Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2024 – 06:06 PM EDT

PDB ID	:	6FMH
Title	:	Crystal structure of the membrane attack complex assembly inhibitor BGA71 from Lyme disease agent Borreliella bavariensis (Native data)
Authors	:	Brangulis, K.; Akopjana, I.; Petrovskis, I.; Kazaks, A.; Tars, K.
Deposited on	:	2018-01-31
Resolution	:	2.80 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

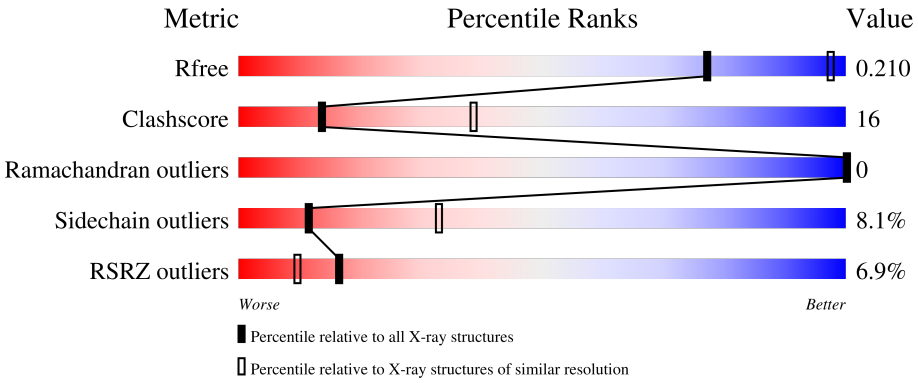
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	194	70% 22% • 6%
1	B	194	61% 27% • • 9%
1	C	194	7% 66% 22% • • 10%
1	D	194	70% 21% • 8%
1	E	194	13% 51% 31% • • 14%

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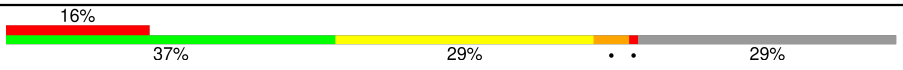
Mol	Chain	Length	Quality of chain
1	F	194	

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 8400 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 membrane attack complex assembly inhibitor BGA71.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	183	Total	C	N	O	S	0	0	0
			1513	954	253	302	4			
1	B	177	Total	C	N	O	S	0	0	0
			1463	925	241	293	4			
1	C	175	Total	C	N	O	S	0	0	0
			1446	915	238	289	4			
1	D	179	Total	C	N	O	S	0	0	0
			1480	934	245	297	4			
1	E	166	Total	C	N	O	S	0	0	0
			1366	863	225	274	4			
1	F	137	Total	C	N	O	S	0	0	0
			1132	719	190	220	3			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLY	-	expression tag	UNP Q6ASF4
A	3	ALA	-	expression tag	UNP Q6ASF4
A	4	MET	-	expression tag	UNP Q6ASF4
A	5	GLY	-	expression tag	UNP Q6ASF4
B	2	GLY	-	expression tag	UNP Q6ASF4
B	3	ALA	-	expression tag	UNP Q6ASF4
B	4	MET	-	expression tag	UNP Q6ASF4
B	5	GLY	-	expression tag	UNP Q6ASF4
C	2	GLY	-	expression tag	UNP Q6ASF4
C	3	ALA	-	expression tag	UNP Q6ASF4
C	4	MET	-	expression tag	UNP Q6ASF4
C	5	GLY	-	expression tag	UNP Q6ASF4
D	2	GLY	-	expression tag	UNP Q6ASF4
D	3	ALA	-	expression tag	UNP Q6ASF4
D	4	MET	-	expression tag	UNP Q6ASF4
D	5	GLY	-	expression tag	UNP Q6ASF4
E	2	GLY	-	expression tag	UNP Q6ASF4

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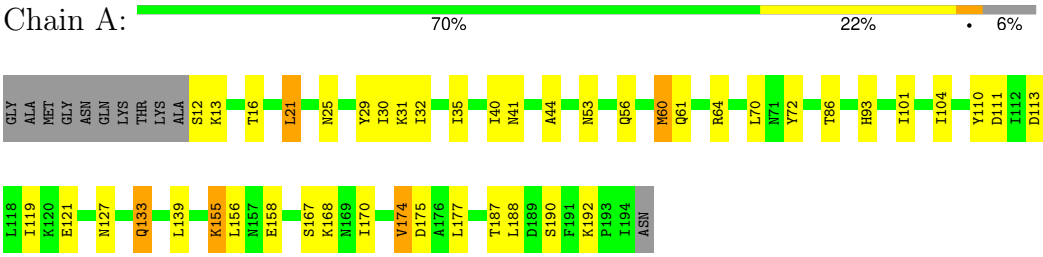
Chain	Residue	Modelled	Actual	Comment	Reference
E	3	ALA	-	expression tag	UNP Q6ASF4
E	4	MET	-	expression tag	UNP Q6ASF4
E	5	GLY	-	expression tag	UNP Q6ASF4
F	2	GLY	-	expression tag	UNP Q6ASF4
F	3	ALA	-	expression tag	UNP Q6ASF4
F	4	MET	-	expression tag	UNP Q6ASF4
F	5	GLY	-	expression tag	UNP Q6ASF4

