



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:39 AM UTC

PDB ID : 8CIT / pdb_00008cit
Title : The FERM domain of human moesin mutant L281R
Authors : Bradshaw, W.J.; Katis, V.L.; Koekemoer, L.; Bountra, C.; von Delft, F.;
Brennan, P.E.
Deposited on : 2023-02-10
Resolution : 2.54 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

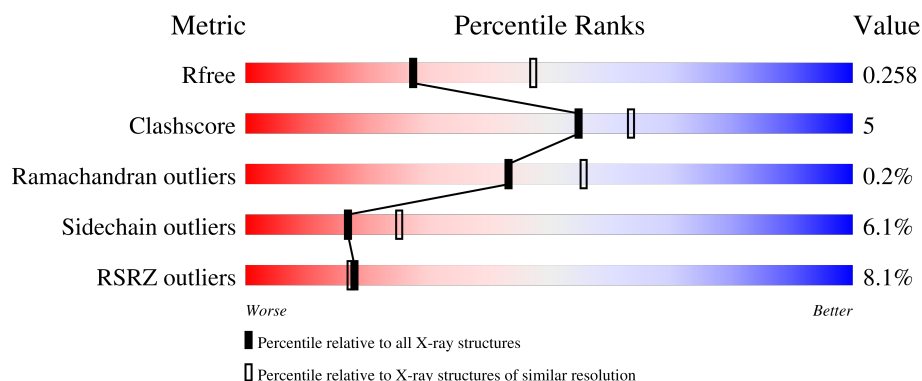
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	347	5% 79% 17% ..
1	B	347	7% 80% 16% ..
1	C	347	11% 79% 14% 5%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 8259 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 Moesin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	336	Total	C	N	O	S	0	0	0
			2749	1766	478	494	11			
1	B	341	Total	C	N	O	S	0	0	0
			2826	1809	490	516	11			
1	C	328	Total	C	N	O	S	0	2	0
			2684	1726	466	483	9			

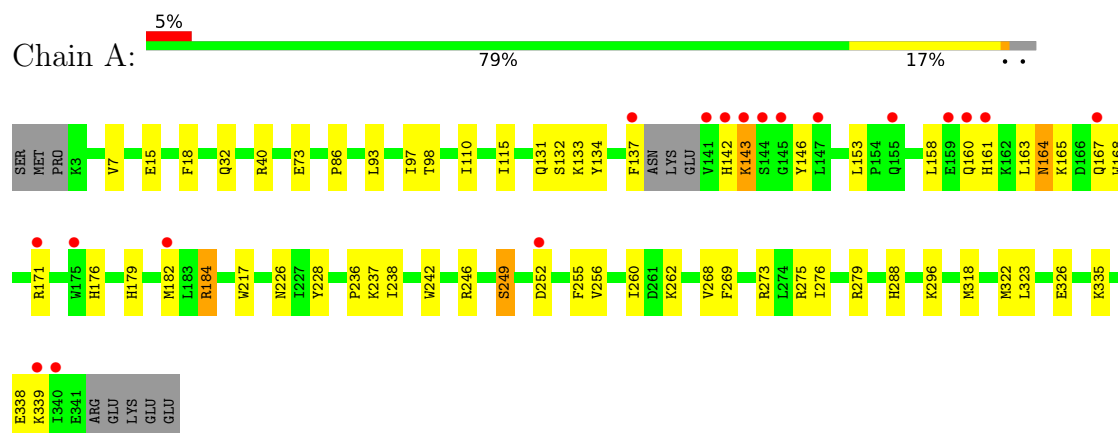
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP P26038
A	281	ARG	LEU	engineered mutation	UNP P26038
B	0	SER	-	expression tag	UNP P26038
B	281	ARG	LEU	engineered mutation	UNP P26038
C	0	SER	-	expression tag	UNP P26038
C	281	ARG	LEU	engineered mutation	UNP P26038

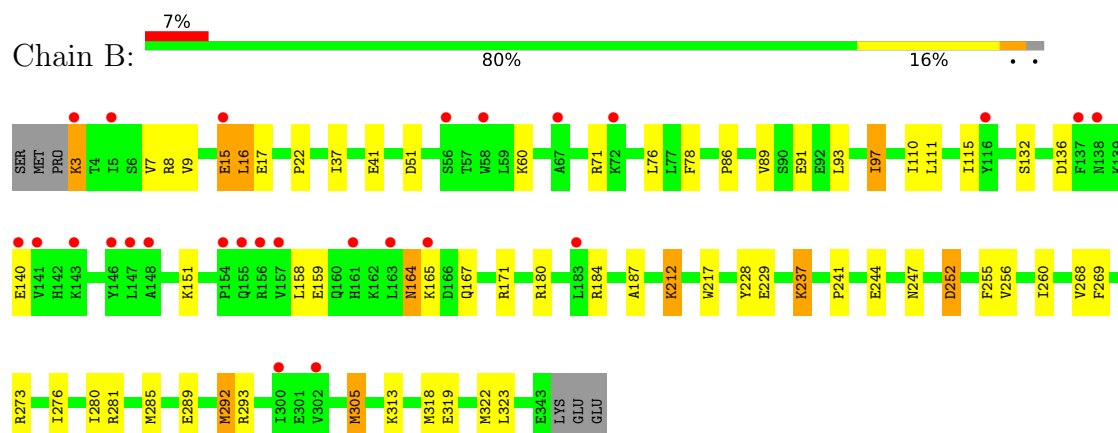
3 Residue-property plots [i](#)

