



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:52 PM GMT

PDB ID : 1MF2
Title : ANTI HIV1 PROTEASE FAB COMPLEX
Authors : Lescar, J.; Bentley, G.A.
Deposited on : 1996-12-27
Resolution : 2.60 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

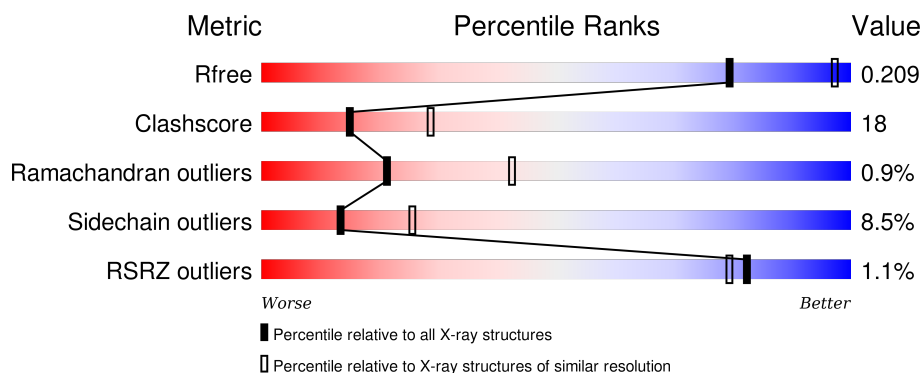
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	L	215	<div> <div>68%</div> <div>29%</div> <div>•</div> </div>
1	M	215	<div> <div>68%</div> <div>29%</div> <div>•</div> </div>
2	H	226	<div> <div>2%</div> <div>61%</div> <div>31%</div> <div>5%</div> <div>•</div> </div>
2	N	226	<div> <div>2%</div> <div>62%</div> <div>29%</div> <div>7%</div> <div>•</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6857 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 MONOCLONAL ANTIBODY F11.2.32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	215	Total	C	N	O	S	0	0	0
			1647	1024	281	334	8			
1	M	215	Total	C	N	O	S	0	0	0
			1647	1024	281	334	8			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	4	LEU	MET	CONFLICT	GB 600718
L	18	ARG	SER	CONFLICT	GB 600718
L	19	ALA	VAL	CONFLICT	GB 600718
L	27C	ASP	GLU	CONFLICT	GB 600718
L	30	LYS	THR	CONFLICT	GB 600718
L	32	PHE	LEU	CONFLICT	GB 600718
L	34	ASN	GLN	CONFLICT	GB 600718
L	36	PHE	TYR	CONFLICT	GB 600718
L	50	ALA	GLY	CONFLICT	GB 600718
L	54	GLN	VAL	CONFLICT	GB 600718
L	55	GLY	GLU	CONFLICT	GB 600718
L	74	HIS	ASN	CONFLICT	GB 600718
L	78	MET	VAL	CONFLICT	GB 600718
L	83	SER	ILE	CONFLICT	GB 600718
L	85	MET	ILE	CONFLICT	GB 600718
L	92	LYS	ARG	CONFLICT	GB 600718
L	93	GLU	LYS	CONFLICT	GB 600718
L	96	TRP	ALA	CONFLICT	GB 600718
L	100	GLY	SER	CONFLICT	GB 600718
M	4	LEU	MET	CONFLICT	GB 600718
M	18	ARG	SER	CONFLICT	GB 600718
M	19	ALA	VAL	CONFLICT	GB 600718
M	27C	ASP	GLU	CONFLICT	GB 600718
M	30	LYS	THR	CONFLICT	GB 600718
M	32	PHE	LEU	CONFLICT	GB 600718

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Chain	Residue	Modelled	Actual	Comment	Reference
M	34	ASN	GLN	CONFLICT	GB 600718
M	36	PHE	TYR	CONFLICT	GB 600718
M	50	ALA	GLY	CONFLICT	GB 600718
M	54	GLN	VAL	CONFLICT	GB 600718
M	55	GLY	GLU	CONFLICT	GB 600718
M	74	HIS	ASN	CONFLICT	GB 600718
M	78	MET	VAL	CONFLICT	GB 600718
M	83	SER	ILE	CONFLICT	GB 600718
M	85	MET	ILE	CONFLICT	GB 600718
M	92	LYS	ARG	CONFLICT	GB 600718
M	93	GLU	LYS	CONFLICT	GB 600718
M	96	TRP	ALA	CONFLICT	GB 600718
M	100	GLY	SER	CONFLICT	GB 600718

