Full wwPDB X-ray Structure Validation Report ⓘ

Feb 3, 2024 – 08:17 PM EST

PDB ID	:	1MN8
Title	:	Structure of Moloney Murine Leukaemia Virus Matrix Protein
Authors	:	Riffel, N.; Harlos, K.; Iourin, O.; Rao, Z.; Kingsman, A.; Stuart, D.; Fry, E.
Deposited on	:	2002-09-05
Resolution	:	1.00 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:

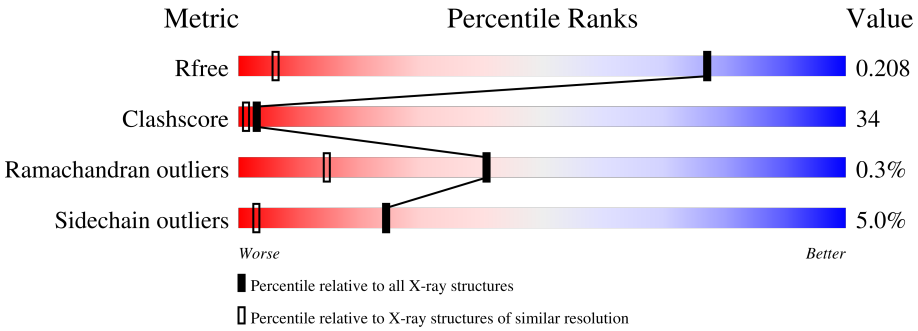
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.00 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	1050 (1.06-0.94)
Clashscore	141614	1117 (1.06-0.94)
Ramachandran outliers	138981	1043 (1.06-0.94)
Sidechain outliers	138945	1045 (1.06-0.94)

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%

Mol	Chain	Length	Quality of chain
1	A	100	67% 20% • • 5%
1	B	100	74% 15% • 7%
1	C	100	72% 14% 8% 6%
1	D	100	67% 20% 8% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3732 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 Core protein p15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	95	Total	C	N	O	S	0	8	0
			797	517	140	139	1			
1	B	93	Total	C	N	O	S	0	6	0
			778	511	134	132	1			
1	C	94	Total	C	N	O	S	0	5	0
			782	509	139	133	1			
1	D	97	Total	C	N	O	S	0	7	0
			806	523	142	140	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ALA	-	cloning artifact	UNP P03332
B	0	ALA	-	cloning artifact	UNP P03332
C	0	ALA	-	cloning artifact	UNP P03332
D	0	ALA	-	cloning artifact	UNP P03332

- Molecule 2 is water.

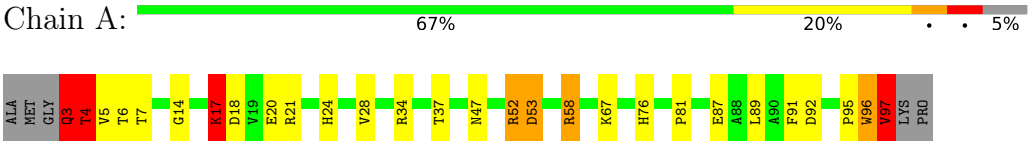
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	163	Total	O	0	0
			163	163		
2	B	136	Total	O	0	0
			136	136		
2	C	116	Total	O	0	0
			116	116		
2	D	154	Total	O	0	0
			154	154		

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. 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. 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.

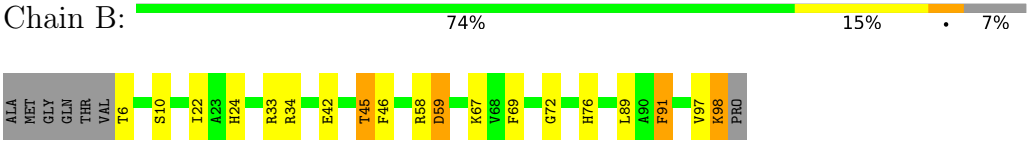
• Molecule 1: Core protein p15

Chain A:



