Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:49 AM GMT

PDB ID	:	2YHF
Title	:	1.9 ANGSTROM CRYSTAL STRUCTURE OF CLEC5A
Authors	:	Watson, A.A.; Lebedev, A.A.; Murshudov, G.M.; Vagin, A.A.; Hall, B.A.; O'Callaghan, C.A.
Deposited on	:	2011-04-30
Resolution	:	1.90 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 ①) were used in the production of this report:

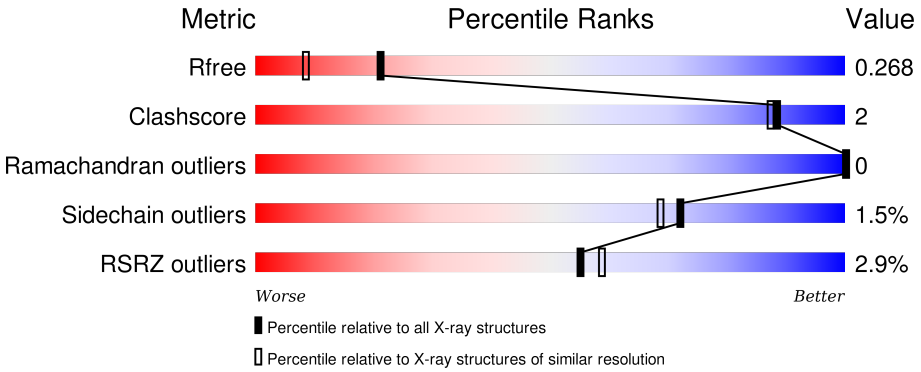
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	118	2% 97%
1	B	118	% 91% 9%
1	C	118	4% 89% 11%
1	D	118	2% 92% 7%
1	E	118	3% 95% 5%

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Mol	Chain	Length	Quality of chain
1	F	118	2% 92% 8%
1	G	118	2% 95% 5%
1	H	118	% 92% 8%
1	I	118	11% 94% 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9688 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 C-TYPE LECTIN DOMAIN FAMILY 5 MEMBER A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	118	Total	C	N	O	S	0	0	0
			972	613	169	183	7			
1	B	118	Total	C	N	O	S	0	0	0
			972	613	169	183	7			
1	C	118	Total	C	N	O	S	0	0	0
			972	613	169	183	7			
1	D	118	Total	C	N	O	S	0	0	0
			972	613	169	183	7			
1	E	118	Total	C	N	O	S	0	0	0
			972	613	169	183	7			
1	F	118	Total	C	N	O	S	0	0	0
			972	613	169	183	7			
1	G	118	Total	C	N	O	S	0	0	0
			972	613	169	183	7			
1	H	118	Total	C	N	O	S	0	0	0
			972	613	169	183	7			
1	I	118	Total	C	N	O	S	0	0	0
			972	613	169	183	7			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	70	MET	VAL	ENGINEERED MUTATION	UNP Q9NY25
B	70	MET	VAL	ENGINEERED MUTATION	UNP Q9NY25
C	70	MET	VAL	ENGINEERED MUTATION	UNP Q9NY25
D	70	MET	VAL	ENGINEERED MUTATION	UNP Q9NY25
E	70	MET	VAL	ENGINEERED MUTATION	UNP Q9NY25
F	70	MET	VAL	ENGINEERED MUTATION	UNP Q9NY25
G	70	MET	VAL	ENGINEERED MUTATION	UNP Q9NY25
H	70	MET	VAL	ENGINEERED MUTATION	UNP Q9NY25
I	70	MET	VAL	ENGINEERED MUTATION	UNP Q9NY25

- Molecule 2 is water.