- Molecule 2 is a protein called MONOCLONAL ANTIBODY F11.2.32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	221	Total	C	N	O	S	0	0	0
			1664	1051	281	323	9			
2	N	221	Total	C	N	O	S	0	0	0
			1664	1051	281	323	9			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	5	VAL	LEU	CONFLICT	GB 600716
H	13	GLN	LYS	CONFLICT	GB 600716
H	18	ARG	LEU	CONFLICT	GB 600716
H	30	MET	SER	CONFLICT	GB 600716
H	31	ARG	ASP	CONFLICT	GB 600716
H	32	PHE	TYR	CONFLICT	GB 600716
H	74	PRO	ALA	CONFLICT	GB 600716
H	89	LEU	MET	CONFLICT	GB 600716
H	95	SER	-	INSERTION	GB 600716
H	96	GLY	-	INSERTION	GB 600716
H	97	GLY	TRP	CONFLICT	GB 600716
H	98	ILE	ASP	CONFLICT	GB 600716
H	99	GLU	THR	CONFLICT	GB 600716
H	100	ARG	THR	CONFLICT	GB 600716
H	100A	TYR	VAL	CONFLICT	GB 600716
H	100B	ASP	SER	CONFLICT	GB 600716
H	100D	THR	HIS	CONFLICT	GB 600716

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Chain	Residue	Modelled	Actual	Comment	Reference
H	182	PRO	THR	CONFLICT	GB 600716
H	183	ARG	TRP	CONFLICT	GB 600716
H	186	GLU	GLN	CONFLICT	GB 600716
N	5	VAL	LEU	CONFLICT	GB 600716
N	13	GLN	LYS	CONFLICT	GB 600716
N	18	ARG	LEU	CONFLICT	GB 600716
N	30	MET	SER	CONFLICT	GB 600716
N	31	ARG	ASP	CONFLICT	GB 600716
N	32	PHE	TYR	CONFLICT	GB 600716
N	74	PRO	ALA	CONFLICT	GB 600716
N	89	LEU	MET	CONFLICT	GB 600716
N	95	SER	-	INSERTION	GB 600716
N	96	GLY	-	INSERTION	GB 600716
N	97	GLY	TRP	CONFLICT	GB 600716
N	98	ILE	ASP	CONFLICT	GB 600716
N	99	GLU	THR	CONFLICT	GB 600716
N	100	ARG	THR	CONFLICT	GB 600716
N	100A	TYR	VAL	CONFLICT	GB 600716
N	100B	ASP	SER	CONFLICT	GB 600716
N	100D	THR	HIS	CONFLICT	GB 600716
N	182	PRO	THR	CONFLICT	GB 600716
N	183	ARG	TRP	CONFLICT	GB 600716
N	186	GLU	GLN	CONFLICT	GB 600716

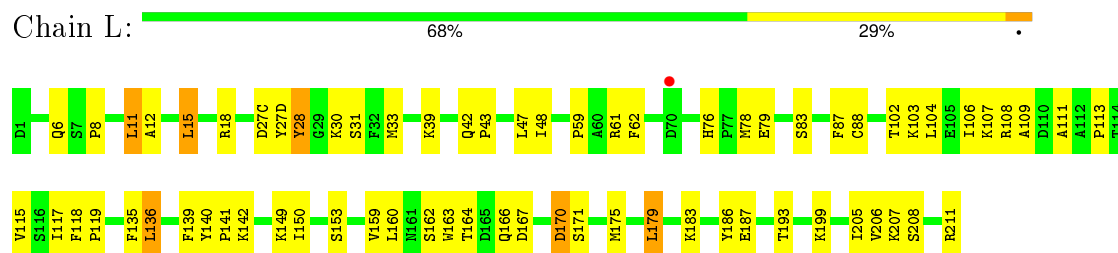
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	H	78	Total O 78 78	0	0
3	L	48	Total O 48 48	0	0
3	M	45	Total O 45 45	0	0
3	N	64	Total O 64 64	0	0

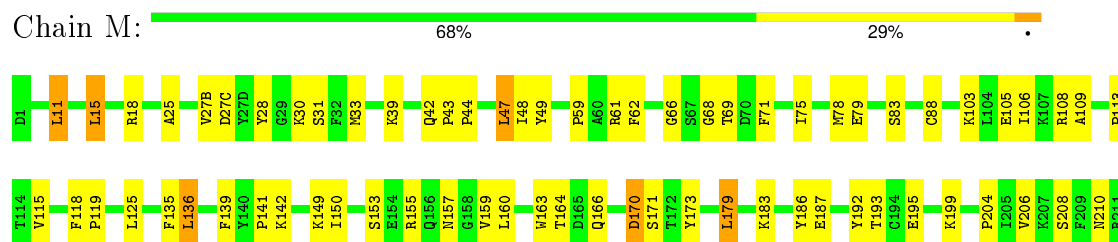
3 Residue-property plots [i](#)

