



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:44 PM UTC

PDB ID : 6FMH / pdb_00006fmh
Title : Crystal structure of the membrane attack complex assembly inhibitor BGA71 from Lyme disease agent *Borrelia burgdorferi* (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:

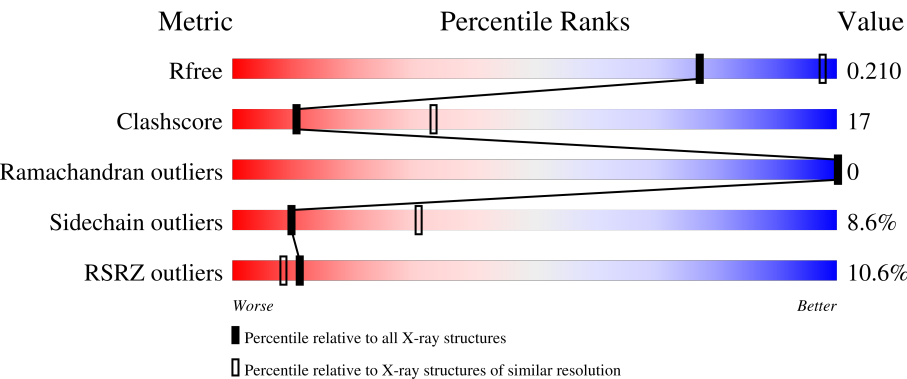
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	194	8% 66%24%6%
1	B	194	2% 60%25%5%9%
1	C	194	8% 66%21%10%
1	D	194	3% 69%21%8%
1	E	194	19% 51%30%14%

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Mol	Chain	Length	Quality of chain
1	F	194	24% 35% 30% 2% 29%

2 Entry composition

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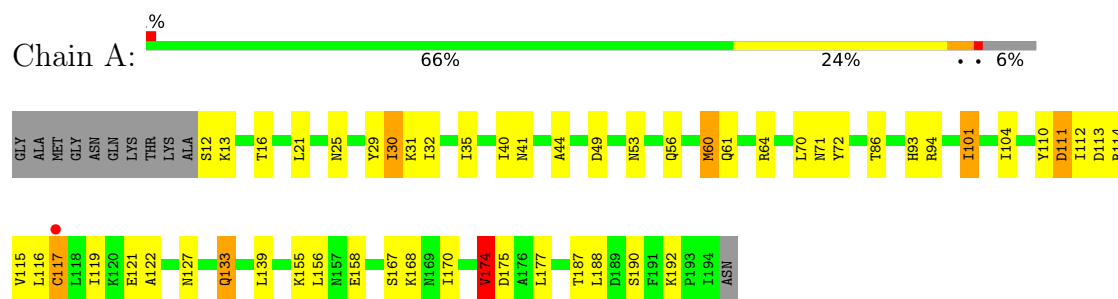
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Chain	Residue	Modelled	Actual	Comment	Reference
E	3	ALA	-	expression tag	UNP Q6ASF4
E	4	MET	-	expression tag	UNP Q6ASF4
E	5	GLY	-	expression tag	UNP Q6ASF4
F	2	GLY	-	expression tag	UNP Q6ASF4
F	3	ALA	-	expression tag	UNP Q6ASF4
F	4	MET	-	expression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

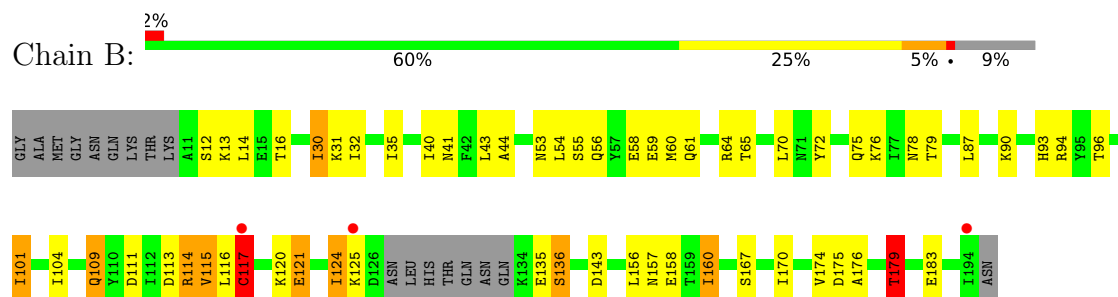
3 Residue-property plots

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.