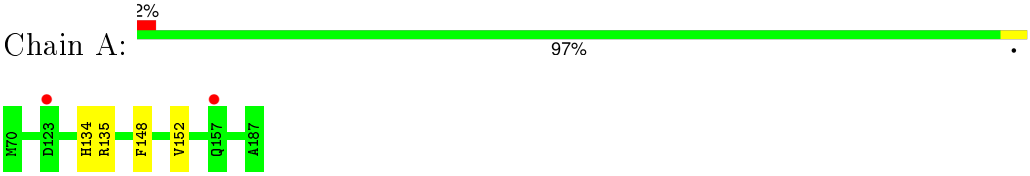
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	144	Total O 144 144	0	0
2	B	128	Total O 128 128	0	0
2	C	97	Total O 97 97	0	0
2	D	98	Total O 98 98	0	0
2	E	113	Total O 113 113	0	0
2	F	98	Total O 98 98	0	0
2	G	98	Total O 98 98	0	0
2	H	108	Total O 108 108	0	0
2	I	56	Total O 56 56	0	0

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.

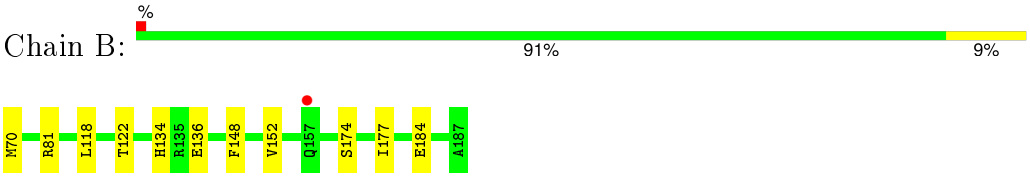
- Molecule 1: C-TYPE LECTIN DOMAIN FAMILY 5 MEMBER A

Chain A:



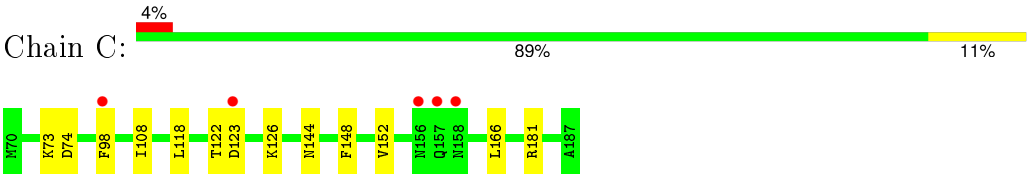
- Molecule 1: C-TYPE LECTIN DOMAIN FAMILY 5 MEMBER A

Chain B:



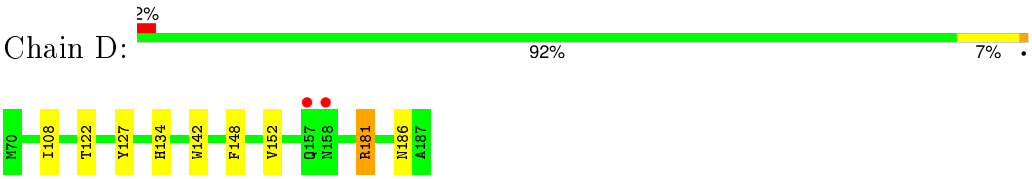
- Molecule 1: C-TYPE LECTIN DOMAIN FAMILY 5 MEMBER A

Chain C:



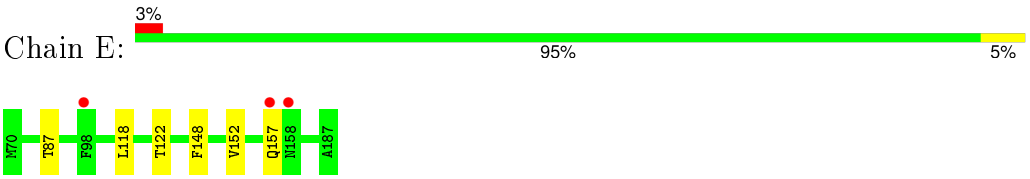
- Molecule 1: C-TYPE LECTIN DOMAIN FAMILY 5 MEMBER A

Chain D:



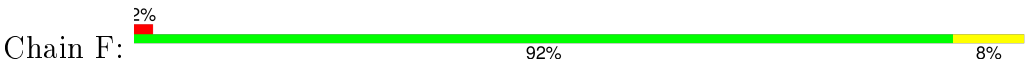
- Molecule 1: C-TYPE LECTIN DOMAIN FAMILY 5 MEMBER A

