



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:43 PM GMT

PDB ID : 1Q94
Title : Structures of HLA-A*1101 in complex with immunodominant nonamer and decamer HIV-1 epitopes clearly reveal the presence of a middle anchor residue
Authors : Li, L.; McNicholl, J.M.; Bouvier, M.
Deposited on : 2003-08-22
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 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

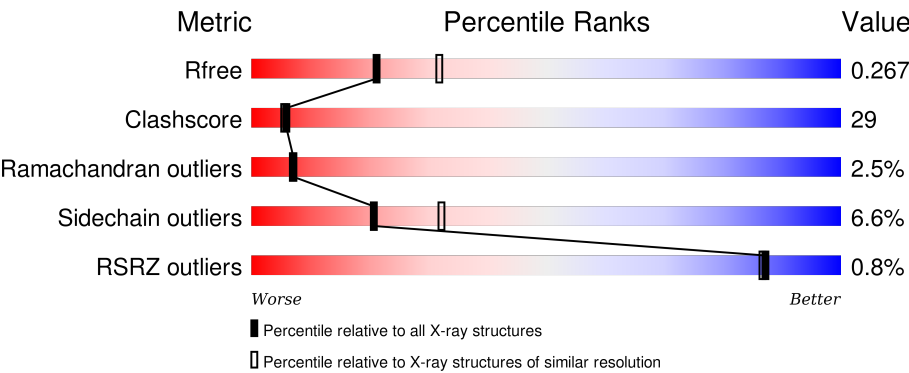
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	275	<div><div></div><div><div></div><div>56%</div><div></div><div>38%</div><div></div><div>5%</div><div></div></div></div>
1	D	275	<div><div></div><div><div></div><div>58%</div><div></div><div>37%</div><div></div><div>5%</div><div></div></div></div>
2	B	100	<div><div>2%</div><div></div><div><div></div><div>54%</div><div></div><div>42%</div><div></div><div></div></div><div></div></div>
2	E	100	<div><div>2%</div><div></div><div><div></div><div>60%</div><div></div><div>32%</div><div></div><div>8%</div><div></div></div><div></div></div>
3	C	9	<div><div></div><div><div></div><div>44%</div><div></div><div>44%</div><div></div><div>11%</div><div></div></div></div>

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Mol	Chain	Length	Quality of chain
3	F	9	 56% 44%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6456 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 HLA class I histocompatibility antigen, A-11 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2245	1394	408	434	9			
1	D	275	Total	C	N	O	S	0	0	0
			2245	1394	408	434	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	E	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	CLONING ARTIFACT	UNP P61769
E	0	MET	-	CLONING ARTIFACT	UNP P61769

- Molecule 3 is a protein called POL polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	S	0	0	0
			70	44	11	14	1			
3	F	9	Total	C	N	O	S	0	0	0
			70	44	11	14	1			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	45	Total	O	0	0
			45	45		

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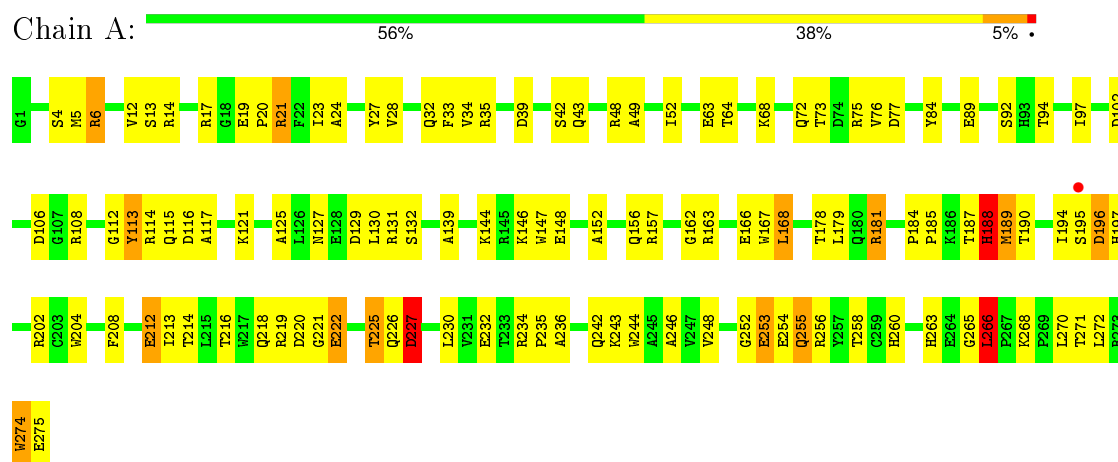
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	20	Total 20	O 20	0	0
4	C	5	Total 5	O 5	0	0
4	D	55	Total 55	O 55	0	0
4	E	21	Total 21	O 21	0	0
4	F	6	Total 6	O 6	0	0

3 Residue-property plots

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.