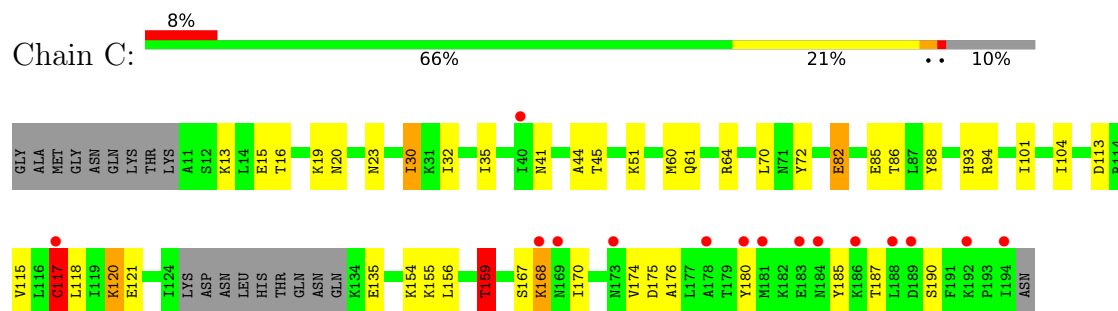
- Molecule 1: membrane attack complex assembly inhibitor BGA71



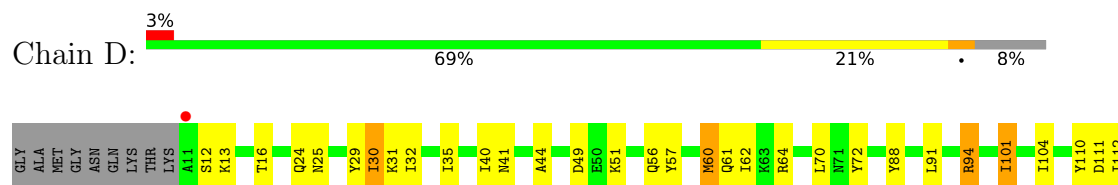
- Molecule 1: membrane attack complex assembly inhibitor BGA71