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

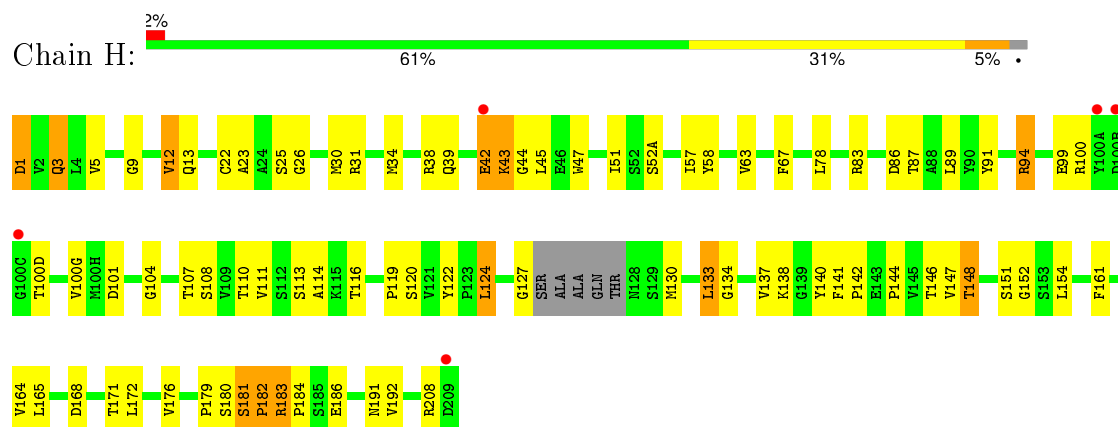
- Molecule 1: MONOCLONAL ANTIBODY F11.2.32



- Molecule 1: MONOCLONAL ANTIBODY F11.2.32

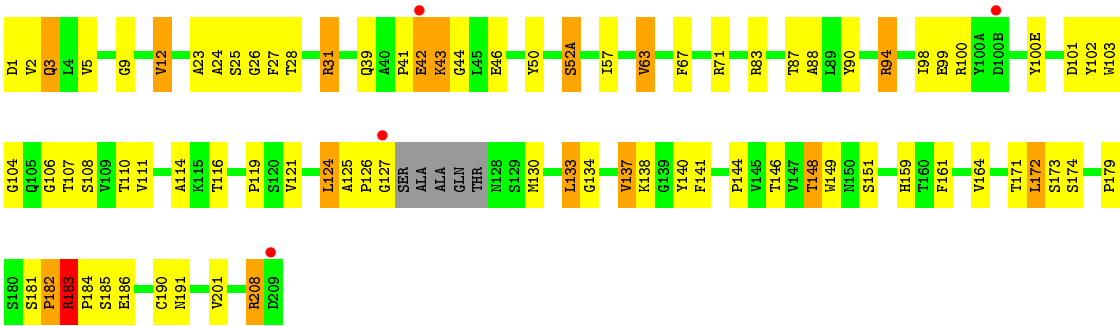


- Molecule 2: MONOCLONAL ANTIBODY F11.2.32



- Molecule 2: MONOCLONAL ANTIBODY F11.2.32





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.60 Å 94.70 Å 70.30 Å 90.00° 105.80° 90.00°	Depositor
Resolution (Å)	8.00 – 2.60 7.99 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.0 (8.00-2.60) 96.7 (7.99-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.58 Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.206 , 0.307 0.200 , 0.209	Depositor DCC
R_{free} test set	1194 reflections (4.86%)	DCC
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 65.3	EDS
Estimated twinning fraction	0.144 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 24613 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6857	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.35% 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 [i](#)

5.1 Standard geometry [i](#)

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	L	0.50	0/1686	0.76	0/2285
1	M	0.51	0/1686	0.77	0/2285
2	H	0.59	0/1706	0.83	1/2323 (0.0%)
2	N	0.55	0/1706	0.82	1/2323 (0.0%)
All	All	0.54	0/6784	0.80	2/9216 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	183	ARG	NE-CZ-NH1	6.27	123.44	120.30
2	H	42	GLU	N-CA-CB	-5.03	101.55	110.60

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	L	1647	0	1567	57	0
1	M	1647	0	1567	58	0
2	H	1664	0	1625	67	0
2	N	1664	0	1625	72	0
3	H	78	0	0	4	0
3	L	48	0	0	1	0
3	M	45	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	N	64	0	0	1	0
All	All	6857	0	6384	229	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 18.