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

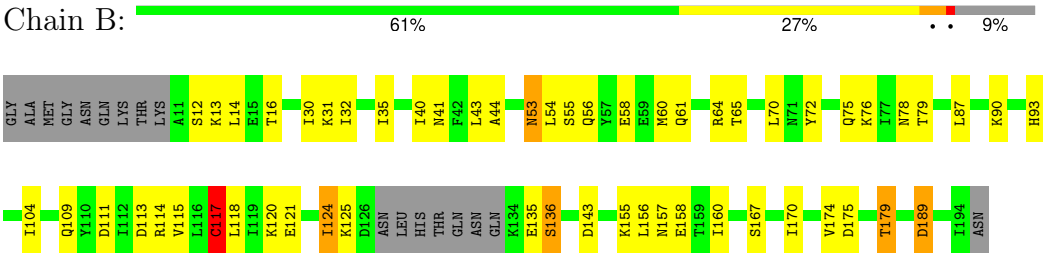
• Molecule 1: membrane attack complex assembly inhibitor BGA71

Chain A:



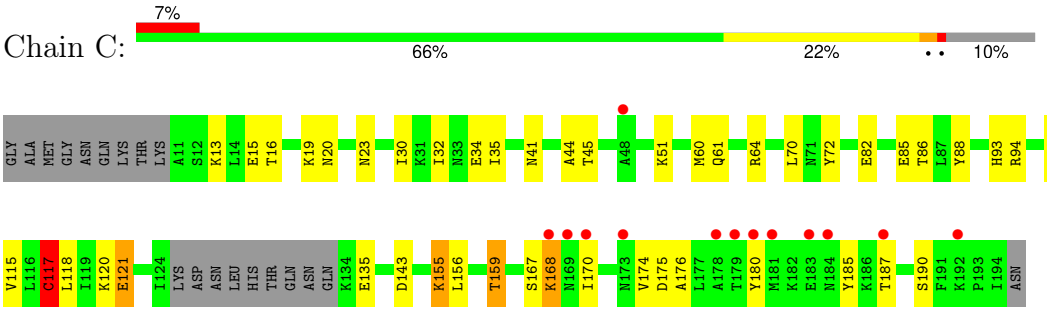
• Molecule 1: membrane attack complex assembly inhibitor BGA71

Chain B:



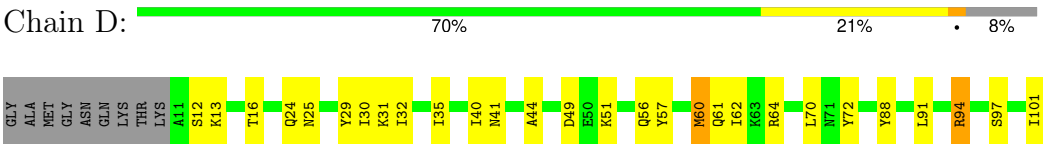
• Molecule 1: membrane attack complex assembly inhibitor BGA71

Chain C:

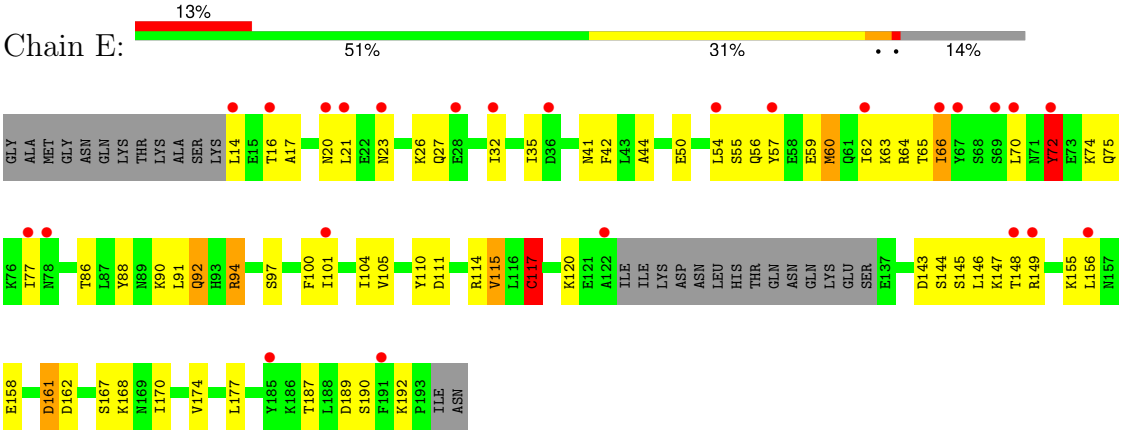


• Molecule 1: membrane attack complex assembly inhibitor BGA71

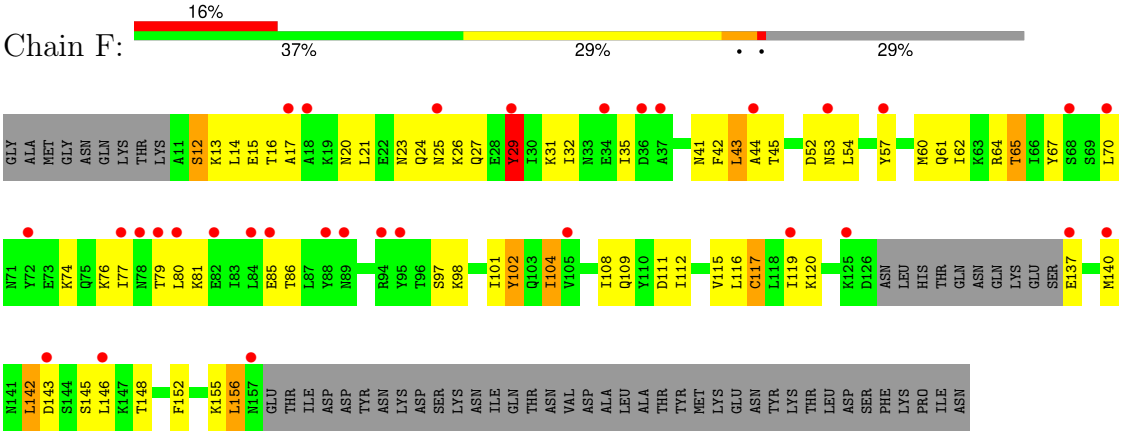
Chain D:



• Molecule 1: membrane attack complex assembly inhibitor BGA71



• Molecule 1: membrane attack complex assembly inhibitor BGA71



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α, β, γ	137.06Å 137.06Å 56.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	68.53 – 2.80 50.95 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.5 (68.53-2.80) 97.5 (50.95-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	7.95 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.190 , 0.216 0.190 , 0.210	Depositor DCC
R_{free} test set	1404 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å^2)	31.1	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 46.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.018 for -h,-k,l 0.159 for h,-h-k,-l 0.017 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	8400	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.64% 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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.87	1/1531 (0.1%)	0.84	1/2057 (0.0%)
1	B	0.96	0/1479	0.88	3/1984 (0.2%)
1	C	0.84	1/1462 (0.1%)	0.86	3/1962 (0.2%)
1	D	0.89	1/1496 (0.1%)	0.82	1/2007 (0.0%)
1	E	0.65	0/1382	0.85	5/1859 (0.3%)
1	F	0.64	0/1143	0.89	3/1531 (0.2%)
All	All	0.83	3/8493 (0.0%)	0.86	16/11400 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	121	GLU	CD-OE1	-6.92	1.18	1.25
1	C	121	GLU	CD-OE1	-5.94	1.19	1.25
1	D	121	GLU	CD-OE2	-5.40	1.19	1.25

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	117	CYS	CA-CB-SG	7.65	127.77	114.00
1	D	94	ARG	CG-CD-NE	6.88	126.25	111.80
1	E	72	TYR	CB-CG-CD1	-6.70	116.98	121.00
1	B	117	CYS	CA-CB-SG	6.41	125.53	114.00
1	F	14	LEU	CA-CB-CG	6.37	129.95	115.30
1	C	143	ASP	CB-CG-OD2	-6.01	112.89	118.30
1	E	14	LEU	CA-CB-CG	5.97	129.03	115.30
1	B	189	ASP	CB-CG-OD2	5.95	123.66	118.30
1	F	142	LEU	CB-CG-CD2	5.95	121.11	111.00
1	B	118	LEU	CB-CG-CD1	-5.64	101.42	111.00
1	F	29	TYR	N-CA-CB	5.49	120.48	110.60
1	E	91	LEU	CA-CB-CG	5.46	127.86	115.30
1	A	21	LEU	CB-CG-CD2	-5.27	102.04	111.00
1	C	143	ASP	CB-CG-OD1	5.21	122.99	118.30
1	E	94	ARG	CG-CD-NE	5.09	122.48	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	117	CYS	CA-CB-SG	5.08	123.15	114.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1513	0	1525	39	1
1	B	1463	0	1479	43	1
1	C	1446	0	1463	45	1
1	D	1480	0	1494	39	1
1	E	1366	0	1357	67	0
1	F	1132	0	1159	66	0
All	All	8400	0	8477	268	2

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 16.