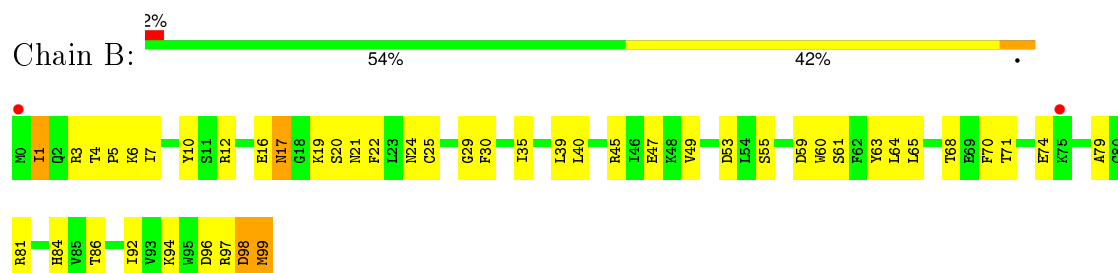
- Molecule 1: HLA class I histocompatibility antigen, A-11 alpha chain



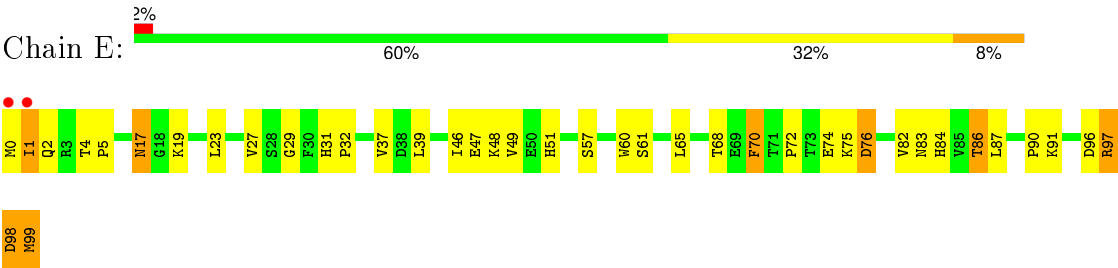
- Molecule 1: HLA class I histocompatibility antigen, A-11 alpha chain



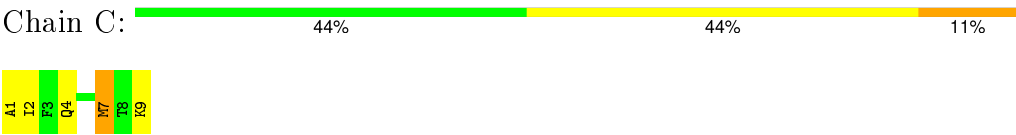
- Molecule 2: Beta-2-microglobulin



● Molecule 2: Beta-2-microglobulin



● Molecule 3: POL polypeptide



● Molecule 3: POL polypeptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	45.40 Å 46.40 Å 124.30 Å 91.50° 97.90° 114.80°	Depositor
Resolution (Å)	23.70 – 2.40 23.71 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.5 (23.70-2.40) 86.0 (23.71-2.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.64 (at 2.31 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.232 , 0.272 0.229 , 0.267	Depositor DCC
R_{free} test set	1639 reflections (4.94%)	DCC
Wilson B-factor (Å ²)	38.0	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 26.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 36125 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6456	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.17% 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, 0.375 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	0.42	0/2306	0.68	1/3129 (0.0%)
1	D	0.41	0/2306	0.68	0/3129
2	B	0.40	0/860	0.67	0/1162
2	E	0.42	0/860	0.65	0/1162
3	C	0.51	0/70	0.72	0/90
3	F	0.51	0/70	0.80	0/90
All	All	0.41	0/6472	0.68	1/8762 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	266	LEU	CA-CB-CG	6.45	130.12	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	113	TYR	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2245	0	2091	154	0
1	D	2245	0	2091	120	0
2	B	837	0	803	50	0
2	E	837	0	803	43	0
3	C	70	0	74	10	0
3	F	70	0	74	7	0
4	A	45	0	0	3	0
4	B	20	0	0	1	0
4	C	5	0	0	3	0
4	D	55	0	0	3	0
4	E	21	0	0	3	0
4	F	6	0	0	1	0
All	All	6456	0	5936	350	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 29.