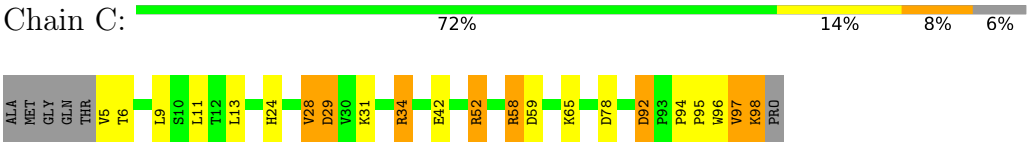
• Molecule 1: Core protein p15

Chain B:



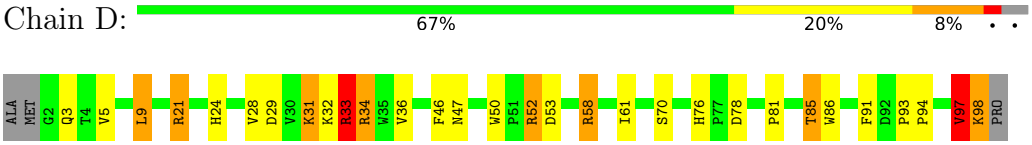
• Molecule 1: Core protein p15

Chain C:



• Molecule 1: Core protein p15

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	33.80Å 49.50Å 50.80Å 71.90° 81.90° 80.00°	Depositor
Resolution (Å)	20.00 – 1.00 46.60 – 1.00	Depositor EDS
% Data completeness (in resolution range)	86.0 (20.00-1.00) 79.2 (46.60-1.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	-0.34 (at 0.99Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.133 , 0.169 0.209 , 0.208	Depositor DCC
R_{free} test set	7088 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	9.6	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 72.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.012 for -h,-l,-k	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	3732	wwPDB-VP
Average B, all atoms (Å^2)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.01% 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	2.79	7/875 (0.8%)	2.63	34/1200 (2.8%)
1	B	0.76	0/841	1.32	12/1155 (1.0%)
1	C	0.79	0/842	1.54	12/1155 (1.0%)
1	D	0.77	0/875	1.39	15/1202 (1.2%)
All	All	1.56	7/3433 (0.2%)	1.81	73/4712 (1.5%)

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	1	2
1	B	1	0
1	C	0	1
1	D	1	0
All	All	3	3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	3	GLN	N-CA	54.08	2.54	1.46
1	A	3	GLN	CD-OE1	34.46	1.99	1.24
1	A	3	GLN	CB-CG	27.64	2.27	1.52
1	A	3	GLN	CA-CB	27.00	2.13	1.53
1	A	3	GLN	CG-CD	-17.44	1.10	1.51
1	A	3	GLN	CA-C	13.63	1.88	1.52
1	A	3	GLN	CD-NE2	11.95	1.62	1.32