All (268) 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:117:CYS:CB	1:B:117:CYS:SG	2.06	1.42
1:C:117:CYS:SG	1:D:117:CYS:CB	2.11	1.38
1:A:117:CYS:SG	1:B:117:CYS:CB	2.11	1.37
1:C:117:CYS:CB	1:D:117:CYS:SG	2.17	1.31
1:E:117:CYS:SG	1:F:117:CYS:CB	2.30	1.18
1:E:117:CYS:CB	1:F:117:CYS:SG	2.36	1.12
1:E:55:SER:HB3	1:F:15:GLU:CG	1.79	1.12
1:E:55:SER:CB	1:F:15:GLU:HG3	1.84	1.07
1:F:43:LEU:HD23	1:F:54:LEU:HD21	1.38	1.04
1:E:57:TYR:CZ	1:F:120:LYS:HD3	1.93	1.03
1:C:170:ILE:HG22	1:C:180:TYR:CE2	1.95	1.01
1:F:111:ASP:O	1:F:115:VAL:HG23	1.60	1.01
1:F:81:LYS:HE3	1:F:85:GLU:OE2	1.60	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:149:ARG:HH11	1:E:149:ARG:HB3	1.26	0.98
1:C:121:GLU:HG3	1:D:110:TYR:OH	1.67	0.94
1:F:65:THR:HG23	1:F:109:GLN:HG2	1.46	0.93
1:C:19:LYS:HE2	1:D:56:GLN:OE1	1.70	0.91
1:E:55:SER:HB3	1:F:15:GLU:HG3	0.92	0.88
1:C:88:TYR:CE2	1:C:94:ARG:NH1	2.41	0.88
1:E:149:ARG:HB3	1:E:149:ARG:NH1	1.88	0.88
1:E:120:LYS:HD3	1:F:57:TYR:CZ	2.09	0.86
1:E:88:TYR:O	1:E:94:ARG:NH1	2.10	0.84
1:B:61:GLN:NE2	1:B:113:ASP:OD2	2.11	0.84
1:F:116:LEU:O	1:F:120:LYS:HG3	1.78	0.82
1:C:170:ILE:HG22	1:C:180:TYR:CD2	2.14	0.82
1:D:88:TYR:O	1:D:94:ARG:NH1	2.13	0.81
1:C:15:GLU:O	1:C:19:LYS:HG3	1.79	0.81
1:B:53:ASN:HD22	1:B:54:LEU:N	1.76	0.81
1:B:79:THR:OG1	1:B:160:ILE:HD12	1.82	0.79
1:F:26:LYS:HA	1:F:29:TYR:CE2	2.17	0.79
1:A:29:TYR:CD1	1:A:60:MET:CE	2.66	0.79
1:D:61:GLN:HE22	1:D:64:ARG:NH1	1.80	0.79
1:C:19:LYS:CE	1:D:56:GLN:OE1	2.31	0.78
1:A:29:TYR:CD1	1:A:60:MET:HE2	2.19	0.78
1:C:61:GLN:NE2	1:C:113:ASP:OD2	2.17	0.78
1:E:144:SER:O	1:E:148:THR:HG23	1.83	0.77
1:E:56:GLN:HG3	1:E:59:GLU:OE2	1.83	0.77
1:E:170:ILE:HD12	1:E:177:LEU:HA	1.68	0.76
1:D:143:ASP:OD2	1:D:147:LYS:HE2	1.85	0.76
1:A:170:ILE:HD12	1:A:177:LEU:HA	1.68	0.76
1:E:54:LEU:HD21	1:E:59:GLU:HB3	1.67	0.76
1:E:143:ASP:OD2	1:E:147:LYS:HE2	1.85	0.75
1:F:81:LYS:CE	1:F:85:GLU:OE2	2.35	0.74
1:C:176:ALA:O	1:C:180:TYR:CD2	2.41	0.73
1:E:149:ARG:HH11	1:E:149:ARG:CB	1.98	0.73
1:A:61:GLN:NE2	1:A:113:ASP:OD2	2.22	0.73
1:D:30:ILE:HG13	1:D:31:LYS:N	2.03	0.73
1:E:189:ASP:OD1	1:E:192:LYS:HE3	1.88	0.73
1:B:40:ILE:HD13	1:B:78:ASN:HD21	1.53	0.72
1:D:170:ILE:HD12	1:D:177:LEU:HA	1.72	0.72
1:E:56:GLN:O	1:E:59:GLU:HG2	1.90	0.72
1:C:170:ILE:CG2	1:C:180:TYR:CD2	2.71	0.72
1:A:29:TYR:HD1	1:A:60:MET:HE2	1.53	0.71
1:E:23:ASN:O	1:E:27:GLN:OE1	2.08	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:TYR:OH	1:B:121:GLU:HG3	1.91	0.71
1:A:61:GLN:HE22	1:A:64:ARG:NH1	1.87	0.70
1:B:90:LYS:HE3	1:B:174:VAL:HG11	1.73	0.70
1:F:35:ILE:HD11	1:F:67:TYR:CD2	2.27	0.70
1:A:30:ILE:HG13	1:A:31:LYS:N	2.05	0.70
1:C:176:ALA:O	1:C:180:TYR:CE2	2.46	0.69
1:F:23:ASN:O	1:F:27:GLN:OE1	2.08	0.69
1:F:65:THR:CG2	1:F:109:GLN:HG2	2.21	0.69
1:B:14:LEU:CD1	1:B:135:GLU:HG3	2.23	0.69
1:B:30:ILE:HG13	1:B:31:LYS:N	2.06	0.68
1:F:137:GLU:N	1:F:137:GLU:OE1	2.27	0.67
1:A:115:VAL:O	1:A:119:ILE:HG13	1.93	0.67
1:F:42:PHE:O	1:F:81:LYS:HD3	1.93	0.67
1:F:35:ILE:HD11	1:F:67:TYR:HD2	1.59	0.66
1:D:24:GLN:OE1	1:D:24:GLN:HA	1.95	0.65
1:F:108:ILE:HD12	1:F:148:THR:HG22	1.77	0.65
1:A:110:TYR:CZ	1:B:121:GLU:HG3	2.31	0.65
1:E:120:LYS:HD3	1:F:57:TYR:CE1	2.31	0.65
1:D:115:VAL:O	1:D:119:ILE:HG13	1.98	0.64
1:B:93:HIS:O	1:B:96:THR:HG22	1.98	0.64
1:E:66:ILE:HD11	1:E:70:LEU:HD12	1.79	0.63
1:C:45:THR:HG21	1:C:85:GLU:HG2	1.78	0.63
1:A:93:HIS:HE1	1:A:175:ASP:OD1	1.80	0.63
1:C:167:SER:O	1:C:170:ILE:HG23	1.99	0.63
1:F:26:LYS:HA	1:F:29:TYR:CD2	2.33	0.63
1:C:88:TYR:CZ	1:C:94:ARG:NH1	2.67	0.62
1:B:65:THR:OG1	1:B:109:GLN:HG2	1.98	0.62
1:C:93:HIS:HE1	1:C:175:ASP:OD1	1.81	0.62
1:F:52:ASP:OD2	1:F:98:LYS:CE	2.48	0.61
1:F:112:ILE:HD11	1:F:145:SER:O	2.00	0.60
1:A:25:ASN:OD1	1:A:64:ARG:NH2	2.34	0.59
1:E:100:PHE:O	1:E:105:VAL:HG12	2.03	0.59
1:E:161:ASP:OD1	1:E:162:ASP:N	2.36	0.58
1:B:167:SER:O	1:B:170:ILE:HG23	2.04	0.58
1:F:52:ASP:OD2	1:F:98:LYS:HE2	2.03	0.57
1:E:70:LEU:HD12	1:E:77:ILE:HG22	1.86	0.57