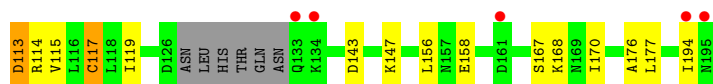


- Molecule 1: membrane attack complex assembly inhibitor BGA71

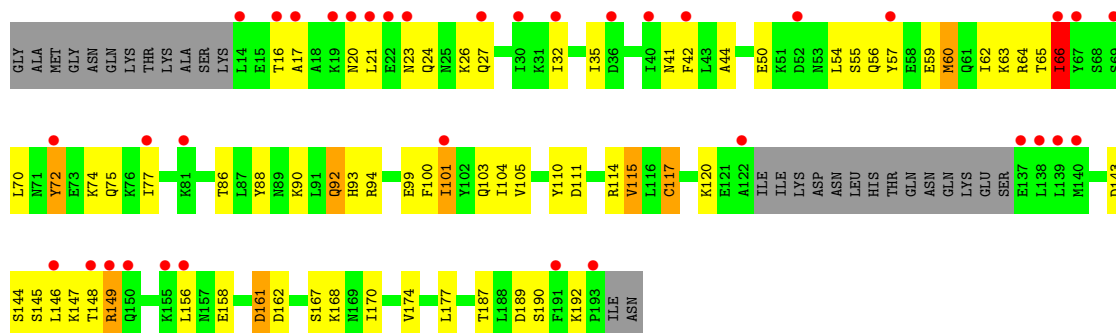


- Molecule 1: membrane attack complex assembly inhibitor BGA71

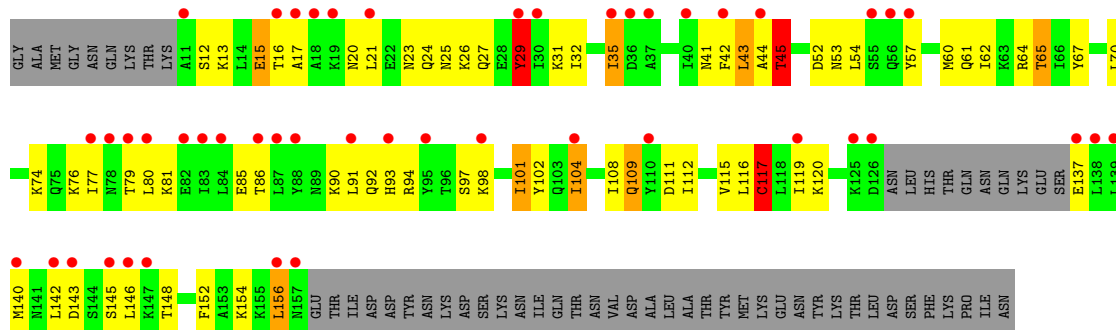




- Molecule 1: membrane attack complex assembly inhibitor BGA71



- Molecule 1: membrane attack complex assembly inhibitor BGA71



4 Data and refinement statistics

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 68.53 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.5 (68.53-2.80) 97.5 (68.53-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ 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 (4.82%)	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)	0.33 , 38.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	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.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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	1.15	5/1531 (0.3%)	1.11	2/2057 (0.1%)
1	B	1.23	7/1479 (0.5%)	1.11	3/1984 (0.2%)
1	C	1.12	1/1462 (0.1%)	1.12	6/1962 (0.3%)
1	D	1.19	3/1496 (0.2%)	1.10	2/2007 (0.1%)
1	E	0.88	0/1382	1.05	1/1859 (0.1%)
1	F	0.91	0/1143	1.10	6/1531 (0.4%)
All	All	1.10	16/8493 (0.2%)	1.10	20/11400 (0.2%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	111	ASP	C-O	-6.34	1.16	1.24
1	D	113	ASP	C-O	-6.14	1.17	1.24
1	A	121	GLU	C-O	-5.87	1.17	1.24
1	A	122	ALA	C-O	-5.87	1.17	1.24
1	C	120	LYS	C-O	-5.57	1.17	1.24
1	D	112	ILE	C-O	-5.57	1.17	1.24
1	B	59	GLU	C-O	-5.52	1.17	1.24
1	B	121	GLU	C-O	-5.32	1.17	1.24
1	A	116	LEU	C-O	-5.19	1.17	1.24
1	B	115	VAL	C-O	-5.18	1.18	1.24
1	B	160	ILE	CA-CB	-5.18	1.48	1.54
1	B	116	LEU	C-O	-5.15	1.18	1.24
1	A	111	ASP	C-O	-5.13	1.18	1.24
1	B	183	GLU	CA-C	5.09	1.59	1.52
1	B	117	CYS	C-O	-5.03	1.17	1.24
1	A	112	ILE	C-O	-5.03	1.18	1.24