All (350) 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:4:SER:HB2	1:A:6:ARG:HH12	1.26	1.01
2:E:96:ASP:HB3	2:E:99:MET:HG2	1.49	0.94
1:A:187:THR:HG22	1:A:188:HIS:N	1.84	0.90
2:B:97:ARG:HG3	2:B:98:ASP:H	1.34	0.89
1:A:35:ARG:NH1	1:A:48:ARG:HH21	1.73	0.87
2:B:20:SER:HA	2:B:71:THR:HG22	1.57	0.86
1:A:6:ARG:HH11	1:A:6:ARG:HG3	1.39	0.86
1:A:21:ARG:HH11	1:A:21:ARG:HG2	1.40	0.86
2:E:17:ASN:HD21	2:E:97:ARG:HH12	1.22	0.86
2:B:4:THR:HA	2:B:86:THR:HG21	1.57	0.84
1:A:187:THR:CG2	1:A:188:HIS:H	1.90	0.83
1:A:6:ARG:HH11	1:A:6:ARG:CG	1.91	0.83
2:B:97:ARG:HG3	2:B:98:ASP:N	1.91	0.83
1:A:162:GLY:O	1:A:166:GLU:HG2	1.78	0.82
1:A:187:THR:CG2	1:A:188:HIS:N	2.43	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:14:ARG:HB2	1:D:17:ARG:HD3	1.61	0.81
1:D:35:ARG:HE	1:D:48:ARG:HD3	1.44	0.80
1:A:187:THR:HG21	1:A:272:LEU:HG	1.62	0.80
1:A:72:GLN:HE22	1:D:75:ARG:HE	1.24	0.80
1:D:219:ARG:HB2	1:D:224:GLN:NE2	1.98	0.79
1:D:190:THR:CG2	1:D:202:ARG:HB3	2.13	0.79
1:D:80:THR:HG21	4:F:11:HOH:O	1.83	0.78
1:A:187:THR:HG21	1:A:272:LEU:CG	2.13	0.78
1:A:163:ARG:HH21	1:A:167:TRP:HE1	1.28	0.78
2:B:97:ARG:CG	2:B:98:ASP:H	1.96	0.77
1:D:19:GLU:HG3	1:D:20:PRO:HD2	1.66	0.77
1:A:202:ARG:HH11	1:A:202:ARG:HG3	1.49	0.77
1:D:171:TYR:OH	3:F:1:ALA:HB2	1.85	0.76
1:A:4:SER:HB2	1:A:6:ARG:NH1	1.98	0.76
1:A:94:THR:HG21	4:B:113:HOH:O	1.86	0.76
1:D:49:ALA:O	1:D:52:ILE:HG22	1.85	0.75
1:A:225:THR:C	1:A:226:GLN:HG3	2.05	0.75
1:A:187:THR:HG22	1:A:188:HIS:H	1.50	0.75
2:E:17:ASN:ND2	2:E:97:ARG:HH12	1.85	0.74
1:A:131:ARG:HH11	1:A:131:ARG:HG3	1.52	0.74
1:A:220:ASP:OD2	1:A:256:ARG:HG3	1.87	0.74
1:D:176:LYS:HG3	1:D:180:GLN:NE2	2.05	0.72
1:D:187:THR:O	1:D:188:HIS:HB3	1.89	0.72
1:A:252:GLY:O	1:A:253:GLU:HG3	1.90	0.72
1:A:187:THR:O	1:A:188:HIS:HB3	1.90	0.71
1:D:190:THR:HG22	1:D:202:ARG:HB3	1.73	0.71
1:A:188:HIS:CD2	1:A:188:HIS:C	2.63	0.71
1:D:99:TYR:HB3	1:D:114:ARG:HD3	1.71	0.71
1:D:14:ARG:C	1:D:17:ARG:HH11	1.95	0.70
1:A:227:ASP:N	1:A:227:ASP:OD2	2.23	0.70
1:D:77:ASP:HA	1:D:80:THR:HG23	1.74	0.70
2:B:16:GLU:HB3	2:B:19:LYS:HD2	1.73	0.70
1:A:266:LEU:HD11	1:A:270:LEU:HG	1.72	0.70
1:D:217:TRP:HA	1:D:258:THR:O	1.91	0.69
1:D:77:ASP:HA	1:D:80:THR:CG2	2.22	0.69
1:D:188:HIS:C	1:D:188:HIS:CD2	2.65	0.69
1:A:222:GLU:OE1	1:A:222:GLU:HA	1.91	0.68
1:A:13:SER:O	1:A:92:SER:HB2	1.93	0.68
1:A:114:ARG:HD3	1:A:156:GLN:NE2	2.09	0.67
2:E:17:ASN:CG	2:E:97:ARG:HH22	1.97	0.67
1:A:12:VAL:HG13	1:A:94:THR:HG22	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:SER:CB	1:A:6:ARG:HH12	2.05	0.67
1:D:76:VAL:O	1:D:80:THR:HG22	1.94	0.67
1:A:202:ARG:HH11	1:A:246:ALA:HB2	1.60	0.67
1:D:35:ARG:NE	1:D:48:ARG:HD3	2.09	0.67
1:D:28:VAL:HG11	1:D:179:LEU:HD13	1.77	0.66
1:A:6:ARG:NH1	1:A:6:ARG:CG	2.56	0.66
1:A:202:ARG:HD3	1:A:244:TRP:CD2	2.30	0.66
1:A:131:ARG:HH12	1:A:157:ARG:NH2	1.92	0.66
1:D:189:MET:HE1	1:D:274:TRP:HB2	1.77	0.66
1:A:152:ALA:HB2	3:C:7:MET:CE	2.26	0.66
1:A:187:THR:CG2	1:A:272:LEU:HG	2.26	0.66
2:E:17:ASN:ND2	2:E:97:ARG:HH22	1.93	0.65
3:C:9:LYS:NZ	4:C:119:HOH:O	2.30	0.65
1:D:215:LEU:CD2	1:D:261:VAL:HG22	2.26	0.65