1:F:25:ASN:OD1	1:F:64:ARG:NH2	2.37	0.57
1:F:112:ILE:HD11	1:F:145:SER:C	2.24	0.57
1:D:61:GLN:NE2	1:D:113:ASP:OD2	2.38	0.57
1:B:79:THR:OG1	1:B:160:ILE:CD1	2.51	0.56
1:C:45:THR:HG21	1:C:85:GLU:CG	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:101:ILE:HA	1:E:105:VAL:CG1	2.35	0.56
1:E:117:CYS:SG	1:F:117:CYS:SG	0.57	0.56
1:B:175:ASP:O	1:B:179:THR:HG23	2.06	0.56
1:C:176:ALA:HB1	1:C:180:TYR:HE2	1.71	0.56
1:C:88:TYR:CD2	1:C:94:ARG:NH1	2.73	0.56
1:D:88:TYR:CD2	1:D:94:ARG:NH1	2.74	0.56
1:F:76:LYS:O	1:F:79:THR:HG22	2.06	0.56
1:C:120:LYS:HE2	1:D:57:TYR:CE1	2.41	0.55
1:A:31:LYS:HE2	1:C:30:ILE:HD12	1.89	0.55
1:C:121:GLU:HG3	1:D:110:TYR:CZ	2.41	0.55
1:F:26:LYS:HG3	1:F:27:GLN:OE1	2.07	0.55
1:E:26:LYS:HG3	1:E:27:GLN:OE1	2.06	0.55
1:F:45:THR:HG21	1:F:85:GLU:HG2	1.88	0.55
1:E:17:ALA:O	1:E:20:ASN:OD1	2.25	0.55
1:A:29:TYR:CD1	1:A:60:MET:HE3	2.40	0.54
1:D:13:LYS:HA	1:D:16:THR:HG22	1.89	0.54
1:F:70:LEU:HD12	1:F:77:ILE:HG22	1.88	0.54
1:F:108:ILE:HD12	1:F:148:THR:CG2	2.38	0.54
1:B:87:LEU:HD13	1:B:96:THR:HG23	1.88	0.54
1:E:66:ILE:HD11	1:E:70:LEU:CD1	2.37	0.54
1:F:12:SER:O	1:F:15:GLU:HB3	2.06	0.54
1:F:21:LEU:HA	1:F:24:GLN:HG2	1.90	0.54
1:D:49:ASP:OD1	1:D:49:ASP:C	2.43	0.54
1:E:42:PHE:HB3	1:E:63:LYS:HZ3	1.73	0.54
1:B:40:ILE:CD1	1:B:78:ASN:HD21	2.21	0.54
1:F:17:ALA:O	1:F:20:ASN:OD1	2.25	0.54
1:D:40:ILE:O	1:D:40:ILE:HG23	2.07	0.54
1:E:111:ASP:O	1:E:115:VAL:HG22	2.08	0.53
1:E:120:LYS:CD	1:F:57:TYR:CZ	2.90	0.53
1:E:148:THR:HG22	1:E:192:LYS:HA	1.89	0.53
1:F:140:MET:O	1:F:143:ASP:OD1	2.27	0.53
1:E:145:SER:O	1:E:148:THR:OG1	2.26	0.53
1:F:112:ILE:HD11	1:F:146:LEU:HA	1.90	0.53
1:E:57:TYR:CE2	1:F:120:LYS:HD3	2.41	0.53
1:E:59:GLU:HG3	1:E:60:MET:N	2.23	0.53
1:B:61:GLN:O	1:B:109:GLN:HG3	2.09	0.52
1:F:35:ILE:CD1	1:F:67:TYR:CD2	2.92	0.52
1:C:176:ALA:HB1	1:C:180:TYR:CE2	2.45	0.52
1:A:111:ASP:O	1:A:115:VAL:HG23	2.10	0.52
1:A:110:TYR:CZ	1:A:114:ARG:HG3	2.46	0.51
1:E:55:SER:CB	1:F:15:GLU:CG	2.65	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:16:THR:HA	1:C:19:LYS:HD2	1.93	0.51
1:E:101:ILE:HA	1:E:105:VAL:HG12	1.92	0.51
1:A:188:LEU:HG	1:A:192:LYS:NZ	2.25	0.51
1:A:32:ILE:HD11	1:A:64:ARG:HA	1.93	0.50
1:F:20:ASN:O	1:F:23:ASN:OD1	2.30	0.50
1:C:121:GLU:HG3	1:D:110:TYR:HH	1.73	0.50
1:D:61:GLN:NE2	1:D:64:ARG:NH1	2.55	0.50
1:D:25:ASN:OD1	1:D:64:ARG:NH2	2.44	0.50
1:E:146:LEU:HA	1:E:149:ARG:HH12	1.76	0.50
1:A:40:ILE:HG23	1:A:40:ILE:O	2.12	0.50
1:B:93:HIS:HE1	1:B:175:ASP:OD1	1.94	0.50
1:C:32:ILE:HD11	1:C:64:ARG:HA	1.94	0.50
1:C:170:ILE:HG21	1:C:180:TYR:CD2	2.47	0.50
1:E:110:TYR:OH	1:E:114:ARG:HD2	2.12	0.50
1:A:110:TYR:HH	1:B:121:GLU:HG3	1.75	0.49
1:B:87:LEU:HD13	1:B:96:THR:CG2	2.42	0.49
1:E:62:ILE:O	1:E:66:ILE:HG22	2.12	0.49
1:E:20:ASN:O	1:E:23:ASN:OD1	2.30	0.49
1:F:23:ASN:HA	1:F:26:LYS:HE2	1.94	0.49
1:F:41:ASN:HB3	1:F:44:ALA:HB2	1.95	0.49
1:A:127:ASN:H	1:A:127:ASN:ND2	2.11	0.48
1:C:170:ILE:CG2	1:C:180:TYR:CE2	2.82	0.48
1:A:110:TYR:OH	1:A:114:ARG:HG3	2.14	0.48
1:D:25:ASN:O	1:D:29:TYR:HD2	1.95	0.48
1:D:35:ILE:HD11	1:D:72:TYR:HB3	1.95	0.48
1:C:35:ILE:HD11	1:C:72:TYR:HB3	1.96	0.48
1:C:176:ALA:C	1:C:180:TYR:CE2	2.87	0.48
1:F:32:ILE:HD11	1:F:64:ARG:HA	1.96	0.48
1:F:20:ASN:OD1	1:F:21:LEU:N	2.47	0.47
1:D:29:TYR:CE1	1:D:60:MET:HE2	2.49	0.47
1:D:88:TYR:CE2	1:D:94:ARG:NH1	2.82	0.47
1:B:53:ASN:HD22	1:B:53:ASN:C	2.11	0.47
1:B:76:LYS:HA	1:B:160:ILE:HD11	1.96	0.47
1:E:32:ILE:HD11	1:E:64:ARG:HA	1.96	0.47
1:A:110:TYR:CD1	1:B:124:ILE:HD12	2.49	0.47
1:C:13:LYS:HD2	1:C:135:GLU:OE2	2.14	0.47
1:B:76:LYS:HA	1:B:160:ILE:CD1	2.45	0.47
1:B:35:ILE:HD11	1:B:72:TYR:HB3	1.97	0.47
1:F:43:LEU:CD2	1:F:54:LEU:HD21	2.27	0.47
1:C:115:VAL:O	1:C:115:VAL:HG22	2.15	0.46
1:E:20:ASN:OD1	1:E:21:LEU:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:66:ILE:CD1	1:E:70:LEU:HD12	2.44	0.46
1:C:159:THR:HG21	1:C:185:TYR:OH	2.14	0.46
1:C:70:LEU:HD21	1:C:156:LEU:HG	1.98	0.46