All (229) 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)
2:N:9:GLY:H	2:N:107:THR:HG21	1.29	0.97
2:H:183:ARG:HD2	2:H:184:PRO:HA	1.46	0.96
2:N:183:ARG:HD2	2:N:184:PRO:HA	1.49	0.94
2:H:9:GLY:H	2:H:107:THR:HG21	1.32	0.93
2:H:183:ARG:HB3	2:H:183:ARG:HH11	1.32	0.92
2:N:119:PRO:HB3	2:N:140:TYR:HB3	1.53	0.89
2:N:181:SER:HB2	2:N:182:PRO:HD3	1.53	0.88
2:N:183:ARG:HH11	2:N:183:ARG:HB3	1.38	0.85
2:N:133:LEU:HD21	2:N:183:ARG:HG3	1.58	0.84
2:H:181:SER:HB2	2:H:182:PRO:HD3	1.58	0.83
2:H:133:LEU:HD21	2:H:183:ARG:HG3	1.64	0.80
2:H:119:PRO:HB3	2:H:140:TYR:HB3	1.62	0.79
2:H:179:PRO:HB2	2:H:182:PRO:HD2	1.64	0.79
2:N:146:THR:HG22	3:N:224:HOH:O	1.83	0.79
2:N:127:GLY:HA2	2:N:208:ARG:NH1	1.98	0.78
2:H:146:THR:HG21	2:N:151:SER:O	1.84	0.78
2:N:179:PRO:HB2	2:N:182:PRO:HD2	1.68	0.76
1:M:119:PRO:HD2	2:N:208:ARG:NH2	2.02	0.75
1:L:150:ILE:HD11	1:L:179:LEU:HD11	1.68	0.74
1:L:43:PRO:HB3	2:H:104:GLY:O	1.89	0.72
2:H:127:GLY:HA2	2:H:208:ARG:NH1	2.03	0.72
2:N:63:VAL:HG13	2:N:67:PHE:HB2	1.72	0.70
2:H:130:MET:HE1	2:N:23:ALA:HB3	1.73	0.69
2:H:151:SER:O	2:N:146:THR:HG21	1.92	0.69
2:H:9:GLY:N	2:H:107:THR:HG21	2.06	0.69
2:N:9:GLY:N	2:N:107:THR:HG21	2.08	0.68
1:M:160:LEU:HD22	2:N:164:VAL:HG11	1.76	0.67
1:L:39:LYS:O	1:L:42:GLN:HB2	1.93	0.67
2:N:179:PRO:O	2:N:182:PRO:HD2	1.95	0.66
2:H:9:GLY:HA3	2:H:107:THR:HG22	1.77	0.66
2:H:179:PRO:HB2	2:H:182:PRO:CD	2.25	0.66
1:M:43:PRO:HB3	2:N:104:GLY:O	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:179:PRO:O	2:H:182:PRO:HD2	1.95	0.65
2:N:9:GLY:HA3	2:N:107:THR:HG22	1.77	0.65
1:L:119:PRO:HD2	2:H:208:ARG:NH2	2.11	0.64
1:M:39:LYS:O	1:M:42:GLN:HB2	1.97	0.64
2:N:1:ASP:O	2:N:26:GLY:HA3	1.97	0.64
1:M:183:LYS:O	1:M:187:GLU:HG3	1.98	0.63
1:L:15:LEU:HD13	1:L:106:ILE:HG23	1.80	0.63
1:M:115:VAL:HG22	1:M:136:LEU:HD13	1.80	0.63
2:N:183:ARG:HD2	2:N:184:PRO:CA	2.26	0.62
2:N:179:PRO:HB2	2:N:182:PRO:CD	2.28	0.62
1:L:183:LYS:O	1:L:187:GLU:HG3	2.00	0.62
2:H:63:VAL:HG13	2:H:67:PHE:HB2	1.80	0.62
2:N:39:GLN:HA	2:N:44:GLY:O	2.00	0.61
2:H:147:VAL:HG22	2:H:192:VAL:HG22	1.82	0.61
1:M:103:LYS:HZ1	1:M:142:LYS:HE3	1.65	0.61
2:H:146:THR:HG22	3:H:230:HOH:O	2.00	0.61
2:H:138:LYS:HB3	2:H:171:THR:HG23	1.83	0.60
1:M:103:LYS:NZ	1:M:142:LYS:HE3	2.17	0.60
1:M:105:GLU:HB2	1:M:166:GLN:OE1	2.00	0.60
1:L:33:MET:SD	1:L:88:CYS:HB2	2.43	0.59
1:L:108:ARG:NH1	1:L:109:ALA:O	2.34	0.59
1:L:18:ARG:HG3	1:L:76:HIS:HB2	1.84	0.59
2:N:127:GLY:HA2	2:N:208:ARG:CZ	2.31	0.59
1:L:59:PRO:HG2	1:L:62:PHE:CD2	2.38	0.58
2:H:114:ALA:HB1	3:H:239:HOH:O	2.04	0.58
1:M:59:PRO:HG2	1:M:62:PHE:HD2	1.69	0.58
