2.14	0.47
1:D:9:LYS:HB3	1:D:10:PRO:CD	2.45	0.47
1:A:37:MET:HB2	1:F:35:ILE:HG22	1.97	0.47
1:D:33:ILE:HD11	1:D:74:VAL:CG1	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ILE:HG23	1:A:129:ILE:HD11	1.97	0.46
1:C:30:TYR:CE1	1:C:80:MET:SD	3.09	0.46
1:B:137:THR:HB	1:E:39:ASN:ND2	2.32	0.45
1:D:106:CYS:SG	1:D:113:VAL:HG13	2.56	0.45
1:A:35:ILE:HG22	1:F:37:MET:HB2	1.98	0.45
1:F:106:CYS:SG	1:F:113:VAL:HG13	2.56	0.45
1:E:51:GLU:HG2	1:E:52:LEU:N	2.31	0.45
1:F:8:ILE:HG12	1:F:21:ILE:HD13	1.99	0.45
1:D:111:LYS:HD2	1:E:146:TRP:HA	1.98	0.45
1:B:107:LEU:HB3	1:C:78:VAL:HG13	1.99	0.45
1:B:92:ASN:HA	1:B:93:PRO:HD3	1.86	0.45
1:E:8:ILE:HD12	1:E:18:VAL:HA	1.98	0.45
1:B:23:LYS:HD2	1:E:20:THR:OG1	2.16	0.44
1:F:31:LYS:HD2	1:F:141:HIS:HB2	1.98	0.44
1:B:143:MET:CE	1:F:143:MET:HG2	2.47	0.44
1:E:106:CYS:SG	1:E:113:VAL:HG13	2.57	0.44
1:B:11:ASP:HB3	1:B:111:LYS:HD3	1.99	0.44
1:E:88:ILE:HG12	1:E:114:ILE:CG2	2.47	0.44
1:A:1:MET:HB3	1:A:76:GLU:HG2	2.00	0.44
1:D:28:LYS:NZ	1:F:106:CYS:O	2.49	0.44
1:B:118:ASP:OD1	1:B:118:ASP:N	2.49	0.43
1:A:147:ILE:HG22	1:C:107:LEU:HD12	1.99	0.43
1:A:6:ILE:HG23	1:A:114:ILE:HD11	2.01	0.43
1:A:35:ILE:O	1:F:36:LYS:HA	2.18	0.43
1:F:92:ASN:ND2	1:F:95:THR:OG1	2.50	0.43
1:A:41:THR:HG22	1:A:44:ILE:HD12	2.01	0.43
1:B:137:THR:HB	1:E:39:ASN:HD21	1.84	0.43
1:F:5:PHE:CZ	1:F:7:MET:HB2	2.53	0.43
1:C:21:ILE:HG12	1:C:101:ILE:HD13	2.00	0.43
1:F:119:SER:OG	1:F:122:SER:HB2	2.19	0.43
1:A:5:PHE:CZ	1:A:7:MET:HB2	2.53	0.42
1:A:97:ASN:ND2	1:A:97:ASN:H	2.16	0.42
1:E:102:ARG:CD	1:E:112:ASN:HB2	2.43	0.42
1:B:106:CYS:O	1:C:28:LYS:HE2	2.20	0.42
1:B:41:THR:O	1:B:45:LEU:HB2	2.20	0.42
1:B:93:PRO:HA	1:B:96:SER:HB3	2.01	0.42
1:D:91:THR:HA	1:D:102:ARG:NH1	2.35	0.42
1:E:5:PHE:CZ	1:E:7:MET:HB2	2.55	0.42
1:D:8:ILE:N	1:D:8:ILE:CD1	2.82	0.42
1:A:107:LEU:HD23	1:B:28:LYS:HE2	2.01	0.42
1:E:108:GLU:OE2	1:E:110:SER:OG	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:69:PRO:HG3	1:D:139:TRP:CE3	2.54	0.42
1:B:139:TRP:CE3	1:E:69:PRO:HG3	2.55	0.42
1:D:15:ARG:CG	1:D:15:ARG:NH1	2.75	0.41
1:A:39:ASN:ND2	1:A:67:LYS:O	2.53	0.41
1:C:142:HIS:CG	1:E:146:TRP:HE1	2.38	0.41
1:C:36:LYS:HA	1:D:35:ILE:O	2.21	0.41
1:E:6:ILE:HG22	1:E:73:MET:HE3	2.02	0.41
1:D:22:ILE:HD11	1:D:73:MET:HE2	2.02	0.41
1:C:129:ILE:HG22	1:C:130:TRP:CD1	2.55	0.41
1:E:47:GLU:OE2	1:E:129:ILE:HD11	2.21	0.41
1:B:39:ASN:HD22	1:B:39:ASN:N	2.19	0.41
1:B:129:ILE:CG1	1:B:130:TRP:CD1	3.04	0.41
1:C:10:PRO:HA	1:C:69:PRO:O	2.21	0.41
1:A:112:ASN:O	1:A:114:ILE:N	2.49	0.41
1:E:145:GLU:HG2	1:E:145:GLU:H	1.68	0.41
1:F:90:GLU:HB2	1:F:96:SER:HB3	2.03	0.41
1:B:57:PHE:HA	1:B:57:PHE:HD1	1.76	0.41
1:D:125:LYS:O	1:D:129:ILE:HG23	2.20	0.41
1:B:11:ASP:HB3	1:B:111:LYS:CD	2.51	0.41
1:E:55:GLN:HE21	1:E:55:GLN:HB3	1.62	0.41
1:E:96:SER:O	1:F:86:LYS:HE2	2.21	0.40
1:B:9:LYS:HB3	1:B:10:PRO:CD	2.51	0.40
1:C:88:ILE:HG12	1:C:114:ILE:CG2	2.50	0.40
1:E:85:ARG:HD2	1:E:117:SER:O	2.21	0.40
1:D:79:ASP:O	1:D:83:GLN:HB2	2.22	0.40
1:C:3:LYS:CG	1:C:76:GLU:HG3	2.50	0.40
1:B:6:ILE:CG2	1:B:114:ILE:HD11	2.50	0.40
1:C:79:ASP:O	1:C:83:GLN:HB2	2.22	0.40
1:C:131:PHE:HD2	1:C:135:GLU:HB3	1.87	0.40
1:B:104:ASP:O	1:C:28:LYS:HD2	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	147/157 (94%)	141 (96%)	5 (3%)	1 (1%)	26	64
1	B	147/157 (94%)	143 (97%)	3 (2%)	1 (1%)	26	64
1	C	147/157 (94%)	141 (96%)	5 (3%)	1 (1%)	26	64
1	D	147/157 (94%)	140 (95%)	5 (3%)	2 (1%)	14	46
1	E	147/157 (94%)	141 (96%)	5 (3%)	1 (1%)	26	64
1	F	147/157 (94%)	141 (96%)	5 (3%)	1 (1%)	26	64
All	All	882/942 (94%)	847 (96%)	28 (3%)	7 (1%)	24	61