1:A:61:GLN:NE2	1:A:64:ARG:NH1	2.60	0.46
1:E:101:ILE:O	1:E:105:VAL:CG1	2.63	0.46
1:C:155:LYS:O	1:C:159:THR:HG23	2.16	0.46
1:D:32:ILE:HD11	1:D:64:ARG:HA	1.98	0.46
1:E:74:LYS:HA	1:E:77:ILE:HG12	1.98	0.46
1:B:40:ILE:HD13	1:B:78:ASN:ND2	2.27	0.46
1:F:76:LYS:HA	1:F:79:THR:HG22	1.98	0.46
1:A:167:SER:O	1:A:168:LYS:HB2	2.15	0.45
1:B:120:LYS:O	1:B:124:ILE:CG1	2.64	0.45
1:D:94:ARG:HH11	1:D:94:ARG:HG3	1.81	0.45
1:A:25:ASN:O	1:A:29:TYR:HD2	1.99	0.45
1:C:86:THR:CG2	1:C:174:VAL:HG12	2.46	0.45
1:A:170:ILE:HG23	1:A:177:LEU:CD1	2.46	0.45
1:D:29:TYR:CD1	1:D:60:MET:HE1	2.51	0.45
1:B:70:LEU:HD21	1:B:156:LEU:HG	1.99	0.45
1:F:115:VAL:O	1:F:119:ILE:HG13	2.17	0.45
1:F:61:GLN:O	1:F:109:GLN:HG3	2.17	0.45
1:B:32:ILE:HD11	1:B:64:ARG:HA	1.99	0.45
1:D:167:SER:O	1:D:168:LYS:HB2	2.17	0.45
1:E:189:ASP:HA	1:E:192:LYS:HE3	1.99	0.45
1:E:54:LEU:HD23	1:E:54:LEU:O	2.17	0.45
1:E:167:SER:O	1:E:168:LYS:HB2	2.16	0.45
1:A:187:THR:HG23	1:A:190:SER:H	1.82	0.45
1:B:40:ILE:CD1	1:B:78:ASN:ND2	2.80	0.45
1:F:31:LYS:O	1:F:35:ILE:HG12	2.17	0.45
1:A:86:THR:HG22	1:A:174:VAL:HG13	1.99	0.44
1:C:20:ASN:O	1:C:23:ASN:OD1	2.35	0.44
1:E:70:LEU:HD21	1:E:156:LEU:HG	1.98	0.44
1:E:57:TYR:OH	1:F:120:LYS:HD3	2.13	0.44
1:C:187:THR:HG23	1:C:190:SER:H	1.82	0.44
1:F:62:ILE:HD12	1:F:101:ILE:HG23	1.99	0.44
1:C:159:THR:CG2	1:C:185:TYR:OH	2.66	0.44
1:A:70:LEU:HD21	1:A:156:LEU:HG	1.98	0.44
1:E:88:TYR:CD2	1:E:94:ARG:NH1	2.86	0.44
1:A:110:TYR:OH	1:B:121:GLU:CG	2.62	0.44
1:D:41:ASN:HB3	1:D:44:ALA:HB2	2.00	0.44
1:E:41:ASN:HB3	1:E:44:ALA:HB2	1.98	0.44
1:F:104:ILE:O	1:F:108:ILE:HG12	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:72:TYR:CG	1:E:72:TYR:O	2.70	0.44
1:B:101:ILE:HD13	1:B:101:ILE:HA	1.87	0.43
1:A:35:ILE:HD11	1:A:72:TYR:HB3	1.99	0.43
1:B:40:ILE:HG23	1:B:40:ILE:O	2.18	0.43
1:C:117:CYS:HG	1:D:117:CYS:CB	2.00	0.43
1:E:23:ASN:HA	1:E:26:LYS:HE2	2.00	0.43
1:E:56:GLN:O	1:E:59:GLU:CG	2.64	0.43
1:E:56:GLN:C	1:E:59:GLU:HG2	2.38	0.43
1:D:94:ARG:O	1:D:97:SER:OG	2.35	0.43
1:E:90:LYS:HD2	1:E:92:GLN:NE2	2.33	0.43
1:F:52:ASP:OD2	1:F:98:LYS:NZ	2.49	0.43
1:F:152:PHE:CZ	1:F:156:LEU:HG	2.53	0.43
1:C:117:CYS:CB	1:D:117:CYS:HG	2.02	0.43
1:D:25:ASN:O	1:D:29:TYR:CD2	2.71	0.43
1:C:41:ASN:HB3	1:C:44:ALA:HB2	1.99	0.43
1:D:170:ILE:HG23	1:D:177:LEU:CD1	2.48	0.43
1:E:101:ILE:O	1:E:105:VAL:HG13	2.17	0.43
1:B:43:LEU:HD13	1:B:54:LEU:HD21	2.00	0.43
1:B:93:HIS:CE1	1:B:175:ASP:OD1	2.72	0.43
1:A:155:LYS:NZ	1:A:158:GLU:OE1	2.42	0.43
1:F:74:LYS:HA	1:F:77:ILE:HG12	2.00	0.43
1:A:41:ASN:HB3	1:A:44:ALA:HB2	2.01	0.43
1:E:187:THR:HG23	1:E:190:SER:H	1.83	0.43
1:B:41:ASN:HB3	1:B:44:ALA:HB2	1.99	0.42
1:E:170:ILE:HG23	1:E:177:LEU:CD1	2.49	0.42
1:F:26:LYS:CG	1:F:27:GLN:OE1	2.67	0.42
1:D:70:LEU:HD21	1:D:156:LEU:HG	1.99	0.42
1:B:111:ASP:O	1:B:115:VAL:HG23	2.18	0.42
1:A:101:ILE:HD13	1:A:101:ILE:HA	1.88	0.42
1:A:25:ASN:O	1:A:29:TYR:CD2	2.72	0.42
1:B:55:SER:OG	1:B:58:GLU:HG3	2.20	0.42
1:E:26:LYS:CG	1:E:27:GLN:OE1	2.68	0.42
1:F:76:LYS:O	1:F:79:THR:CG2	2.67	0.42
1:A:13:LYS:HD3	1:A:139:LEU:CD1	2.49	0.42
1:B:13:LYS:HD2	1:B:135:GLU:OE2	2.20	0.42
1:B:75:GLN:O	1:B:79:THR:HG23	2.19	0.41
1:E:86:THR:CG2	1:E:174:VAL:HG12	2.50	0.41
1:F:146:LEU:O	1:F:146:LEU:HD23	2.20	0.41
1:E:35:ILE:HG13	1:E:72:TYR:HD1	1.86	0.41
1:F:62:ILE:HD11	1:F:101:ILE:O	2.21	0.41
1:F:102:TYR:O	1:F:102:TYR:CD1	2.74	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:62:ILE:HD13	1:D:101:ILE:HG23	2.03	0.41
1:C:168:LYS:HD3	1:C:180:TYR:CE1	2.56	0.41
1:C:170:ILE:HG22	1:C:180:TYR:HE2	1.69	0.41
1:E:59:GLU:CG	1:E:60:MET:N	2.84	0.41
1:E:146:LEU:HA	1:E:149:ARG:NH1	2.35	0.41
1:D:24:GLN:OE1	1:D:24:GLN:CA	2.65	0.41
1:F:156:LEU:HD23	1:F:156:LEU:HA	1.95	0.41
1:B:157:ASN:HA	1:B:160:ILE:HG22	2.03	0.40
1:F:62:ILE:CD1	1:F:101:ILE:O	2.69	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:GLN:O	1:B:136:SER:OG[2_885]	1.76	0.44
1:C:175:ASP:OD2	1:D:51:LYS:NZ[3_595]	1.83	0.37