2:H:116:THR:HA	2:H:141:PHE:O	2.03	0.57
2:N:148:THR:HG22	2:N:191:ASN:OD1	2.04	0.57
1:M:141:PRO:CD	1:M:199:LYS:HD3	2.34	0.57
1:L:167:ASP:OD2	1:L:170:ASP:HB2	2.05	0.57
2:N:138:LYS:HB3	2:N:171:THR:HG23	1.84	0.57
2:H:9:GLY:HA3	2:H:107:THR:CG2	2.35	0.57
1:M:59:PRO:HG2	1:M:62:PHE:CD2	2.39	0.57
1:L:113:PRO:HG2	1:L:205:ILE:HD12	1.86	0.57
2:H:12:VAL:O	2:H:111:VAL:HA	2.04	0.57
2:H:183:ARG:HB3	2:H:183:ARG:NH1	2.13	0.56
2:H:23:ALA:HB3	2:N:130:MET:HE1	1.86	0.56
2:H:148:THR:HG22	2:H:191:ASN:HB2	1.86	0.56
2:N:2:VAL:HG21	2:N:102:TYR:CE2	2.40	0.56
1:M:108:ARG:NH1	1:M:109:ALA:O	2.38	0.56
1:L:59:PRO:HG2	1:L:62:PHE:HD2	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:183:ARG:CB	2:H:183:ARG:HH11	2.14	0.55
2:H:114:ALA:HB3	2:H:141:PHE:CE2	2.42	0.55
2:N:3:GLN:HB3	2:N:25:SER:HB2	1.88	0.55
1:M:193:THR:OG1	1:M:208:SER:HB3	2.07	0.54
1:L:149:LYS:HB2	1:L:193:THR:HB	1.88	0.54
2:N:116:THR:HA	2:N:141:PHE:O	2.08	0.54
1:L:30:LYS:HE2	2:H:100:ARG:HH21	1.71	0.54
1:L:78:MET:HE2	1:L:106:ILE:HD13	1.90	0.54
2:N:41:PRO:HD2	2:N:42:GLU:OE2	2.07	0.54
1:L:27(C):ASP:HB3	1:L:31:SER:OG	2.08	0.54
2:H:39:GLN:HA	2:H:44:GLY:O	2.09	0.53
1:L:47:LEU:O	1:L:48:ILE:HD12	2.07	0.53
1:M:155:ARG:HD2	1:M:157:ASN:O	2.07	0.53
1:M:142:LYS:HB3	1:M:173:TYR:CG	2.44	0.53
1:M:27(C):ASP:HA	1:M:31:SER:HA	1.91	0.53
1:M:195:GLU:HG3	1:M:204:PRO:HB2	1.91	0.52
1:M:108:ARG:HG3	1:M:109:ALA:O	2.10	0.52
2:N:28:THR:OG1	2:N:31:ARG:HG3	2.08	0.52
2:H:183:ARG:HD2	2:H:184:PRO:CA	2.28	0.52
2:N:181:SER:HB2	2:N:182:PRO:CD	2.33	0.52
2:H:181:SER:CB	2:H:182:PRO:HD3	2.37	0.52
1:L:115:VAL:HG22	1:L:136:LEU:HD13	1.92	0.52
2:H:100(G):VAL:HA	3:H:251:HOH:O	2.10	0.51
2:H:51:ILE:HG13	2:H:57:ILE:HG12	1.91	0.51
1:M:150:ILE:HD11	1:M:179:LEU:HD11	1.91	0.51
2:H:181:SER:HB2	2:H:182:PRO:CD	2.36	0.51
1:L:108:ARG:HG3	1:L:109:ALA:O	2.11	0.51
1:M:125:LEU:O	1:M:183:LYS:HD2	2.11	0.51
1:M:33:MET:SD	1:M:88:CYS:HB2	2.51	0.51
1:L:15:LEU:HD11	1:L:106:ILE:HD12	1.92	0.50
2:H:38:ARG:NH1	2:H:86:ASP:HA	2.26	0.50
1:L:118:PHE:HE1	1:L:135:PHE:CD2	2.29	0.50
2:H:34:MET:HB3	2:H:78:LEU:HD22	1.93	0.50
2:H:13:GLN:HE22	2:H:113:SER:HA	1.76	0.50
2:H:124:LEU:HB2	2:H:134:GLY:C	2.32	0.50
1:M:15:LEU:HD13	1:M:106:ILE:HG23	1.92	0.50
2:N:87:THR:O	2:N:88:ALA:HB2	2.12	0.50
2:H:119:PRO:CB	2:H:140:TYR:HB3	2.38	0.49
2:N:39:GLN:O	2:N:88:ALA:HB1	2.11	0.49
1:L:162:SER:O	1:L:175:MET:HA	2.12	0.49
1:L:141:PRO:CD	1:L:199:LYS:HD3	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:27(C):ASP:HB3	1:L:31:SER:HA	1.93	0.49
2:N:9:GLY:HA3	2:N:107:THR:CG2	2.43	0.49
1:L:78:MET:HE1	1:L:104:LEU:HG	1.95	0.49