2:B:16:GLU:HG2	2:B:19:LYS:HZ3	1.62	0.65
2:B:16:GLU:HG2	2:B:19:LYS:NZ	2.11	0.65
1:A:189:MET:HG3	1:A:190:THR:N	2.13	0.64
1:D:144:LYS:O	1:D:148:GLU:HG3	1.98	0.64
1:D:129:ASP:CG	1:D:131:ARG:HG2	2.18	0.64
3:C:4:GLN:HG2	4:C:17:HOH:O	1.98	0.63
1:A:28:VAL:HG11	1:A:179:LEU:HD13	1.80	0.63
1:A:187:THR:HG21	1:A:272:LEU:CD2	2.28	0.63
1:D:63:GLU:OE2	3:F:1:ALA:N	2.31	0.63
1:D:227:ASP:OD1	1:D:248:VAL:HB	1.98	0.63
1:A:184:PRO:HB3	1:A:265:GLY:O	1.99	0.63
1:A:49:ALA:O	1:A:52:ILE:HG22	1.98	0.62
1:D:129:ASP:OD1	1:D:131:ARG:HG2	1.99	0.62
1:D:138:MET:O	1:D:141:GLN:HB2	1.99	0.62
1:A:131:ARG:HH22	1:A:157:ARG:NH2	1.98	0.62
1:A:121:LYS:HG3	2:B:1:ILE:HD12	1.81	0.61
2:E:4:THR:HA	2:E:86:THR:HG21	1.81	0.61
1:A:131:ARG:HH22	1:A:157:ARG:HH22	1.48	0.60
1:D:218:GLN:NE2	1:D:260:HIS:HD2	1.99	0.60
1:A:114:ARG:HH11	1:A:156:GLN:HE22	1.49	0.60
2:E:87:LEU:HD22	2:E:91:LYS:HE3	1.82	0.60
1:A:181:ARG:C	1:A:181:ARG:HD2	2.22	0.60
2:E:96:ASP:O	2:E:99:MET:HB2	2.02	0.60
1:D:189:MET:CE	1:D:201:LEU:HG	2.31	0.60
1:A:260:HIS:CE1	1:A:271:THR:HG23	2.36	0.60
1:A:131:ARG:NH1	1:A:131:ARG:HG3	2.17	0.60
2:E:0:MET:HG2	2:E:1:ILE:N	2.17	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:219:ARG:HH11	1:D:219:ARG:HG3	1.65	0.60
1:A:131:ARG:NH2	1:A:157:ARG:NH1	2.50	0.59
1:A:21:ARG:NH1	1:A:21:ARG:HG2	2.12	0.59
1:D:94:THR:HG23	4:D:281:HOH:O	2.01	0.59
1:D:114:ARG:HH11	1:D:114:ARG:HG3	1.67	0.59
1:A:181:ARG:HH11	1:A:181:ARG:HG2	1.66	0.59
1:A:75:ARG:NH1	1:D:72:GLN:NE2	2.51	0.59
1:A:63:GLU:OE1	3:C:1:ALA:HB1	2.02	0.59
1:D:253:GLU:HB2	1:D:256:ARG:HH11	1.67	0.59
1:D:219:ARG:CZ	1:D:256:ARG:HD3	2.34	0.58
2:E:84:HIS:HB3	2:E:86:THR:HG22	1.84	0.58
1:A:187:THR:HG21	1:A:272:LEU:HD21	1.86	0.58
1:D:219:ARG:NE	1:D:256:ARG:HD3	2.17	0.58
1:A:5:MET:HB2	1:A:168:LEU:HG	1.86	0.58
1:A:212:GLU:OE1	1:A:213:ILE:N	2.37	0.58
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.38	0.57
2:B:10:TYR:O	2:B:24:ASN:HB2	2.03	0.57
1:A:188:HIS:C	1:A:188:HIS:HD2	2.05	0.57
2:B:7:ILE:HD12	2:B:7:ILE:N	2.19	0.57
1:A:152:ALA:HB2	3:C:7:MET:HE1	1.85	0.57
2:B:12:ARG:NH1	2:B:22:PHE:CD1	2.72	0.57
2:B:40:LEU:CD2	2:B:45:ARG:HG2	2.35	0.57
2:E:47:GLU:O	2:E:48:LYS:HB3	2.03	0.57
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.39	0.57
2:E:97:ARG:C	2:E:99:MET:H	2.08	0.57
1:D:189:MET:HE2	1:D:274:TRP:CD1	2.40	0.57
1:A:127:ASN:ND2	1:A:132:SER:OG	2.36	0.56
1:A:21:ARG:NH1	1:A:23:ILE:HD11	2.19	0.56
1:D:5:MET:HB2	1:D:168:LEU:HG	1.87	0.56
1:D:55:GLU:OE2	1:D:170:ARG:NH1	2.32	0.56
1:A:21:ARG:HE	1:A:39:ASP:HB2	1.70	0.56
1:A:195:SER:HA	4:A:279:HOH:O	2.05	0.56
1:A:121:LYS:CG	2:B:1:ILE:HD12	2.35	0.56
1:D:121:LYS:HE3	2:E:1:ILE:CD1	2.36	0.56
1:A:189:MET:HG2	1:A:274:TRP:HD1	1.70	0.56
1:A:14:ARG:HB3	1:A:17:ARG:HB2	1.89	0.55
1:A:187:THR:HG23	1:A:188:HIS:H	1.71	0.55
1:D:189:MET:HE1	1:D:201:LEU:HG	1.89	0.55
1:A:14:ARG:NH2	1:A:19:GLU:O	2.36	0.55
1:D:229:GLU:O	1:D:244:TRP:CZ3	2.60	0.55
1:A:208:PHE:CE2	1:A:242:GLN:HA	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:163:ARG:HG3	4:D:303:HOH:O	2.07	0.55
1:D:15:PRO:N	1:D:17:ARG:HH11	2.05	0.55
1:D:94:THR:HG21	4:E:117:HOH:O	2.06	0.55