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	3	GLN	N-CA-CB	-53.50	14.30	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	3	GLN	CB-CA-C	-27.57	55.27	110.40
1	A	3	GLN	N-CA-C	-25.86	41.17	111.00
1	A	3	GLN	CA-CB-CG	-19.98	69.45	113.40
1	C	58[A]	ARG	CD-NE-CZ	19.27	150.58	123.60
1	C	58[B]	ARG	CD-NE-CZ	19.27	150.58	123.60
1	A	3	GLN	OE1-CD-NE2	-17.37	81.95	121.90
1	D	34	ARG	CD-NE-CZ	15.43	145.20	123.60
1	A	3	GLN	CG-CD-NE2	13.42	148.91	116.70
1	A	52[A]	ARG	NE-CZ-NH1	-11.03	114.79	120.30
1	A	52[B]	ARG	NE-CZ-NH1	-11.03	114.79	120.30
1	A	34	ARG	NE-CZ-NH1	-10.45	115.07	120.30
1	A	21	ARG	NE-CZ-NH2	-10.38	115.11	120.30
1	D	52	ARG	NE-CZ-NH2	10.01	125.31	120.30
1	C	28	VAL	CG1-CB-CG2	9.43	125.98	110.90
1	A	3	GLN	CA-C-N	-9.08	97.23	117.20
1	C	42	GLU	OE1-CD-OE2	8.95	134.03	123.30
1	A	96	TRP	C-N-CA	7.97	141.64	121.70
1	A	18	ASP	CB-CG-OD1	7.96	125.47	118.30
1	C	29	ASP	CB-CG-OD2	7.96	125.46	118.30
1	A	52[A]	ARG	CA-CB-CG	7.72	130.38	113.40
1	A	52[B]	ARG	CA-CB-CG	7.72	130.38	113.40
1	A	53[A]	ASP	CB-CG-OD2	7.64	125.18	118.30
1	A	53[B]	ASP	CB-CG-OD2	7.64	125.18	118.30
1	D	98	LYS	CB-CA-C	7.40	125.20	110.40
1	A	52[A]	ARG	NH1-CZ-NH2	7.34	127.48	119.40
1	A	52[B]	ARG	NH1-CZ-NH2	7.34	127.48	119.40
1	B	45[A]	THR	OG1-CB-CG2	7.33	126.86	110.00
1	B	45[B]	THR	OG1-CB-CG2	7.33	126.86	110.00
1	B	69	PHE	CB-CG-CD1	7.26	125.88	120.80
1	D	9	LEU	CB-CG-CD1	7.18	123.20	111.00
1	B	91[A]	PHE	CB-CG-CD1	-7.09	115.84	120.80
1	B	91[B]	PHE	CB-CG-CD1	-7.09	115.84	120.80
1	B	34	ARG	NE-CZ-NH2	7.07	123.83	120.30
1	A	18	ASP	CB-CG-OD2	-7.03	111.97	118.30
1	B	33	ARG	NE-CZ-NH1	-6.92	116.84	120.30
1	B	34	ARG	NE-CZ-NH1	-6.84	116.88	120.30
1	C	92	ASP	CB-CG-OD2	6.84	124.46	118.30
1	C	97	VAL	C-N-CA	6.83	138.77	121.70
1	D	52	ARG	NE-CZ-NH1	-6.82	116.89	120.30
1	B	59	ASP	CB-CG-OD1	6.68	124.31	118.30
1	D	97	VAL	C-N-CA	-6.65	105.07	121.70
1	A	58[A]	ARG	NE-CZ-NH2	6.60	123.60	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	58[B]	ARG	NE-CZ-NH2	6.60	123.60	120.30
1	A	95	PRO	O-C-N	-6.59	112.15	122.70
1	A	97	VAL	N-CA-C	6.18	127.70	111.00
1	D	46	PHE	CB-CG-CD1	-6.14	116.50	120.80
1	A	17[A]	LYS	CG-CD-CE	6.13	130.30	111.90
1	A	17[B]	LYS	CG-CD-CE	6.13	130.30	111.90
1	A	3	GLN	O-C-N	6.13	132.51	122.70
1	D	46	PHE	CB-CG-CD2	6.03	125.02	120.80
1	C	58[A]	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	C	58[B]	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	D	52	ARG	CD-NE-CZ	5.91	131.87	123.60
1	B	97	VAL	CA-CB-CG2	5.91	119.76	110.90
1	A	91	PHE	CB-CG-CD2	5.75	124.83	120.80
1	C	28	VAL	N-CA-CB	-5.58	99.21	111.50
1	B	46	PHE	CB-CG-CD1	-5.52	116.94	120.80
1	A	53[A]	ASP	CB-CG-OD1	-5.40	113.44	118.30
1	A	53[B]	ASP	CB-CG-OD1	-5.40	113.44	118.30
1	A	52[A]	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	A	52[B]	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	C	34	ARG	NE-CZ-NH2	5.33	122.97	120.30
1	D	33	ARG	NE-CZ-NH1	-5.32	117.64	120.30
1	D	78	ASP	CB-CG-OD2	-5.25	113.57	118.30
1	A	34	ARG	NH1-CZ-NH2	5.24	125.16	119.40
1	A	91	PHE	CB-CG-CD1	-5.23	117.14	120.80
1	B	46	PHE	CB-CG-CD2	5.23	124.46	120.80
1	D	21[A]	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	D	21[B]	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	D	85	THR	CA-CB-OG1	5.06	119.63	109.00
1	D	28	VAL	CG1-CB-CG2	5.02	118.93	110.90
1	C	52	ARG	NE-CZ-NH2	-5.01	117.80	120.30

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	3	GLN	CA
1	B	45[A]	THR	CB
1	D	85	THR	CB

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	3	GLN	Mainchain

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Mol	Chain	Res	Type	Group
1	A	96	TRP	Peptide
1	C	97	VAL	Peptide

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	797	0	769	100	0
1	B	778	0	759	27	0
1	C	782	0	762	34	1
1	D	806	0	786	77	0
2	A	163	0	0	12	5
2	B	136	0	0	10	2
2	C	116	0	0	13	2
2	D	154	0	0	17	2
All	All	3732	0	3076	217	7

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 34.