2:H:152:GLY:HA2	3:H:212:HOH:O	2.13	0.49
1:L:78:MET:CE	1:L:106:ILE:HD13	2.43	0.48
2:H:47:TRP:HH2	2:H:58:TYR:HB3	1.78	0.48
2:N:127:GLY:HA2	2:N:208:ARG:HH12	1.74	0.48
1:M:78:MET:CE	1:M:106:ILE:HD13	2.43	0.48
1:M:66:GLY:HA3	1:M:71:PHE:HA	1.95	0.48
1:L:15:LEU:CD1	1:L:106:ILE:HD12	2.43	0.48
2:H:3:GLN:HB3	2:H:25:SER:HB2	1.96	0.47
2:N:43:LYS:HD3	2:N:46:GLU:HB3	1.96	0.47
1:M:142:LYS:HB3	1:M:173:TYR:CD2	2.49	0.47
2:N:121:VAL:HG21	2:N:201:VAL:CG2	2.44	0.47
2:N:90:TYR:O	2:N:106:GLY:HA2	2.14	0.47
1:L:118:PHE:CE1	1:L:135:PHE:CD2	3.03	0.47
1:L:186:TYR:CE2	1:L:211:ARG:HD2	2.49	0.47
2:H:124:LEU:HB2	2:H:134:GLY:O	2.15	0.47
2:H:94:ARG:NH1	2:H:101:ASP:OD2	2.48	0.47
1:L:6:GLN:OE1	1:L:87:PHE:HA	2.15	0.47
1:M:119:PRO:HD2	2:N:208:ARG:CZ	2.45	0.47
1:L:139:PHE:HE1	1:L:142:LYS:HA	1.79	0.47
1:L:30:LYS:HE2	2:H:100:ARG:NH2	2.30	0.47
1:L:118:PHE:HE1	1:L:135:PHE:HD2	1.62	0.47
1:L:79:GLU:OE2	1:M:61:ARG:HD3	2.15	0.47
2:H:179:PRO:C	2:H:182:PRO:HD2	2.35	0.46
2:N:87:THR:HG23	2:N:110:THR:HA	1.95	0.46
2:N:161:PHE:HD2	2:N:173:SER:O	1.97	0.46
1:L:160:LEU:HD22	2:H:164:VAL:HG11	1.97	0.46
1:M:25:ALA:O	1:M:69:THR:HG23	2.15	0.46
2:H:87:THR:HG23	2:H:110:THR:HA	1.98	0.46
1:L:193:THR:CG2	1:L:206:VAL:HG13	2.46	0.46
2:N:12:VAL:O	2:N:111:VAL:HA	2.16	0.46
1:M:136:LEU:N	1:M:136:LEU:HD22	2.31	0.46
1:M:118:PHE:HE1	1:M:135:PHE:CD2	2.34	0.46
1:M:27(C):ASP:CB	1:M:31:SER:HA	2.46	0.45
2:N:159:HIS:O	2:N:174:SER:HA	2.15	0.45
3:L:259:HOH:O	2:N:1:ASP:HB2	2.16	0.45
1:M:186:TYR:HA	1:M:192:TYR:OH	2.17	0.45
2:H:127:GLY:HA2	2:H:208:ARG:CZ	2.45	0.45
1:L:18:ARG:HB2	1:L:76:HIS:CD2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:115:VAL:HG22	1:L:136:LEU:CD1	2.46	0.45
2:N:172:LEU:C	2:N:172:LEU:HD23	2.37	0.45
1:M:47:LEU:O	1:M:48:ILE:HD12	2.16	0.45
1:M:108:ARG:HG3	1:M:108:ARG:HH11	1.81	0.45
1:L:61:ARG:HD3	1:M:79:GLU:OE2	2.17	0.45
1:M:170:ASP:O	1:M:171:SER:HB2	2.16	0.45
1:M:49:TYR:CZ	2:N:98:ILE:HD13	2.51	0.45
2:H:43:LYS:HE3	2:H:43:LYS:HB2	1.74	0.45
1:L:11:LEU:HD13	1:L:104:LEU:HD13	1.98	0.45
2:N:67:PHE:CD1	2:N:67:PHE:N	2.85	0.44
2:N:119:PRO:CB	2:N:140:TYR:HB3	2.37	0.44
1:M:149:LYS:HB2	1:M:193:THR:HB	1.99	0.44
1:L:164:THR:HG23	2:H:161:PHE:CD1	2.53	0.44
1:M:78:MET:HE3	1:M:106:ILE:HD13	1.99	0.44
1:M:30:LYS:HD3	2:N:100(E):TYR:HD1	1.82	0.44
1:L:27(D):TYR:O	1:L:28:TYR:CB	2.65	0.44
1:L:139:PHE:CE1	1:L:142:LYS:HA	2.53	0.44
2:H:138:LYS:CB	2:H:171:THR:HG23	2.48	0.43
1:L:108:ARG:NH1	1:L:111:ALA:HB2	2.33	0.43
1:M:11:LEU:O	1:M:105:GLU:HG2	2.19	0.43
1:M:193:THR:CG2	1:M:206:VAL:HG13	2.47	0.43
1:M:118:PHE:CE1	1:M:135:PHE:CD2	3.05	0.43