1:D:15:PRO:HD2	1:D:17:ARG:HH12	1.72	0.54
1:A:194:ILE:HD12	1:A:248:VAL:HG13	1.88	0.54
1:D:212:GLU:HA	1:D:212:GLU:OE1	2.07	0.54
2:B:96:ASP:HB3	2:B:99:MET:HG3	1.87	0.54
2:E:17:ASN:CG	2:E:97:ARG:NH2	2.61	0.54
1:A:35:ARG:NH1	1:A:48:ARG:NH2	2.51	0.54
2:E:84:HIS:ND1	2:E:86:THR:HB	2.22	0.54
1:A:6:ARG:CZ	1:A:113:TYR:OH	2.56	0.54
1:D:225:THR:O	1:D:226:GLN:HG2	2.08	0.54
1:D:33:PHE:CD2	1:D:34:VAL:HG13	2.42	0.54
1:A:263:HIS:HB3	1:A:266:LEU:HD23	1.90	0.53
1:A:202:ARG:NH1	1:A:246:ALA:HB2	2.22	0.53
2:E:47:GLU:O	2:E:47:GLU:HG3	2.07	0.53
1:A:106:ASP:OD1	1:A:108:ARG:HD2	2.08	0.53
1:A:20:PRO:HG2	1:A:75:ARG:HG2	1.91	0.53
1:D:152:ALA:HB2	3:F:7:MET:SD	2.49	0.53
1:A:260:HIS:HE1	1:A:271:THR:HG23	1.71	0.53
1:A:35:ARG:HH12	1:A:48:ARG:HH21	1.55	0.53
1:A:220:ASP:CG	1:A:256:ARG:HG3	2.28	0.53
1:D:1:GLY:CA	1:D:105:PRO:HG3	2.39	0.53
2:E:37:VAL:HG22	2:E:82:VAL:HG22	1.89	0.53
1:D:168:LEU:O	1:D:172:LEU:HG	2.09	0.53
1:D:212:GLU:O	1:D:263:HIS:HD2	1.92	0.53
2:B:16:GLU:CG	2:B:19:LYS:NZ	2.72	0.52
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.90	0.52
1:A:4:SER:CB	1:A:6:ARG:NH1	2.68	0.52
2:B:17:ASN:ND2	2:B:74:GLU:HG2	2.24	0.52
2:E:83:ASN:ND2	2:E:90:PRO:HG3	2.23	0.52
1:D:135:ALA:HB1	1:D:140:ALA:CB	2.40	0.52
1:D:7:TYR:OH	3:F:1:ALA:HB3	2.10	0.52
1:D:55:GLU:CD	1:D:170:ARG:HH12	2.12	0.52
1:D:19:GLU:HG3	1:D:20:PRO:CD	2.37	0.52
1:D:258:THR:HB	1:D:260:HIS:CE1	2.45	0.52
2:E:23:LEU:HB2	2:E:70:PHE:CE1	2.45	0.52
2:E:96:ASP:HB3	2:E:99:MET:CG	2.31	0.51
1:A:131:ARG:NH2	1:A:157:ARG:HH12	2.08	0.51
1:D:226:GLN:O	1:D:227:ASP:HB2	2.10	0.51
1:D:188:HIS:HD2	1:D:189:MET:N	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ARG:NH1	1:D:72:GLN:HE22	2.09	0.51
1:D:126:LEU:HD11	1:D:130:LEU:HA	1.91	0.51
1:D:121:LYS:HG3	2:E:1:ILE:HD13	1.92	0.51
1:D:20:PRO:HG3	1:D:75:ARG:HG3	1.92	0.51
2:E:72:PRO:HB2	2:E:97:ARG:HH21	1.76	0.50
1:D:1:GLY:HA2	1:D:105:PRO:HG3	1.94	0.50
1:A:266:LEU:CD1	1:A:270:LEU:HG	2.40	0.50
1:D:135:ALA:HB1	1:D:140:ALA:HB3	1.92	0.50
1:D:117:ALA:HB2	2:E:60:TRP:CE2	2.47	0.50
1:A:188:HIS:NE2	1:A:204:TRP:CD1	2.79	0.50
1:A:219:ARG:N	1:A:222:GLU:O	2.31	0.50
1:D:181:ARG:HD2	1:D:181:ARG:C	2.32	0.50
2:B:25:CYS:HB2	2:B:39:LEU:HD21	1.93	0.49
1:A:163:ARG:HA	1:A:166:GLU:CG	2.42	0.49
1:A:266:LEU:HD12	1:A:268:LYS:O	2.13	0.49
1:A:13:SER:OG	1:A:92:SER:HA	2.11	0.49
1:A:114:ARG:HH11	1:A:156:GLN:NE2	2.11	0.49
1:A:274:TRP:O	1:A:275:GLU:HG3	2.11	0.49
1:D:99:TYR:CB	1:D:114:ARG:HD3	2.41	0.49
1:A:196:ASP:C	1:A:196:ASP:OD1	2.51	0.49
2:B:49:VAL:HG22	2:B:68:THR:HB	1.94	0.49
2:B:59:ASP:O	2:B:60:TRP:HB2	2.12	0.49
1:D:201:LEU:O	1:D:246:ALA:HA	2.13	0.49
1:A:20:PRO:CG	1:A:75:ARG:HG2	2.43	0.49
1:D:15:PRO:HD2	1:D:17:ARG:NH1	2.27	0.48
1:A:117:ALA:HB2	2:B:60:TRP:CZ2	2.47	0.48
1:D:9:TYR:CE1	1:D:70:GLN:HG2	2.48	0.48
2:E:31:HIS:ND1	2:E:32:PRO:HA	2.28	0.48
1:D:22:PHE:CG	1:D:71:SER:HB3	2.48	0.48
1:D:82:ARG:HD2	1:D:88:SER:O	2.13	0.48
4:A:289:HOH:O	3:C:4:GLN:HA	2.13	0.48
1:A:24:ALA:O	1:A:35:ARG:HA	2.14	0.48
1:D:19:GLU:CG	1:D:20:PRO:HD2	2.39	0.48
1:A:13:SER:HA	1:A:20:PRO:HB3	1.95	0.48
1:A:196:ASP:OD1	1:A:197:HIS:N	2.47	0.48