All (217) 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:3:GLN:N	1:A:3:GLN:C	1.67	1.47
1:A:17[B]:LYS:HZ1	1:A:20:GLU:CD	1.21	1.43
1:A:3:GLN:C	1:A:3:GLN:CB	1.87	1.43
1:A:3:GLN:C	1:A:3:GLN:CA	1.88	1.42
1:C:95:PRO:O	1:C:98:LYS:CB	1.69	1.41
1:A:3:GLN:N	1:A:4:THR:N	1.75	1.35
1:A:17[B]:LYS:NZ	1:A:20:GLU:CD	1.83	1.30
1:A:3:GLN:CB	1:D:52:ARG:HE	1.43	1.30
1:D:3:GLN:NE2	1:D:5[A]:VAL:HG12	1.49	1.27
1:B:6:THR:HG22	1:B:10[A]:SER:OG	1.18	1.26
1:A:3:GLN:CB	1:A:3:GLN:CA	2.13	1.26
1:A:3:GLN:HB3	1:D:52:ARG:NE	1.50	1.25
1:D:52:ARG:HB3	2:D:2322:HOH:O	1.07	1.21

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:THR:CG2	1:B:10[A]:SER:OG	1.89	1.20
1:A:3:GLN:CB	1:A:4:THR:N	2.04	1.19
1:A:3:GLN:N	1:A:3:GLN:O	1.74	1.17
1:C:58[A]:ARG:HD3	2:C:2120:HOH:O	1.30	1.16
1:C:95:PRO:O	1:C:98:LYS:HB3	1.35	1.16
1:B:59:ASP:CG	2:B:2222:HOH:O	1.85	1.15
1:C:31:LYS:HD2	1:C:34:ARG:NH1	1.60	1.14
1:D:97:VAL:O	1:D:98:LYS:O	1.66	1.14
1:A:3:GLN:CB	1:A:3:GLN:CG	2.27	1.12
1:C:95:PRO:O	1:C:98:LYS:HB2	1.36	1.12
1:A:17[B]:LYS:NZ	1:A:20:GLU:OE1	1.81	1.11
1:A:4:THR:HG23	1:A:6[B]:THR:HG22	1.30	1.10
1:A:3:GLN:HB3	1:D:52:ARG:HE	0.99	1.10
1:A:4:THR:HG23	1:A:6[A]:THR:HG23	1.31	1.10
1:A:17[B]:LYS:HZ1	1:A:20:GLU:CG	1.65	1.09
1:A:3:GLN:CB	1:A:3:GLN:O	2.02	1.08
1:A:3:GLN:N	1:A:3:GLN:HB3	1.40	1.07
1:C:31:LYS:HD2	1:C:34:ARG:HH11	1.06	1.05
1:A:3:GLN:HB2	1:A:4:THR:N	1.67	1.04
1:A:4:THR:HG23	1:A:6[B]:THR:CG2	1.88	1.04
1:C:94:PRO:O	1:C:98:LYS:HA	1.56	1.04
1:A:3:GLN:N	1:D:52:ARG:HE	1.57	1.01
1:B:59:ASP:OD2	2:B:2222:HOH:O	1.72	1.01
1:A:3:GLN:OE1	1:A:3:GLN:CD	1.99	1.00
1:C:24:HIS:HD2	1:D:24:HIS:CD2	1.79	1.00
1:C:24:HIS:CD2	1:D:24:HIS:HD2	1.79	0.99
1:A:4:THR:CG2	1:A:6[B]:THR:CG2	2.40	0.99
1:B:72:GLY:HA2	2:B:2258:HOH:O	1.59	0.99
1:B:22:ILE:HD13	1:B:89[B]:LEU:CD1	1.95	0.97
1:C:98:LYS:HD3	1:C:98:LYS:C	1.84	0.96
1:C:5:VAL:HG13	1:C:5:VAL:O	1.67	0.94
1:D:3:GLN:HE21	1:D:5[A]:VAL:HG12	1.24	0.94
1:D:98:LYS:HB2	2:D:2405:HOH:O	1.66	0.94
1:D:33:ARG:O	2:D:2077:HOH:O	1.83	0.94
1:A:4:THR:CG2	1:A:6[A]:THR:HG23	1.98	0.94
1:A:87:GLU:OE2	2:A:2316:HOH:O	1.85	0.93
1:A:17[A]:LYS:HZ2	1:A:20:GLU:HB2	1.34	0.92