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	VAL	CB-CA-C	-7.29	102.31	112.14
1	F	29	TYR	N-CA-CB	7.07	120.48	109.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	94	ARG	CG-CD-NE	6.48	126.25	112.00
1	C	30	ILE	N-CA-CB	-6.44	103.02	110.55
1	C	117	CYS	N-CA-C	-6.33	104.43	111.71
1	C	117	CYS	CA-CB-SG	5.81	127.77	114.40
1	F	35	ILE	CB-CA-C	5.78	119.28	111.88
1	C	159	THR	N-CA-CB	5.75	118.67	110.16
1	F	117	CYS	N-CA-C	-5.64	105.27	111.82
1	B	114	ARG	N-CA-C	5.63	117.87	111.11
1	F	45	THR	CB-CA-C	-5.62	101.13	110.68
1	B	176	ALA	N-CA-C	-5.61	105.25	111.36
1	F	15	GLU	N-CA-CB	-5.35	102.25	110.01
1	A	174	VAL	N-CA-CB	5.27	117.71	110.54
1	F	154	LYS	N-CA-CB	5.13	117.59	109.94
1	E	66	ILE	N-CA-C	5.09	115.32	110.53
1	D	176	ALA	N-CA-C	-5.08	105.74	111.28
1	B	179	THR	CB-CA-C	5.06	119.45	110.85
1	C	82	GLU	CB-CG-CD	5.05	121.18	112.60
1	C	154	LYS	N-CA-C	-5.00	105.72	111.07

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	1513	0	1525	43	1
1	B	1463	0	1479	44	1
1	C	1446	0	1463	45	1
1	D	1480	0	1494	38	1
1	E	1366	0	1357	72	0
1	F	1132	0	1159	70	0
All	All	8400	0	8477	283	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 17.

All (283) 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
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:A:29:TYR:CD1	1:A:60:MET:HE3	2.14	0.83
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:143:ASP:OD2	1:E:147:LYS:HE2	1.85	0.75
1:D:29:TYR:CD1	1:D:60:MET:HE2	2.22	0.74
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:D:29:TYR:CE1	1:D:60:MET:HE2	2.24	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:E:23:ASN:O	1:E:27:GLN:OE1	2.08	0.71
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:93:HIS:HE1	1:C:175:ASP:OD1	1.81	0.62
1:F:112:ILE:HD11	1:F:145:SER:C	2.24	0.61
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:D:49:ASP:OD1	1:D:49:ASP:C	2.43	0.58
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:D:61:GLN:NE2	1:D:113:ASP:OD2	2.38	0.57
1:A:29:TYR:CE1	1:A:60:MET:HE3	2.39	0.56
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
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: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: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:A:29:TYR:HD1	1:A:60:MET:HE2	1.72	0.54
1:D:40:ILE:O	1:D:40:ILE:HG23	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:42:PHE:HB3	1:E:63:LYS:HZ1	1.74	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:C:176:ALA:C	1:C:180:TYR:CE2	2.87	0.52
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:29:TYR:HD1	1:A:60:MET:CE	2.19	0.51
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
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: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:53:ASN:HD22	1:B:53:ASN:C	2.11	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: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:E:56:GLN:C	1:E:59:GLU:HG2	2.38	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:88:TYR:CE2	1:D:94:ARG:NH1	2.82	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
1:E:66:ILE:CD1	1:E:70:LEU:HD12	2.44	0.46
1:A:155:LYS:NZ	1:A:158:GLU:OE1	2.42	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: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: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: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:B:93:HIS:O	1:B:94:ARG:C	2.58	0.45
1:F:31:LYS:O	1:F:35:ILE:HG12	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:167:SER:O	1:A:168:LYS:HB2	2.15	0.44
1:C:187:THR:HG23	1:C:190:SER:H	1.82	0.44
1:E:99:GLU:HG2	1:E:103:GLN:CD	2.42	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
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:D:61:GLN:CD	1:D:64:ARG:HH11	2.26	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:D:167:SER:O	1:D:168:LYS:HB2	2.17	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: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:E:56:GLN:HA	1:E:59:GLU:CD	2.43	0.43
1:F:25:ASN:CG	1:F:64:ARG:NH2	2.77	0.43
1:C:93:HIS:O	1:C:94:ARG:C	2.59	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:29:TYR:HA	1:A:60:MET:HE1	2.01	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:A:93:HIS:O	1:A:94:ARG:C	2.60	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:161:ASP:OD1	1:E:161:ASP:C	2.62	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:E:143:ASP:CG	1:E:147:LYS:HE2	2.45	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:F:93:HIS:O	1:F:94:ARG:C	2.63	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:F:92:GLN:H	1:F:92:GLN:CD	2.28	0.41
1:B:53:ASN:C	1:B:53:ASN:ND2	2.75	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
1:D:62:ILE:HD13	1:D:101:ILE:HG23	2.03	0.41
1:A:49:ASP:OD1	1:A:49:ASP:C	2.62	0.41
1:A:61:GLN:CD	1:A:64:ARG:HH11	2.29	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:F:90:LYS:O	1:F:91:LEU:C	2.64	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:E:93:HIS:O	1:E:94:ARG:C	2.64	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%)	158 (92%)	13 (8%)	12	36
1	B	165/179 (92%)	149 (90%)	16 (10%)	8	25
1	C	163/179 (91%)	154 (94%)	9 (6%)	19	51