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	181/194 (93%)	175 (97%)	6 (3%)	0	100	100
1	B	173/194 (89%)	168 (97%)	5 (3%)	0	100	100
1	C	171/194 (88%)	168 (98%)	3 (2%)	0	100	100
1	D	175/194 (90%)	168 (96%)	7 (4%)	0	100	100
1	E	162/194 (84%)	157 (97%)	5 (3%)	0	100	100
1	F	133/194 (69%)	128 (96%)	5 (4%)	0	100	100
All	All	995/1164 (86%)	964 (97%)	31 (3%)	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	171/179 (96%)	160 (94%)	11 (6%)	17 45
1	B	165/179 (92%)	149 (90%)	16 (10%)	8 24
1	C	163/179 (91%)	153 (94%)	10 (6%)	18 48
1	D	167/179 (93%)	160 (96%)	7 (4%)	30 63
1	E	152/179 (85%)	137 (90%)	15 (10%)	8 23
1	F	126/179 (70%)	109 (86%)	17 (14%)	4 11
All	All	944/1074 (88%)	868 (92%)	76 (8%)	11 33

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	SER
1	A	16	THR
1	A	21	LEU
1	A	53	ASN
1	A	56	GLN
1	A	60	MET
1	A	104	ILE
1	A	117	CYS
1	A	133	GLN
1	A	155	LYS
1	A	174	VAL
1	B	12	SER
1	B	16	THR
1	B	53	ASN
1	B	56	GLN
1	B	60	MET
1	B	104	ILE
1	B	114	ARG
1	B	117	CYS
1	B	124	ILE
1	B	125	LYS
1	B	136	SER