1:D:21[B]:ARG:HD3	2:D:1962:HOH:O	1.67	0.92
1:D:3:GLN:CG	1:D:5[A]:VAL:HG12	2.01	0.90
1:A:3:GLN:HB3	1:D:52:ARG:CZ	2.01	0.89
1:D:3:GLN:CD	1:D:5[A]:VAL:HG12	1.92	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:THR:HG22	1:B:10[A]:SER:HG	1.24	0.89
1:D:3:GLN:NE2	1:D:5[A]:VAL:CG1	2.35	0.89
1:A:3:GLN:C	1:A:3:GLN:HB2	1.71	0.88
1:A:3:GLN:CD	2:A:2342:HOH:O	2.10	0.88
1:A:3:GLN:CA	1:A:3:GLN:CG	2.51	0.88
1:A:4:THR:CG2	1:A:6[B]:THR:HG23	2.02	0.87
1:B:22:ILE:CD1	1:B:89[B]:LEU:CD1	2.51	0.87
1:C:5:VAL:HG13	2:C:2027:HOH:O	1.73	0.86
1:A:3:GLN:CB	1:D:52:ARG:NE	2.16	0.86
1:A:4:THR:CG2	1:A:6[B]:THR:HG22	2.04	0.86
1:A:3:GLN:N	1:A:3:GLN:CB	0.71	0.86
1:C:31:LYS:CD	1:C:34:ARG:NH1	2.38	0.85
1:A:3:GLN:O	1:A:3:GLN:CG	2.24	0.85
1:B:22:ILE:HD13	1:B:89[B]:LEU:HD13	1.57	0.84
1:A:3:GLN:CA	1:A:3:GLN:CD	2.47	0.82
1:A:6[A]:THR:HG22	2:D:2108:HOH:O	1.80	0.82
1:D:32:LYS:HB3	1:D:33:ARG:HH21	1.45	0.82
1:C:31:LYS:CD	1:C:34:ARG:HH11	1.93	0.81
1:A:17[A]:LYS:NZ	1:A:20:GLU:HB2	1.94	0.81
1:A:3:GLN:CA	1:A:4:THR:N	2.43	0.80
1:D:53[B]:ASP:OD1	2:D:2254:HOH:O	2.00	0.80
1:C:58[A]:ARG:HD2	2:C:2120:HOH:O	1.42	0.79
1:A:37:THR:HG21	2:A:2396:HOH:O	1.84	0.78
1:A:17[B]:LYS:HZ2	1:A:20:GLU:CD	1.87	0.77
1:D:70[B]:SER:OG	2:D:2469:HOH:O	2.01	0.77
1:B:22:ILE:CD1	1:B:89[B]:LEU:HD13	2.14	0.76
1:A:3:GLN:CG	2:A:2342:HOH:O	2.33	0.76
1:A:47:ASN:ND2	2:A:2293:HOH:O	2.18	0.75
1:A:3:GLN:CG	1:D:52:ARG:HD2	2.17	0.75
1:C:95:PRO:C	1:C:98:LYS:CB	2.55	0.74
1:C:65[B]:LYS:CE	2:C:1907:HOH:O	2.28	0.74
1:A:3:GLN:N	1:A:4:THR:CA	2.51	0.73
1:B:22:ILE:HD13	1:B:89[B]:LEU:HD12	1.71	0.73
1:A:4:THR:HG21	1:A:6[B]:THR:CG2	2.20	0.72
1:A:5:VAL:HB	1:D:3:GLN:HG2	1.72	0.71
1:B:67:LYS:NZ	1:B:76:HIS:HE1	1.87	0.71
1:A:3:GLN:N	1:A:3:GLN:CA	2.54	0.71
1:A:17[A]:LYS:NZ	1:A:20:GLU:CB	2.54	0.70
1:D:81:PRO:O	1:D:85:THR:HG23	1.92	0.70
1:A:67:LYS:NZ	1:A:76:HIS:HE1	1.89	0.70
1:A:92:ASP:OD1	2:A:2444:HOH:O	2.08	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:GLU:O	1:B:45[A]:THR:HG22	1.92	0.69
1:C:5:VAL:O	1:C:5:VAL:CG1	2.39	0.69
1:A:67:LYS:HZ2	1:A:76:HIS:HE1	1.42	0.68
1:D:47:ASN:HB2	2:D:2168:HOH:O	1.95	0.67