2:N:94:ARG:NH1	2:N:101:ASP:OD2	2.50	0.43
2:H:89:LEU:HD23	2:H:91:TYR:CZ	2.53	0.43
2:N:125:ALA:HA	2:N:126:PRO:HD3	1.87	0.43
1:L:170:ASP:O	1:L:171:SER:HB2	2.17	0.43
1:M:113:PRO:HB3	1:M:139:PHE:HB3	2.00	0.43
2:H:1:ASP:O	2:H:26:GLY:HA3	2.19	0.43
1:M:44:PRO:HD2	2:N:103:TRP:CE3	2.54	0.43
2:N:124:LEU:HB2	2:N:134:GLY:C	2.40	0.42
1:M:139:PHE:HE1	1:M:142:LYS:HA	1.85	0.42
2:H:5:VAL:HB	2:N:130:MET:HE1	2.01	0.42
2:H:67:PHE:N	2:H:67:PHE:CD1	2.87	0.42
1:L:166:GLN:HG2	1:L:171:SER:HA	2.02	0.42
1:M:27(B):VAL:O	1:M:27(B):VAL:HG12	2.18	0.42
1:L:113:PRO:HB3	1:L:139:PHE:HB3	2.01	0.42
2:N:149:TRP:CH2	2:N:190:CYS:HB3	2.54	0.42
1:L:149:LYS:HA	1:L:153:SER:O	2.20	0.42
1:L:193:THR:HG23	1:L:206:VAL:HG13	2.02	0.42
2:N:137:VAL:O	2:N:137:VAL:HG22	2.20	0.42
1:M:136:LEU:CD2	1:M:136:LEU:N	2.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:120:SER:HB3	2:H:122:TYR:CZ	2.54	0.41
1:L:107:LYS:HA	1:L:140:TYR:OH	2.19	0.41
2:N:179:PRO:C	2:N:182:PRO:HD2	2.40	0.41
2:N:148:THR:HG22	2:N:191:ASN:HB2	2.03	0.41
1:M:27(C):ASP:HB3	1:M:31:SER:OG	2.21	0.41
2:N:43:LYS:HE3	2:N:43:LYS:HB2	1.71	0.41
2:H:39:GLN:HB2	2:H:45:LEU:HD23	2.03	0.41
1:M:149:LYS:HA	1:M:153:SER:O	2.20	0.41
1:M:18:ARG:HA	1:M:75:ILE:O	2.21	0.41
2:H:130:MET:CE	2:N:3:GLN:HE22	2.33	0.41
2:H:130:MET:CE	2:N:5:VAL:HB	2.50	0.41
2:N:114:ALA:HB3	2:N:141:PHE:CE2	2.56	0.41
1:M:103:LYS:HD2	3:M:217:HOH:O	2.21	0.41
1:L:27(C):ASP:CB	1:L:31:SER:HA	2.51	0.41
1:L:47:LEU:C	1:L:48:ILE:HD12	2.41	0.41
2:N:121:VAL:HG21	2:N:201:VAL:HG21	2.02	0.41
2:N:50:TYR:O	2:N:57:ILE:HA	2.21	0.41
1:L:8:PRO:O	1:L:102:THR:HG23	2.20	0.41
1:L:117:ILE:HG22	1:L:207:LYS:HG3	2.02	0.41
2:H:127:GLY:HA2	2:H:208:ARG:HH12	1.80	0.41
1:L:103:LYS:NZ	1:L:142:LYS:HE3	2.35	0.41
2:H:148:THR:HG22	2:H:191:ASN:OD1	2.21	0.41
1:M:30:LYS:HE2	2:N:100:ARG:HH21	1.86	0.41
2:N:24:ALA:HB1	2:N:27:PHE:CE1	2.56	0.41
2:N:121:VAL:CG2	2:N:201:VAL:HG21	2.50	0.41
1:M:164:THR:HG23	2:N:161:PHE:CD1	2.56	0.40
1:M:141:PRO:HD3	1:M:199:LYS:HD3	2.03	0.40
1:L:103:LYS:HE3	1:L:103:LYS:HB3	1.84	0.40
1:M:78:MET:HE3	1:M:106:ILE:CD1	2.51	0.40
1:M:30:LYS:HD3	2:N:100(E):TYR:CD1	2.57	0.40
2:N:52(A):SER:HA	2:N:71:ARG:NH1	2.37	0.40
2:H:165:LEU:HD11	2:H:168:ASP:HA	2.04	0.40
2:H:154:LEU:HD13	2:H:176:VAL:HG21	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	L	213/215 (99%)	204 (96%)	7 (3%)	2 (1%)	21	42
1	M	213/215 (99%)	202 (95%)	9 (4%)	2 (1%)	21	42
2	H	217/226 (96%)	203 (94%)	12 (6%)	2 (1%)	21	42
2	N	217/226 (96%)	203 (94%)	12 (6%)	2 (1%)	21	42
All	All	860/882 (98%)	812 (94%)	40 (5%)	8 (1%)	21	42