2:E:57:SER:HB2	4:E:103:HOH:O	2.12	0.48
2:E:17:ASN:HD22	2:E:74:GLU:N	2.12	0.48
2:B:16:GLU:CB	2:B:19:LYS:HD2	2.41	0.48
1:A:6:ARG:NH1	1:A:113:TYR:OH	2.46	0.47
1:D:14:ARG:CA	1:D:17:ARG:HH11	2.28	0.47
1:A:84:TYR:HB3	1:A:139:ALA:HB1	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:229:GLU:O	1:D:244:TRP:HZ3	1.97	0.47
2:B:17:ASN:N	2:B:17:ASN:OD1	2.36	0.47
1:A:230:LEU:HD11	1:A:243:LYS:NZ	2.29	0.47
1:A:146:LYS:HE3	1:D:84:TYR:OH	2.14	0.47
1:A:202:ARG:NE	2:B:99:MET:OXT	2.47	0.47
1:A:72:GLN:HE22	1:D:75:ARG:NE	2.04	0.47
2:B:98:ASP:O	2:B:99:MET:C	2.53	0.47
1:D:7:TYR:CE1	3:F:2:ILE:HG22	2.50	0.47
2:B:16:GLU:CG	2:B:19:LYS:HD2	2.45	0.47
1:D:184:PRO:HB3	1:D:265:GLY:O	2.14	0.47
1:A:178:THR:O	1:A:181:ARG:HG3	2.14	0.47
2:B:21:ASN:OD1	2:B:22:PHE:N	2.35	0.47
2:E:51:HIS:HA	2:E:65:LEU:O	2.15	0.47
1:A:63:GLU:OE2	3:C:1:ALA:HB3	2.15	0.47
2:B:29:GLY:HA2	2:B:61:SER:HB2	1.96	0.47
1:D:14:ARG:CB	1:D:17:ARG:HD3	2.40	0.47
1:A:167:TRP:CD1	3:C:1:ALA:N	2.82	0.46
1:A:225:THR:C	1:A:226:GLN:CG	2.81	0.46
1:D:189:MET:HE3	1:D:201:LEU:HG	1.96	0.46
1:D:218:GLN:NE2	1:D:260:HIS:CD2	2.81	0.46
2:B:40:LEU:HD11	2:B:81:ARG:HB2	1.96	0.46
1:A:181:ARG:NH1	1:A:181:ARG:HG2	2.29	0.46
1:D:14:ARG:HA	1:D:17:ARG:NH1	2.31	0.46
1:A:260:HIS:HA	1:A:270:LEU:O	2.15	0.46
1:A:144:LYS:O	1:A:148:GLU:HG3	2.16	0.46
1:A:163:ARG:HA	1:A:166:GLU:HG3	1.96	0.46
2:B:92:ILE:HG22	2:B:92:ILE:O	2.14	0.46
2:E:46:ILE:HG22	2:E:47:GLU:N	2.31	0.46
2:E:39:LEU:HD12	2:E:49:VAL:HG13	1.98	0.46
2:E:39:LEU:HD13	2:E:68:THR:HG22	1.98	0.46
1:A:115:GLN:HG2	1:A:125:ALA:HB1	1.98	0.46
1:D:35:ARG:HE	1:D:48:ARG:CD	2.20	0.45
1:A:218:GLN:CD	1:A:260:HIS:NE2	2.70	0.45
1:A:232:GLU:OE1	2:B:6:LYS:HD2	2.15	0.45
2:E:4:THR:CG2	2:E:5:PRO:HD2	2.46	0.45
1:D:188:HIS:C	1:D:188:HIS:HD2	2.15	0.45
1:D:43:GLN:NE2	1:D:43:GLN:HA	2.31	0.45
1:A:202:ARG:HD2	2:B:99:MET:OXT	2.17	0.45
1:D:218:GLN:CD	1:D:260:HIS:CD2	2.89	0.45
1:D:229:GLU:N	1:D:246:ALA:O	2.49	0.45
1:A:226:GLN:C	1:A:227:ASP:OD2	2.55	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:23:LEU:HB2	2:E:70:PHE:CD1	2.52	0.45
2:E:17:ASN:ND2	2:E:97:ARG:NH1	2.60	0.45
1:A:195:SER:C	1:A:197:HIS:N	2.70	0.45
1:D:4:SER:HB3	1:D:102:ASP:OD2	2.17	0.45
1:D:129:ASP:OD1	1:D:131:ARG:N	2.42	0.45
1:D:178:THR:O	1:D:181:ARG:HG3	2.17	0.45
1:A:202:ARG:HH11	1:A:202:ARG:CG	2.24	0.44
1:D:134:THR:HA	4:D:297:HOH:O	2.16	0.44
1:A:271:THR:O	1:A:271:THR:HG22	2.18	0.44
1:D:220:ASP:HB2	1:D:256:ARG:HG2	1.99	0.44
1:A:219:ARG:HG2	1:A:219:ARG:O	2.17	0.44
2:E:29:GLY:HA2	2:E:61:SER:CB	2.47	0.44
2:B:97:ARG:CG	2:B:98:ASP:N	2.57	0.44
1:A:234:ARG:HD2	2:B:10:TYR:CE1	2.52	0.44
2:B:3:ARG:O	2:B:30:PHE:HA	2.18	0.44
2:E:75:LYS:O	2:E:76:ASP:C	2.55	0.44
2:E:98:ASP:C	2:E:99:MET:OXT	2.56	0.44
1:A:185:PRO:HD2	1:A:266:LEU:HD22	1.99	0.44
2:B:97:ARG:O	2:B:98:ASP:C	2.55	0.44
1:A:131:ARG:NH2	1:A:157:ARG:CZ	2.81	0.44
1:A:131:ARG:NH1	1:A:157:ARG:CZ	2.81	0.44
1:D:35:ARG:CD	1:D:48:ARG:HD3	2.48	0.43
2:E:27:VAL:O	2:E:27:VAL:HG23	2.18	0.43
1:D:188:HIS:CD2	1:D:189:MET:N	2.86	0.43
1:A:219:ARG:CG	1:A:219:ARG:O	2.65	0.43
1:A:75:ARG:NH2	4:A:315:HOH:O	2.51	0.43