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	113	VAL
1	B	113	VAL
1	C	113	VAL
1	D	113	VAL
1	E	113	VAL
1	F	113	VAL
1	D	54	ASP

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	132/139 (95%)	113 (86%)	19 (14%)	4	16
1	B	132/139 (95%)	104 (79%)	28 (21%)	1	5
1	C	132/139 (95%)	99 (75%)	33 (25%)	1	2
1	D	132/139 (95%)	98 (74%)	34 (26%)	0	2
1	E	132/139 (95%)	95 (72%)	37 (28%)	0	1
1	F	132/139 (95%)	109 (83%)	23 (17%)	2	10
All	All	792/834 (95%)	618 (78%)	174 (22%)	1	4

All (174) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ARG
1	A	50	LYS
1	A	54	ASP
1	A	59	LYS
1	A	61	LEU
1	A	78	VAL
1	A	80	MET
1	A	82	LYS
1	A	85	ARG
1	A	90	GLU
1	A	91	THR
1	A	97	ASN
1	A	113	VAL
1	A	122	SER
1	A	129	ILE
1	A	134	GLU
1	A	138	GLN
1	A	140	LYS
1	A	144	LYS
1	B	6	ILE
1	B	20	THR
1	B	28	LYS
1	B	36	LYS
1	B	45	LEU
1	B	50	LYS
1	B	51	GLU
1	B	53	SER
1	B	57	PHE
1	B	59	LYS
1	B	64	TYR
1	B	66	SER
1	B	67	LYS
1	B	85	ARG
1	B	86	LYS
1	B	90	GLU
1	B	94	LEU
1	B	95	THR
1	B	96	SER
1	B	98	THR
1	B	112	ASN
1	B	118	ASP
1	B	122	SER

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Mol	Chain	Res	Type
1	B	129	ILE
1	B	140	LYS
1	B	143	MET
1	B	144	LYS
1	B	149	SER
1	C	9	LYS
1	C	15	ARG
1	C	23	LYS
1	C	28	LYS
1	C	43	GLU
1	C	47	GLU
1	C	50	LYS
1	C	51	GLU
1	C	52	LEU
1	C	54	ASP
1	C	55	GLN
1	C	57	PHE
1	C	61	LEU
1	C	67	LYS
1	C	73	MET
1	C	78	VAL
1	C	82	LYS
1	C	86	LYS
1	C	90	GLU
1	C	91	THR
1	C	94	LEU
1	C	97	ASN
1	C	108	GLU
1	C	122	SER
1	C	125	LYS
1	C	129	ILE
1	C	134	GLU
1	C	135	GLU
1	C	136	LEU
1	C	137	THR
1	C	140	LYS
1	C	143	MET
1	C	144	LYS
1	D	6	ILE
1	D	13	VAL
1	D	15	ARG
1	D	20	THR