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	167/179 (93%)	157 (94%)	10 (6%)	17	47
1	E	152/179 (85%)	136 (90%)	16 (10%)	6	22
1	F	126/179 (70%)	109 (86%)	17 (14%)	4	13
All	All	944/1074 (88%)	863 (91%)	81 (9%)	10	31

All (81) 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	30	ILE
1	A	53	ASN
1	A	56	GLN
1	A	60	MET
1	A	71	ASN
1	A	101	ILE
1	A	104	ILE
1	A	117	CYS
1	A	133	GLN
1	A	174	VAL
1	B	12	SER
1	B	16	THR
1	B	30	ILE
1	B	56	GLN
1	B	60	MET
1	B	101	ILE
1	B	104	ILE
1	B	109	GLN
1	B	114	ARG
1	B	117	CYS
1	B	124	ILE
1	B	125	LYS
1	B	136	SER
1	B	143	ASP
1	B	158	GLU
1	B	179	THR
1	C	51	LYS
1	C	60	MET
1	C	82	GLU
1	C	101	ILE

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Mol	Chain	Res	Type
1	C	104	ILE
1	C	117	CYS
1	C	118	LEU
1	C	159	THR
1	C	168	LYS
1	D	12	SER
1	D	30	ILE
1	D	60	MET
1	D	91	LEU
1	D	101	ILE
1	D	104	ILE
1	D	114	ARG
1	D	117	CYS
1	D	158	GLU
1	D	194	ILE
1	E	16	THR
1	E	24	GLN
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	101	ILE
1	E	104	ILE
1	E	115	VAL
1	E	117	CYS
1	E	149	ARG
1	E	158	GLU
1	E	161	ASP
1	F	13	LYS
1	F	16	THR
1	F	29	TYR
1	F	43	LEU
1	F	45	THR
1	F	53	ASN
1	F	60	MET
1	F	65	THR
1	F	80	LEU
1	F	86	THR
1	F	97	SER

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Mol	Chain	Res	Type
1	F	101	ILE
1	F	104	ILE
1	F	109	GLN
1	F	117	CYS
1	F	142	LEU
1	F	156	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	78	ASN
1	A	93	HIS
1	A	127	ASN
1	A	131	GLN
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	C	169	ASN
1	D	20	ASN
1	D	41	ASN
1	D	151	ASN
1	E	24	GLN
1	E	33	ASN
1	E	41	ASN
1	E	78	ASN
1	E	92	GLN
1	E	109	GLN
1	E	151	ASN
1	E	169	ASN
1	F	41	ASN
1	F	53	ASN
1	F	56	GLN
1	F	61	GLN

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Mol	Chain	Res	Type
1	F	78	ASN
1	F	109	GLN
1	F	151	ASN

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 oligosaccharides 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

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, 95th 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%)	-0.23	1 (0%) 87 82	9, 22, 45, 83	0
1	B	177/194 (91%)	-0.22	3 (1%) 69 60	6, 21, 50, 76	0
1	C	175/194 (90%)	0.16	15 (8%) 16 12	7, 25, 95, 125	0
1	D	179/194 (92%)	-0.23	6 (3%) 48 39	6, 20, 52, 86	0
1	E	166/194 (85%)	1.09	36 (21%) 2 2	28, 60, 101, 119	0
1	F	137/194 (70%)	1.60	47 (34%) 1 1	41, 79, 116, 144	0
All	All	1017/1164 (87%)	0.30	108 (10%) 11 8	6, 29, 96, 144	0