1:A:202:ARG:HG3	1:A:202:ARG:NH1	2.25	0.43
1:A:260:HIS:HD1	1:A:271:THR:HA	1.83	0.43
2:B:16:GLU:CG	2:B:19:LYS:HZ2	2.31	0.43
1:A:19:GLU:HB2	1:A:20:PRO:HD2	2.01	0.43
1:D:226:GLN:O	1:D:227:ASP:CB	2.66	0.43
2:E:19:LYS:HG3	4:E:114:HOH:O	2.18	0.43
2:B:96:ASP:O	2:B:99:MET:HG3	2.19	0.42
1:D:191:HIS:CD2	1:D:254:GLU:OE2	2.72	0.42
1:D:150:ALA:HB1	3:F:7:MET:CE	2.49	0.42
1:D:143:THR:HG21	3:F:9:LYS:HG3	2.01	0.42
1:A:131:ARG:HH12	1:A:157:ARG:CZ	2.32	0.42
1:A:73:THR:O	1:A:76:VAL:HG22	2.19	0.42
1:D:99:TYR:HB3	1:D:114:ARG:CD	2.46	0.42
1:A:147:TRP:CD1	3:C:7:MET:HE2	2.55	0.42
2:B:55:SER:HB3	2:B:63:TYR:CZ	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:PRO:HG2	2:B:65:LEU:HD13	2.01	0.42
2:E:17:ASN:ND2	2:E:97:ARG:NH2	2.65	0.42
1:D:185:PRO:HD3	1:D:263:HIS:ND1	2.35	0.42
1:D:22:PHE:CD2	1:D:71:SER:HB3	2.54	0.42
1:D:209:TYR:HA	1:D:210:PRO:O	2.19	0.42
1:A:236:ALA:CB	1:A:242:GLN:HE21	2.32	0.42
1:A:97:ILE:HD11	4:C:119:HOH:O	2.20	0.42
1:D:227:ASP:O	1:D:247:VAL:HA	2.20	0.42
1:A:121:LYS:HE3	2:B:1:ILE:HD12	2.02	0.42
1:D:98:MET:CE	1:D:115:GLN:HE21	2.32	0.42
2:E:29:GLY:HA2	2:E:61:SER:OG	2.20	0.42
2:E:97:ARG:C	2:E:99:MET:N	2.73	0.42
1:A:35:ARG:HD3	2:B:53:ASP:OD1	2.20	0.42
2:B:84:HIS:HB3	2:B:86:THR:HG22	2.01	0.42
1:D:195:SER:OG	1:D:197:HIS:HB2	2.20	0.42
1:D:230:LEU:HG	1:D:243:LYS:HE3	2.01	0.42
1:D:126:LEU:HD21	1:D:130:LEU:HD22	2.02	0.41
2:B:39:LEU:HA	2:B:39:LEU:HD23	1.84	0.41
1:A:129:ASP:OD1	1:A:131:ARG:N	2.52	0.41
1:A:42:SER:O	1:A:43:GLN:HB2	2.20	0.41
1:A:4:SER:HB3	1:A:102:ASP:OD2	2.21	0.41
2:E:84:HIS:C	2:E:86:THR:N	2.73	0.41
1:A:76:VAL:HG23	1:A:77:ASP:N	2.35	0.41
1:A:112:GLY:O	1:A:130:LEU:HD21	2.20	0.41
1:A:236:ALA:HB2	1:A:242:GLN:HE21	1.85	0.41
1:A:131:ARG:HH22	1:A:157:ARG:CZ	2.33	0.41
1:D:191:HIS:HB3	1:D:274:TRP:CZ2	2.56	0.41
1:D:263:HIS:H	1:D:266:LEU:HD11	1.85	0.41
2:E:29:GLY:HA2	2:E:61:SER:HB2	2.01	0.41
1:D:192:HIS:HA	1:D:193:PRO:HD3	1.84	0.41
2:B:79:ALA:HB2	2:B:94:LYS:HA	2.03	0.41
1:D:61:ASP:HB3	1:D:65:ARG:HH21	1.85	0.41
1:A:27:TYR:CE2	1:A:32:GLN:HB2	2.55	0.41
2:B:84:HIS:ND1	2:B:86:THR:HG22	2.36	0.41
1:A:64:THR:O	1:A:68:LYS:HG2	2.21	0.41
1:A:255:GLN:C	1:A:255:GLN:OE1	2.59	0.40
1:A:272:LEU:N	1:A:272:LEU:HD22	2.36	0.40
1:A:204:TRP:HZ2	2:B:99:MET:OXT	2.03	0.40
1:A:63:GLU:OE1	3:C:2:ILE:HG12	2.20	0.40
1:D:14:ARG:O	1:D:17:ARG:HD2	2.21	0.40
1:D:224:GLN:NE2	1:D:257:TYR:HE2	2.18	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:GLY:O	1:A:222:GLU:OE1	2.39	0.40
1:A:222:GLU:CA	1:A:222:GLU:OE1	2.64	0.40
2:B:17:ASN:HD22	2:B:74:GLU:HG2	1.86	0.40
1:A:214:THR:HG22	1:A:216:THR:HG23	2.03	0.40
1:A:6:ARG:NH1	1:A:6:ARG:HG2	2.36	0.40
1:D:263:HIS:O	1:D:266:LEU:HG	2.21	0.40
1:D:43:GLN:HE21	1:D:43:GLN:CA	2.34	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	273/275 (99%)	248 (91%)	21 (8%)	4 (2%)	13	17
1	D	273/275 (99%)	248 (91%)	17 (6%)	8 (3%)	6	5
2	B	98/100 (98%)	85 (87%)	10 (10%)	3 (3%)	5	4
2	E	98/100 (98%)	86 (88%)	8 (8%)	4 (4%)	3	3
3	C	7/9 (78%)	7 (100%)	0	0	100	100
3	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
All	All	756/768 (98%)	680 (90%)	57 (8%)	19 (2%)	7	7