1:D:33:ARG:CA	2:D:2077:HOH:O	2.42	0.66
1:D:3:GLN:NE2	1:D:5[B]:VAL:HG23	2.09	0.66
1:C:58[A]:ARG:HD2	2:C:2173:HOH:O	1.92	0.66
1:D:3:GLN:HG3	1:D:5[A]:VAL:HG12	1.78	0.66
1:D:24:HIS:HE1	1:D:29:ASP:OD1	1.78	0.66
1:A:3:GLN:HB3	1:D:52:ARG:NH1	2.10	0.65
1:C:24:HIS:HE1	1:C:29:ASP:OD2	1.80	0.65
1:D:97:VAL:C	1:D:98:LYS:O	2.35	0.65
1:A:17[A]:LYS:HZ2	1:A:20:GLU:CB	2.10	0.64
1:C:5:VAL:CG1	2:C:2027:HOH:O	2.39	0.63
1:D:33:ARG:HA	2:D:2077:HOH:O	1.98	0.63
1:D:58:ARG:NH1	1:D:91:PHE:CE2	2.66	0.63
1:A:5:VAL:O	1:D:3:GLN:HG2	1.98	0.63
1:B:6:THR:HG22	1:B:10[B]:SER:HB2	1.79	0.63
1:D:3:GLN:CG	1:D:5[A]:VAL:CG1	2.75	0.62
1:A:4:THR:HG22	2:A:2308:HOH:O	1.93	0.62
1:B:59:ASP:CB	2:B:2222:HOH:O	2.32	0.62
1:D:93:PRO:HB3	1:D:97:VAL:HG22	1.80	0.62
1:D:3:GLN:HE21	1:D:5[A]:VAL:CG1	2.07	0.62
1:A:3:GLN:HB2	1:A:4:THR:CA	2.29	0.62
1:A:3:GLN:O	1:A:3:GLN:HB2	1.78	0.61
1:D:52:ARG:CB	2:D:2322:HOH:O	1.90	0.61
1:A:3:GLN:N	1:A:3:GLN:CG	2.63	0.60
1:B:67:LYS:HZ1	1:B:76:HIS:HE1	1.49	0.60
1:C:95:PRO:C	1:C:98:LYS:HB3	2.18	0.59
1:B:22:ILE:HD12	1:B:89[B]:LEU:CD1	2.33	0.58
1:A:97:VAL:HG23	2:A:2073:HOH:O	2.03	0.58
1:A:3:GLN:O	1:A:3:GLN:CD	2.41	0.58
1:D:33:ARG:CA	1:D:33:ARG:NE	2.68	0.57
1:D:50:TRP:CZ3	1:D:61[B]:ILE:HD13	2.39	0.57
1:D:33:ARG:HA	1:D:33:ARG:NE	2.20	0.56
1:A:3:GLN:OE1	1:A:3:GLN:NE2	2.39	0.56
1:A:4:THR:HG21	1:A:6[B]:THR:HG23	1.82	0.56
1:C:98:LYS:N	2:C:2462:HOH:O	2.09	0.56
1:C:65[B]:LYS:HE3	2:C:1907:HOH:O	1.95	0.55
1:A:17[B]:LYS:NZ	1:A:20:GLU:CG	2.50	0.54
1:D:33:ARG:HH22	1:D:36:VAL:HG21	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:95:PRO:C	1:C:98:LYS:HB2	2.19	0.54
1:A:24:HIS:HD2	2:A:2100:HOH:O	1.91	0.53
1:C:78:ASP:OD1	2:C:2443:HOH:O	2.18	0.53
1:D:98:LYS:C	2:D:2225:HOH:O	2.46	0.53
1:B:67:LYS:NZ	1:B:76:HIS:CE1	2.73	0.53
1:D:76:HIS:HE1	2:D:2144:HOH:O	1.91	0.53
1:A:3:GLN:N	1:D:52:ARG:NE	2.41	0.53
1:B:76:HIS:HD2	2:B:1952:HOH:O	1.92	0.53
1:D:33:ARG:HA	1:D:33:ARG:CZ	2.39	0.53
1:A:76:HIS:HD2	2:A:1969:HOH:O	1.92	0.52
1:A:58[B]:ARG:HG2	2:A:2177:HOH:O	2.09	0.52
1:A:67:LYS:NZ	1:A:76:HIS:CE1	2.73	0.52
1:A:3:GLN:HG3	1:D:52:ARG:HD2	1.91	0.51
1:A:5:VAL:HB	1:D:3:GLN:CG	2.38	0.51