Chain E:



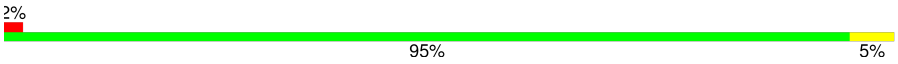
- Molecule 1: C-TYPE LECTIN DOMAIN FAMILY 5 MEMBER A

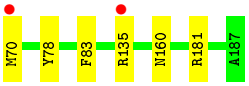
Chain F:



• Molecule 1: C-TYPE LECTIN DOMAIN FAMILY 5 MEMBER A

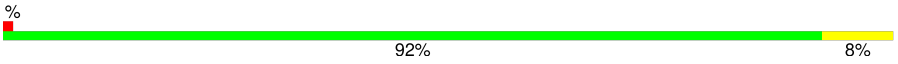
Chain G:





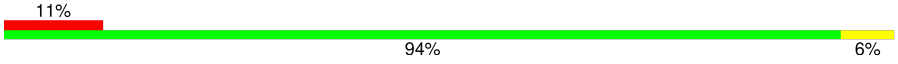
• Molecule 1: C-TYPE LECTIN DOMAIN FAMILY 5 MEMBER A

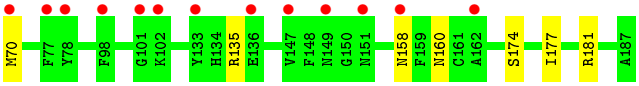
Chain H:



• Molecule 1: C-TYPE LECTIN DOMAIN FAMILY 5 MEMBER A

Chain I:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α, β, γ	109.11Å 109.11Å 84.88Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	94.49 – 1.90 32.92 – 1.90	Depositor EDS
% Data completeness (in resolution range)	87.2 (94.49-1.90) 83.5 (32.92-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.82 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.3.0002	Depositor
R, R_{free}	0.216 , 0.267 0.216 , 0.268	Depositor DCC
R_{free} test set	3728 reflections (5.31%)	DCC
Wilson B-factor (Å^2)	20.6	Xtriage
Anisotropy	0.372	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 24.4	EDS
Estimated twinning fraction	0.043 for -h,-k,l 0.048 for h,-h-k,-l 0.177 for -k,-h,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.27$	Xtriage
Outliers	0 of 73967 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	9688	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 48.70 % of the origin peak, indicating pseudo translational symmetry.The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 8.2880e-05.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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	0.45	0/995	0.54	0/1340
1	B	0.42	0/995	0.53	0/1340
1	C	0.44	0/995	0.55	0/1340
1	D	0.40	0/995	0.53	0/1340
1	E	0.41	0/995	0.52	0/1340
1	F	0.40	0/995	0.52	0/1340
1	G	0.40	0/995	0.50	0/1340
1	H	0.40	0/995	0.51	0/1340
1	I	0.38	0/995	0.47	0/1340
All	All	0.41	0/8955	0.52	0/12060

There are no bond length outliers.

There are no bond angle outliers.

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	972	0	920	5	0
1	B	972	0	920	5	0
1	C	972	0	920	6	0
1	D	972	0	920	8	0
1	E	972	0	920	2	0
1	F	972	0	920	6	0
1	G	972	0	920	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	972	0	920	6	0
1	I	972	0	920	4	0
2	A	144	0	0	2	0
2	B	128	0	0	0	0
2	C	97	0	0	0	0
2	D	98	0	0	2	0
2	E	113	0	0	0	0
2	F	98	0	0	0	0
2	G	98	0	0	0	0
2	H	108	0	0	0	0
2	I	56	0	0	0	0
All	All	9688	0	8280	42	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 2.