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	187	THR
1	A	188	HIS
1	A	254	GLU
1	D	88	SER
1	D	188	HIS

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Mol	Chain	Res	Type
2	E	2	GLN
2	E	76	ASP
2	E	97	ARG
1	A	227	ASP
1	A	253	GLU
1	D	136	ALA
1	D	227	ASP
2	B	98	ASP
2	E	98	ASP
2	B	47	GLU
1	D	220	ASP
1	D	264	GLU
1	D	221	GLY
2	B	35	ILE

5.3.2 Protein sidechains ⓘ

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	232/232 (100%)	215 (93%)	17 (7%)	17	27
1	D	232/232 (100%)	216 (93%)	16 (7%)	19	30
2	B	95/95 (100%)	90 (95%)	5 (5%)	28	44
2	E	95/95 (100%)	90 (95%)	5 (5%)	28	44
3	C	8/8 (100%)	7 (88%)	1 (12%)	6	7
3	F	8/8 (100%)	8 (100%)	0	100	100
All	All	670/670 (100%)	626 (93%)	44 (7%)	21	32

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	21	ARG
1	A	89	GLU
1	A	116	ASP

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Mol	Chain	Res	Type
1	A	168	LEU
1	A	181	ARG
1	A	188	HIS
1	A	189	MET
1	A	196	ASP
1	A	212	GLU
1	A	222	GLU
1	A	225	THR
1	A	227	ASP
1	A	255	GLN
1	A	258	THR
1	A	266	LEU
1	A	274	TRP
2	B	1	ILE
2	B	17	ASN
2	B	64	LEU
2	B	70	PHE
2	B	99	MET
3	C	7	MET
1	D	6	ARG
1	D	17	ARG
1	D	35	ARG
1	D	80	THR
1	D	88	SER
1	D	111	ARG
1	D	157	ARG
1	D	168	LEU
1	D	181	ARG
1	D	188	HIS
1	D	197	HIS
1	D	212	GLU
1	D	223	ASP
1	D	258	THR
1	D	266	LEU
1	D	274	TRP
2	E	1	ILE
2	E	17	ASN
2	E	70	PHE
2	E	86	THR
2	E	99	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	72	GLN
1	A	96	GLN
1	A	127	ASN
1	A	141	GLN
1	A	156	GLN
1	A	174	ASN
1	A	192	HIS
1	A	224	GLN
1	A	242	GLN
2	B	24	ASN
1	D	32	GLN
1	D	43	GLN
1	D	72	GLN
1	D	96	GLN
1	D	115	GLN
1	D	180	GLN
1	D	188	HIS
1	D	224	GLN
2	E	17	ASN
2	E	83	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	275/275 (100%)	-0.34	1 (0%) 93 93	19, 36, 57, 62	0
1	D	275/275 (100%)	-0.37	1 (0%) 93 93	19, 37, 56, 61	0
2	B	100/100 (100%)	-0.35	2 (2%) 68 68	22, 42, 58, 63	0
2	E	100/100 (100%)	-0.39	2 (2%) 68 68	21, 37, 55, 62	0
3	C	9/9 (100%)	-0.37	0 100 100	27, 32, 36, 40	0
3	F	9/9 (100%)	-0.38	0 100 100	25, 30, 37, 38	0
All	All	768/768 (100%)	-0.36	6 (0%) 87 87	19, 37, 57, 63	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	195	SER	5.0
1	A	195	SER	4.3
2	E	0	MET	4.0
2	B	0	MET	2.4
2	B	75	LYS	2.2
2	E	1	ILE	2.1

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.