1:A:7:THR:HB	1:A:53[A]:ASP:HB3	1.92	0.51
1:A:17[A]:LYS:HZ1	1:A:20:GLU:CB	2.22	0.51
2:B:2455:HOH:O	1:C:52:ARG:HD3	2.11	0.51
1:A:17[A]:LYS:HZ1	1:A:20:GLU:HB3	1.76	0.50
1:D:33:ARG:HH22	1:D:36:VAL:CG2	2.25	0.49
1:D:94:PRO:HD2	1:D:97:VAL:HG13	1.94	0.49
1:D:97:VAL:O	1:D:98:LYS:C	2.28	0.49
1:B:67:LYS:HZ2	1:B:76:HIS:CE1	2.31	0.49
1:B:42:GLU:O	1:B:45[B]:THR:HG22	2.13	0.48
1:C:11[A]:LEU:HG	1:C:96:TRP:CE3	2.48	0.48
1:D:34:ARG:NH2	2:D:2219:HOH:O	2.42	0.48
1:D:3:GLN:HE22	1:D:5[B]:VAL:HG23	1.78	0.48
1:D:50:TRP:HZ3	1:D:61[B]:ILE:HD13	1.78	0.48
1:A:3:GLN:C	1:A:3:GLN:CG	2.81	0.47
1:B:24:HIS:HD2	2:B:2116:HOH:O	1.96	0.47
1:B:59:ASP:HB2	2:B:2222:HOH:O	2.08	0.47
1:C:92:ASP:HB2	2:C:2386:HOH:O	2.14	0.47
1:D:31:LYS:HA	1:D:31:LYS:HE3	1.97	0.47
2:C:2437:HOH:O	1:D:21[A]:ARG:HD2	2.14	0.46
1:D:3:GLN:NE2	1:D:5[B]:VAL:CG2	2.78	0.46
1:D:32:LYS:HB3	1:D:33:ARG:NH2	2.22	0.46
1:A:3:GLN:HG2	1:D:52:ARG:HD2	1.97	0.46
1:A:3:GLN:O	1:A:4:THR:HB	2.08	0.46
1:A:3:GLN:HB3	1:D:52:ARG:HH11	1.79	0.45
1:B:98:LYS:HD2	1:B:98:LYS:HA	1.43	0.45
1:C:24:HIS:CD2	1:D:24:HIS:CD2	2.71	0.45
1:D:33:ARG:C	2:D:2077:HOH:O	2.35	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:61[B]:ILE:HD12	1:D:86:TRP:HB3	1.99	0.45
1:A:17[B]:LYS:NZ	1:A:20:GLU:CB	2.81	0.43
1:C:59:ASP:OD1	2:C:2299:HOH:O	2.20	0.43
1:C:29:ASP:HB2	1:D:21[B]:ARG:NH2	2.34	0.43
1:A:28:VAL:HB	1:A:81:PRO:HB3	2.01	0.43
1:B:22:ILE:CD1	1:B:89[B]:LEU:HD12	2.39	0.43
1:D:61[B]:ILE:CD1	1:D:86:TRP:HB3	2.49	0.42
1:A:17[A]:LYS:NZ	1:A:20:GLU:HB3	2.30	0.42
1:A:67:LYS:HZ3	1:A:76:HIS:CE1	2.37	0.42
1:B:58:ARG:NH1	2:B:2079:HOH:O	2.52	0.42
1:D:24:HIS:CE1	1:D:29:ASP:OD1	2.67	0.42
1:A:3:GLN:O	1:A:3:GLN:OE1	2.39	0.41
1:B:58:ARG:NH2	2:B:2286:HOH:O	2.52	0.41
1:C:65[B]:LYS:NZ	2:C:1907:HOH:O	1.76	0.41
1:D:33:ARG:NE	1:D:33:ARG:N	2.68	0.41
1:D:53[B]:ASP:CG	2:D:2254:HOH:O	2.55	0.41
1:D:33:ARG:NH2	1:D:36:VAL:CG2	2.84	0.41
1:D:52:ARG:CG	2:D:2322:HOH:O	2.44	0.41
1:C:29:ASP:HB2	1:D:21[B]:ARG:HH21	1.86	0.41
1:A:14:GLY:O	1:D:33:ARG:NE	2.54	0.41
1:A:17[A]:LYS:CD	2:A:2243:HOH:O	2.67	0.41
1:D:50:TRP:CZ3	1:D:61[B]:ILE:CD1	3.04	0.41