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Mol	Chain	Res	Type
1	D	28	LYS
1	D	33	ILE
1	D	36	LYS
1	D	38	LEU
1	D	41	THR
1	D	42	GLU
1	D	50	LYS
1	D	51	GLU
1	D	52	LEU
1	D	54	ASP
1	D	57	PHE
1	D	59	LYS
1	D	61	LEU
1	D	64	TYR
1	D	67	LYS
1	D	81	VAL
1	D	82	LYS
1	D	83	GLN
1	D	85	ARG
1	D	86	LYS
1	D	90	GLU
1	D	94	LEU
1	D	95	THR
1	D	98	THR
1	D	112	ASN
1	D	125	LYS
1	D	132	LYS
1	D	134	GLU
1	D	135	GLU
1	D	138	GLN
1	E	3	LYS
1	E	8	ILE
1	E	9	LYS
1	E	23	LYS
1	E	27	LYS
1	E	36	LYS
1	E	45	LEU
1	E	46	LYS
1	E	47	GLU
1	E	50	LYS
1	E	51	GLU
1	E	52	LEU

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Mol	Chain	Res	Type
1	E	55	GLN
1	E	57	PHE
1	E	60	ASN
1	E	61	LEU
1	E	66	SER
1	E	67	LYS
1	E	80	MET
1	E	85	ARG
1	E	86	LYS
1	E	91	THR
1	E	96	SER
1	E	97	ASN
1	E	107	LEU
1	E	111	LYS
1	E	118	ASP
1	E	122	SER
1	E	129	ILE
1	E	134	GLU
1	E	135	GLU
1	E	136	LEU
1	E	137	THR
1	E	138	GLN
1	E	142	HIS
1	E	143	MET
1	E	145	GLU
1	F	1	MET
1	F	7	MET
1	F	23	LYS
1	F	27	LYS
1	F	36	LYS
1	F	43	GLU
1	F	52	LEU
1	F	54	ASP
1	F	55	GLN
1	F	67	LYS
1	F	70	VAL
1	F	82	LYS
1	F	85	ARG
1	F	94	LEU
1	F	98	THR
1	F	110	SER
1	F	112	ASN

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Mol	Chain	Res	Type
1	F	118	ASP
1	F	122	SER
1	F	129	ILE
1	F	138	GLN
1	F	140	LYS
1	F	143	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	97	ASN
1	B	39	ASN
1	B	112	ASN
1	C	128	ASN
1	C	141	HIS
1	D	83	GLN
1	D	92	ASN
1	D	112	ASN
1	E	39	ASN
1	E	55	GLN
1	E	83	GLN
1	E	128	ASN
1	E	138	GLN
1	F	39	ASN
1	F	112	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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^{th} 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^2)	Q<0.9
1	A	149/157 (94%)	-0.13	3 (2%)	68	44	17, 39, 120, 151	0
1	B	149/157 (94%)	-0.02	4 (2%)	58	32	16, 40, 139, 151	0
1	C	149/157 (94%)	-0.12	4 (2%)	58	32	16, 39, 139, 151	0
1	D	149/157 (94%)	-0.02	4 (2%)	58	32	17, 40, 142, 151	0
1	E	149/157 (94%)	-0.16	1 (0%)	89	75	16, 39, 140, 151	0
1	F	149/157 (94%)	-0.11	3 (2%)	68	44	16, 39, 119, 151	0
All	All	894/942 (94%)	-0.10	19 (2%)	67	42	16, 40, 139, 151	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	54	ASP	4.0
1	B	56	PRO	3.5
1	A	54	ASP	3.3
1	C	54	ASP	3.3
1	E	55	GLN	3.3
1	A	56	PRO	3.2
1	D	48	HIS	2.8
1	C	55	GLN	2.7
1	C	91	THR	2.6
1	A	57	PHE	2.5
1	D	55	GLN	2.5
1	B	96	SER	2.4
1	C	57	PHE	2.3
1	F	59	LYS	2.3
1	F	60	ASN	2.2
1	B	107	LEU	2.1
1	B	38	LEU	2.1
1	D	95	THR	2.1
1	D	94	LEU	2.0

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

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.