All (42) 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:H:81:ARG:HH12	1:H:186:ASN:HD22	1.11	0.97
1:I:135:ARG:HD3	1:I:160:ASN:HD21	1.34	0.89
1:H:81:ARG:NH1	1:H:186:ASN:HD22	1.86	0.71
1:I:135:ARG:HD3	1:I:160:ASN:ND2	2.05	0.71
1:H:81:ARG:HH12	1:H:186:ASN:ND2	1.88	0.71
1:A:148:PHE:CE2	1:A:152:VAL:HG23	2.36	0.60
1:C:98:PHE:HB3	1:C:181:ARG:CZ	2.34	0.58
1:B:148:PHE:CE2	1:B:152:VAL:HG23	2.40	0.56
1:A:134:HIS:HE1	2:A:2094:HOH:O	1.90	0.54
1:F:118:LEU:O	1:F:122:THR:HG23	2.07	0.53
1:I:174:SER:HB3	1:I:177:ILE:HG12	1.90	0.53
1:D:181:ARG:HG2	1:D:181:ARG:HH11	1.73	0.53
1:F:148:PHE:CE2	1:F:152:VAL:HG23	2.43	0.53
1:A:135:ARG:HG2	1:F:92:TRP:CE3	2.44	0.53
1:F:174:SER:HB3	1:F:177:ILE:HG12	1.92	0.51
1:A:135:ARG:HG2	1:F:92:TRP:CZ3	2.46	0.49
1:D:148:PHE:CE2	1:D:152:VAL:HG23	2.48	0.49
1:B:118:LEU:O	1:B:122:THR:HG23	2.13	0.49
1:C:73:LYS:O	1:C:74:ASP:HB2	2.14	0.47
1:H:134:HIS:CE1	1:H:141:ARG:HB2	2.49	0.47
1:G:135:ARG:HH21	1:G:160:ASN:HD21	1.63	0.46
1:E:118:LEU:O	1:E:122:THR:HG23	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:148:PHE:CE2	1:E:152:VAL:HG23	2.54	0.43
1:F:134:HIS:CE1	1:F:141:ARG:HB2	2.53	0.43
1:H:81:ARG:NH1	1:H:186:ASN:ND2	2.58	0.43
1:D:181:ARG:CG	1:D:181:ARG:HH11	2.32	0.43
1:C:118:LEU:O	1:C:122:THR:HG23	2.18	0.42
1:B:174:SER:HB3	1:B:177:ILE:HG12	2.01	0.42
1:D:186:ASN:ND2	2:D:2025:HOH:O	2.50	0.42
1:D:122:THR:HB	1:D:127:TYR:CE1	2.55	0.42
1:A:134:HIS:CE1	2:A:2094:HOH:O	2.70	0.42
1:B:81:ARG:NE	1:B:184:GLU:OE2	2.53	0.42
1:C:108:ILE:CD1	1:C:144:ASN:HB3	2.50	0.42
1:I:174:SER:HB3	1:I:177:ILE:CG1	2.51	0.41
1:B:134:HIS:HD2	1:B:136:GLU:OE2	2.04	0.41
1:C:148:PHE:CE2	1:C:152:VAL:HG23	2.56	0.41
1:C:126:LYS:HE3	1:C:166:LEU:HD23	2.02	0.41
1:H:75:TRP:CD2	1:H:84:PHE:HB2	2.56	0.41
1:D:148:PHE:HE2	1:D:152:VAL:HG23	1.84	0.41
1:D:134:HIS:HE1	2:D:2073:HOH:O	2.04	0.41
1:D:108:ILE:HD12	1:D:142:TRP:HB3	2.03	0.40
1:G:78:TYR:HB3	1:G:83:PHE:HE2	1.87	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	116/118 (98%)	108 (93%)	8 (7%)	0	100	100
1	B	116/118 (98%)	110 (95%)	6 (5%)	0	100	100
1	C	116/118 (98%)	107 (92%)	9 (8%)	0	100	100
1	D	116/118 (98%)	111 (96%)	5 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	116/118 (98%)	110 (95%)	6 (5%)	0	100	100
1	F	116/118 (98%)	111 (96%)	5 (4%)	0	100	100
1	G	116/118 (98%)	111 (96%)	5 (4%)	0	100	100
1	H	116/118 (98%)	113 (97%)	3 (3%)	0	100	100
1	I	116/118 (98%)	108 (93%)	8 (7%)	0	100	100
All	All	1044/1062 (98%)	989 (95%)	55 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	106/106 (100%)	106 (100%)	0	100	100
1	B	106/106 (100%)	105 (99%)	1 (1%)	84	83
1	C	106/106 (100%)	105 (99%)	1 (1%)	84	83
1	D	106/106 (100%)	105 (99%)	1 (1%)	84	83
1	E	106/106 (100%)	104 (98%)	2 (2%)	65	59
1	F	106/106 (100%)	105 (99%)	1 (1%)	84	83
1	G	106/106 (100%)	104 (98%)	2 (2%)	65	59
1	H	106/106 (100%)	103 (97%)	3 (3%)	51	41
1	I	106/106 (100%)	103 (97%)	3 (3%)	51	41
All	All	954/954 (100%)	940 (98%)	14 (2%)	72	69