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	23	ASN	5.1
1	E	193	PRO	4.8
1	E	20	ASN	4.7
1	F	139	LEU	4.5
1	F	80	LEU	4.3
1	C	194	ILE	4.3
1	F	11	ALA	4.1
1	E	122	ALA	4.1
1	F	138	LEU	4.0
1	E	40	ILE	4.0
1	F	140	MET	3.9
1	F	146	LEU	3.8
1	F	126	ASP	3.8
1	F	42	PHE	3.7
1	E	66	ILE	3.7
1	F	29	TYR	3.7
1	F	57	TYR	3.6
1	E	101	ILE	3.5
1	F	30	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	E	57	TYR	3.4
1	F	40	ILE	3.3
1	C	181	MET	3.3
1	C	186	LYS	3.2
1	B	117	CYS	3.2
1	E	156	LEU	3.2
1	D	194	ILE	3.1
1	C	180	TYR	3.1
1	E	14	LEU	3.1
1	F	56	GLN	3.0
1	B	125	LYS	3.0
1	F	95	TYR	3.0
1	F	125	LYS	2.9
1	D	133	GLN	2.9
1	E	17	ALA	2.9
1	F	36	ASP	2.9
1	F	86	THR	2.9
1	F	142	LEU	2.9
1	F	84	LEU	2.9
1	A	117	CYS	2.8
1	F	17	ALA	2.8
1	F	110	TYR	2.8
1	E	191	PHE	2.8
1	C	183	GLU	2.8
1	E	138	LEU	2.8
1	E	149	ARG	2.7
1	F	137	GLU	2.7
1	C	188	LEU	2.7
1	B	194	ILE	2.7
1	C	117	CYS	2.7
1	F	157	ASN	2.7
1	F	87	LEU	2.7
1	E	148	THR	2.7
1	D	11	ALA	2.6
1	E	19	LYS	2.6
1	F	143	ASP	2.6
1	D	134	LYS	2.6
1	C	40	ILE	2.6
1	E	16	THR	2.6
1	E	139	LEU	2.5
1	F	83	ILE	2.5
1	F	37	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	E	81	LYS	2.5
1	E	77	ILE	2.5
1	F	79	THR	2.5
1	F	78	ASN	2.5
1	F	77	ILE	2.4
1	C	178	ALA	2.4
1	F	35	ILE	2.4
1	F	147	LYS	2.4
1	F	16	THR	2.4
1	E	22	GLU	2.3
1	F	156	LEU	2.3
1	F	19	LYS	2.3
1	F	145	SER	2.3
1	E	42	PHE	2.2
1	C	184	ASN	2.2
1	E	27	GLN	2.2
1	F	18	ALA	2.2
1	E	52	ASP	2.2
1	C	173	ASN	2.2
1	E	146	LEU	2.2
1	F	91	LEU	2.2
1	E	137	GLU	2.2
1	F	93	HIS	2.2
1	D	161	ASP	2.2
1	F	119	ILE	2.1
1	C	169	ASN	2.1
1	E	155	LYS	2.1
1	E	21	LEU	2.1
1	F	88	TYR	2.1
1	E	32	ILE	2.1
1	E	36	ASP	2.1
1	F	55	SER	2.1
1	E	150	GLN	2.1
1	F	98	LYS	2.1
1	F	21	LEU	2.1
1	C	168	LYS	2.1
1	E	140	MET	2.1
1	D	195	ASN	2.1
1	E	30	ILE	2.1
1	F	44	ALA	2.1
1	C	189	ASP	2.0
1	C	192	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	E	67	TYR	2.0
1	E	69	SER	2.0
1	F	104	ILE	2.0
1	F	82	GLU	2.0
1	E	72	TYR	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 oligosaccharides 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.