All (7) 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)
2:A:2052:HOH:O	2:A:2290:HOH:O[1_455]	0.69	1.51
2:B:2198:HOH:O	2:C:2142:HOH:O[1_545]	1.57	0.63
2:B:2112:HOH:O	2:C:2142:HOH:O[1_545]	1.78	0.42
1:C:5:VAL:CB	2:A:2280:HOH:O[1_556]	1.88	0.32
2:A:2340:HOH:O	2:D:2180:HOH:O[1_645]	2.11	0.09
2:A:2281:HOH:O	2:A:2290:HOH:O[1_455]	2.15	0.05
2:A:2257:HOH:O	2:D:2199:HOH:O[1_645]	2.19	0.01

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	101/100 (101%)	99 (98%)	1 (1%)	1 (1%)	15 2
1	B	97/100 (97%)	97 (100%)	0	0	100 100
1	C	97/100 (97%)	97 (100%)	0	0	100 100
1	D	102/100 (102%)	102 (100%)	0	0	100 100
All	All	397/400 (99%)	395 (100%)	1 (0%)	1 (0%)	41 14

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	THR

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	94/89 (106%)	87 (93%)	7 (7%)	13 1
1	B	90/89 (101%)	87 (97%)	3 (3%)	38 8
1	C	90/89 (101%)	85 (94%)	5 (6%)	21 2
1	D	94/89 (106%)	89 (95%)	5 (5%)	22 2
All	All	368/356 (103%)	348 (95%)	20 (5%)	24 2

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	4	THR
1	A	17[A]	LYS
1	A	17[B]	LYS
1	A	52[A]	ARG
1	A	52[B]	ARG
1	A	97	VAL
1	B	91[A]	PHE
1	B	91[B]	PHE
1	B	98	LYS
1	C	6	THR
1	C	9	LEU
1	C	13	LEU
1	C	28	VAL
1	C	98	LYS
1	D	9	LEU
1	D	31	LYS
1	D	33	ARG
1	D	58	ARG
1	D	97	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	HIS
1	A	74	HIS
1	A	76	HIS
1	B	15	HIS
1	B	24	HIS
1	B	76	HIS
1	C	15	HIS
1	C	24	HIS
1	C	47	ASN
1	C	63	GLN
1	D	3	GLN
1	D	15	HIS
1	D	24	HIS
1	D	74	HIS
1	D	76	HIS

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.