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	70	MET
1	C	123	ASP
1	D	181	ARG
1	E	87	THR

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Mol	Chain	Res	Type
1	E	157	GLN
1	F	157	GLN
1	G	70	MET
1	G	181	ARG
1	H	97	ASP
1	H	116	LYS
1	H	181	ARG
1	I	70	MET
1	I	158	ASN
1	I	181	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	134	HIS
1	B	134	HIS
1	B	157	GLN
1	B	158	ASN
1	C	134	HIS
1	C	160	ASN
1	D	134	HIS
1	D	186	ASN
1	E	134	HIS
1	F	134	HIS
1	F	186	ASN
1	G	154	ASN
1	H	134	HIS
1	H	149	ASN
1	H	186	ASN
1	I	134	HIS
1	I	158	ASN
1	I	160	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	118/118 (100%)	-0.10	2 (1%) 73 76	11, 16, 25, 30	0
1	B	118/118 (100%)	-0.08	1 (0%) 87 88	12, 17, 25, 31	0
1	C	118/118 (100%)	-0.06	5 (4%) 40 44	10, 17, 26, 32	0
1	D	118/118 (100%)	-0.06	2 (1%) 73 76	14, 20, 26, 35	0
1	E	118/118 (100%)	-0.01	3 (2%) 61 64	13, 20, 26, 32	0
1	F	118/118 (100%)	-0.09	2 (1%) 73 76	12, 20, 26, 30	0
1	G	118/118 (100%)	0.12	2 (1%) 73 76	18, 24, 33, 36	0
1	H	118/118 (100%)	0.02	1 (0%) 87 88	16, 22, 30, 35	0
1	I	118/118 (100%)	0.82	13 (11%) 7 8	22, 34, 45, 48	0
All	All	1062/1062 (100%)	0.06	31 (2%) 55 59	10, 20, 35, 48	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	157	GLN	4.4
1	E	98	PHE	4.3
1	I	70	MET	4.1
1	I	149	ASN	3.5
1	D	158	ASN	3.3
1	C	158	ASN	3.1
1	F	98	PHE	3.1
1	G	70	MET	3.0
1	D	157	GLN	3.0
1	I	98	PHE	2.9
1	I	101	GLY	2.8
1	I	147	VAL	2.8
1	C	156	ASN	2.8
1	I	78	TYR	2.8
1	I	151	ASN	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	123	ASP	2.7
1	G	135	ARG	2.6
1	B	157	GLN	2.5
1	A	157	GLN	2.5
1	I	136	GLU	2.4
1	H	139	ARG	2.4
1	I	133	TYR	2.4
1	F	157	GLN	2.3
1	I	158	ASN	2.3
1	I	77	PHE	2.3
1	I	162	ALA	2.2
1	E	158	ASN	2.1
1	A	123	ASP	2.1
1	I	102	LYS	2.1
1	C	98	PHE	2.1
1	E	157	GLN	2.0

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

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

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.