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Mol	Chain	Res	Type
1	B	143	ASP
1	B	155	LYS
1	B	158	GLU
1	B	179	THR
1	B	189	ASP
1	C	34	GLU
1	C	51	LYS
1	C	60	MET
1	C	82	GLU
1	C	104	ILE
1	C	117	CYS
1	C	118	LEU
1	C	155	LYS
1	C	159	THR
1	C	168	LYS
1	D	12	SER
1	D	60	MET
1	D	91	LEU
1	D	104	ILE
1	D	114	ARG
1	D	117	CYS
1	D	194	ILE
1	E	16	THR
1	E	50	GLU
1	E	60	MET
1	E	65	THR
1	E	66	ILE
1	E	72	TYR
1	E	75	GLN
1	E	92	GLN
1	E	97	SER
1	E	104	ILE
1	E	115	VAL
1	E	117	CYS
1	E	155	LYS
1	E	158	GLU
1	E	161	ASP
1	F	12	SER
1	F	13	LYS
1	F	16	THR
1	F	29	TYR
1	F	43	LEU

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Mol	Chain	Res	Type
1	F	53	ASN
1	F	60	MET
1	F	65	THR
1	F	80	LEU
1	F	86	THR
1	F	97	SER
1	F	102	TYR
1	F	104	ILE
1	F	117	CYS
1	F	142	LEU
1	F	155	LYS
1	F	156	LEU

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

Mol	Chain	Res	Type
1	A	56	GLN
1	A	93	HIS
1	A	127	ASN
1	A	151	ASN
1	B	20	ASN
1	B	23	ASN
1	B	53	ASN
1	B	78	ASN
1	B	93	HIS
1	B	109	GLN
1	C	41	ASN
1	C	92	GLN
1	C	93	HIS
1	C	151	ASN
1	D	20	ASN
1	D	151	ASN
1	E	24	GLN
1	E	92	GLN
1	E	151	ASN
1	F	41	ASN
1	F	53	ASN
1	F	56	GLN
1	F	109	GLN
1	F	151	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 monosaccharides 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	183/194 (94%)	-0.51	0 100 100	9, 22, 45, 83	0
1	B	177/194 (91%)	-0.51	0 100 100	6, 21, 50, 76	0
1	C	175/194 (90%)	-0.04	13 (7%) 14 8	7, 25, 95, 125	0
1	D	179/194 (92%)	-0.50	0 100 100	6, 20, 52, 86	0
1	E	166/194 (85%)	0.58	25 (15%) 2 1	28, 60, 101, 119	0
1	F	137/194 (70%)	1.24	32 (23%) 0 0	41, 79, 116, 144	0
All	All	1017/1164 (87%)	-0.02	70 (6%) 16 10	6, 29, 96, 144	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	80	LEU	7.3
1	F	146	LEU	7.0
1	E	23	ASN	6.3
1	E	20	ASN	5.3
1	F	37	ALA	5.0
1	F	95	TYR	4.7
1	F	36	ASP	4.7
1	F	140	MET	4.7
1	E	66	ILE	4.6
1	C	180	TYR	4.4
1	F	78	ASN	4.4
1	F	44	ALA	4.4
1	F	84	LEU	4.3
1	F	29	TYR	3.9
1	E	191	PHE	3.8
1	F	105	VAL	3.8
1	E	14	LEU	3.6
1	E	149	ARG	3.5
1	F	18	ALA	3.4

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Mol	Chain	Res	Type	RSRZ
1	F	70	LEU	3.4
1	C	184	ASN	3.4
1	C	181	MET	3.4
1	E	101	ILE	3.3
1	C	168	LYS	3.3
1	E	77	ILE	3.2
1	F	82	GLU	3.1
1	F	57	TYR	3.1
1	C	173	ASN	3.1
1	F	68	SER	3.1
1	F	34	GLU	3.0
1	F	77	ILE	3.0
1	F	72	TYR	3.0
1	F	85	GLU	3.0
1	F	119	ILE	2.9
1	E	32	ILE	2.8
1	E	148	THR	2.8
1	E	57	TYR	2.8
1	C	192	LYS	2.8
1	C	187	THR	2.8
1	E	70	LEU	2.8
1	E	72	TYR	2.8
1	E	122	ALA	2.8
1	F	157	ASN	2.6
1	C	178	ALA	2.6
1	E	36	ASP	2.5
1	F	143	ASP	2.5
1	F	17	ALA	2.5
1	E	21	LEU	2.5
1	C	183	GLU	2.5
1	F	94	ARG	2.4
1	C	179	THR	2.4
1	E	156	LEU	2.4
1	C	170	ILE	2.3
1	F	89	ASN	2.3
1	C	169	ASN	2.3
1	E	69	SER	2.3
1	F	88	TYR	2.3
1	F	137	GLU	2.3
1	E	67	TYR	2.2
1	E	78	ASN	2.2
1	E	16	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	54	LEU	2.2
1	E	62	ILE	2.2
1	F	79	THR	2.1
1	F	53	ASN	2.1
1	F	125	LYS	2.1
1	E	185	TYR	2.1
1	C	48	ALA	2.0
1	E	28	GLU	2.0
1	F	25	ASN	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.