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	12	ALA
1	L	28	TYR
2	H	43	LYS
1	M	28	TYR
2	N	43	LYS
2	H	30	MET
2	N	208	ARG
1	M	68	GLY

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	L	185/187 (99%)	176 (95%)	9 (5%)	31	57
1	M	185/187 (99%)	175 (95%)	10 (5%)	27	52

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	186/192 (97%)	162 (87%)	24 (13%)	5	9
2	N	186/192 (97%)	166 (89%)	20 (11%)	8	15
All	All	742/758 (98%)	679 (92%)	63 (8%)	13	25

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	11	LEU
1	L	15	LEU
1	L	83	SER
1	L	136	LEU
1	L	159	VAL
1	L	163	TRP
1	L	170	ASP
1	L	179	LEU
1	L	208	SER
2	H	1	ASP
2	H	3	GLN
2	H	12	VAL
2	H	22	CYS
2	H	31	ARG
2	H	42	GLU
2	H	52(A)	SER
2	H	83	ARG
2	H	94	ARG
2	H	99	GLU
2	H	100(D)	THR
2	H	108	SER
2	H	124	LEU
2	H	133	LEU
2	H	137	VAL
2	H	142	PRO
2	H	144	PRO
2	H	148	THR
2	H	172	LEU
2	H	180	SER
2	H	181	SER
2	H	182	PRO
2	H	183	ARG
2	H	186	GLU
1	M	11	LEU

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Mol	Chain	Res	Type
1	M	15	LEU
1	M	47	LEU
1	M	83	SER
1	M	136	LEU
1	M	159	VAL
1	M	163	TRP
1	M	170	ASP
1	M	179	LEU
1	M	210	ASN
2	N	3	GLN
2	N	12	VAL
2	N	31	ARG
2	N	42	GLU
2	N	52(A)	SER
2	N	63	VAL
2	N	83	ARG
2	N	94	ARG
2	N	99	GLU
2	N	108	SER
2	N	124	LEU
2	N	133	LEU
2	N	137	VAL
2	N	144	PRO
2	N	148	THR
2	N	172	LEU
2	N	182	PRO
2	N	183	ARG
2	N	185	SER
2	N	186	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	53	ASN
1	L	76	HIS
1	L	137	ASN
1	L	189	HIS
2	H	3	GLN
2	H	13	GLN
1	M	53	ASN
1	M	54	GLN
1	M	76	HIS

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Mol	Chain	Res	Type
1	M	124	GLN
1	M	137	ASN
2	N	3	GLN
2	N	159	HIS

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 [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 [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	L	215/215 (100%)	-0.42	1 (0%) 91 90	3, 16, 36, 50	0
1	M	215/215 (100%)	-0.46	0 100 100	3, 14, 37, 53	0
2	H	221/226 (97%)	-0.50	5 (2%) 64 57	3, 12, 34, 70	0
2	N	221/226 (97%)	-0.47	4 (1%) 71 66	3, 14, 35, 67	0
All	All	872/882 (98%)	-0.46	10 (1%) 82 79	3, 14, 37, 70	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	N	209	ASP	5.1
2	H	209	ASP	4.8
2	H	100(A)	TYR	3.5
2	H	42	GLU	3.4
2	H	100(B)	ASP	3.0
2	N	127	GLY	2.5
2	H	100(C)	GLY	2.4
2	N	42	GLU	2.3
2	N	100(B)	ASP	2.1
1	L	70	ASP	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.