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

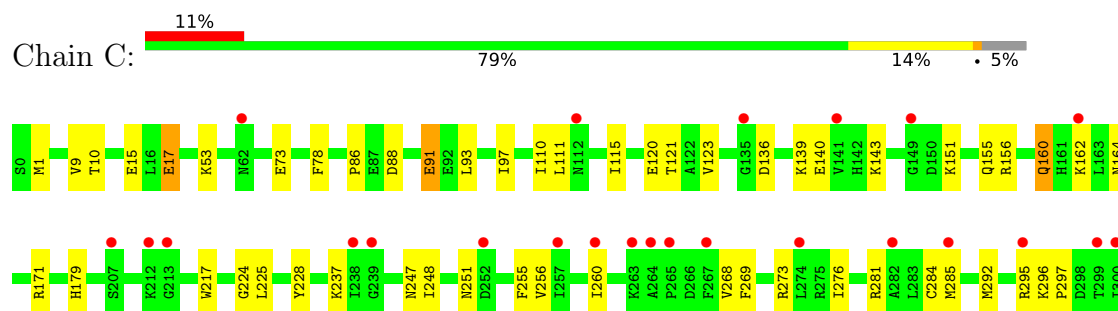
• Molecule 1: Moesin



• Molecule 1: Moesin



• Molecule 1: Moesin



E301	E311	Q317	M318	E319	R320	A321	M322	L323	E324	M325	E326	E327	LYS	LYS	ARG	GLU	MET	ALA	GLU	LYS	LYS	GLU	LYS	ILE	GLU	ARG	GLU	LYS	GLU	GLU
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.70Å 74.05Å 115.08Å 90.00° 95.63° 90.00°	Depositor
Resolution (Å)	62.18 – 2.54 62.18 – 2.54	Depositor EDS
% Data completeness (in resolution range)	58.5 (62.18-2.54) 58.5 (62.18-2.54)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 2.55Å)	Xtriage
Refinement program	BUSTER 2.10.4	Depositor
R, R_{free}	0.244 , 0.287 (Not available) , 0.258	Depositor DCC
R_{free} test set	1173 reflections (2.89%)	wwPDB-VP
Wilson B-factor (Å ²)	28.0	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 30.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	8259	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 4.33% 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	0.81	0/2806	1.17	1/3776 (0.0%)
1	B	0.82	2/2884 (0.1%)	1.18	13/3877 (0.3%)
1	C	0.81	1/2749 (0.0%)	1.19	9/3707 (0.2%)
All	All	0.81	3/8439 (0.0%)	1.18	23/11360 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	251	ASN	CA-C	9.62	1.57	1.52
1	B	292	MET	SD-CE	6.58	1.96	1.79
1	B	305	MET	SD-CE	-6.45	1.63	1.79

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	51	ASP	CA-CB-CG	7.11	119.71	112.60
1	C	17	GLU	CB-CG-CD	6.75	124.08	112.60
1	C	155	GLN	CA-C-N	6.70	129.92	120.28
1	C	155	GLN	C-N-CA	6.70	129.92	120.28
1	C	248	ILE	N-CA-CB	6.23	119.52	111.67
1	B	15	GLU	CB-CG-CD	6.16	123.06	112.60
1	B	187	ALA	CA-C-N	6.01	128.70	120.46
1	B	187	ALA	C-N-CA	6.01	128.70	120.46
1	C	164	ASN	CA-CB-CG	5.96	118.56	112.60
1	B	252	ASP	CA-CB-CG	5.86	118.46	112.60
1	C	247	ASN	CA-CB-CG	5.85	118.45	112.60
1	A	246	ARG	N-CA-C	-5.62	105.07	111.14
1	B	247	ASN	CA-CB-CG	5.55	118.15	112.60
1	C	73	GLU	N-CA-C	5.49	117.75	109.41
1	C	91	GLU	N-CA-C	5.37	116.82	110.97
1	B	164	ASN	CA-CB-CG	5.34	117.94	112.60
1	C	224	GLY	N-CA-C	5.27	117.70	110.58
1	B	293	ARG	CA-C-N	5.20	127.25	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	293	ARG	C-N-CA	5.20	127.25	120.28
1	B	97	ILE	CA-C-N	5.18	127.64	120.29
1	B	97	ILE	C-N-CA	5.18	127.64	120.29
1	B	89	VAL	CA-C-N	5.11	128.86	120.63
1	B	89	VAL	C-N-CA	5.11	128.86	120.63

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	2749	0	2737	35	0
1	B	2826	0	2836	26	0
1	C	2684	0	2670	20	0
All	All	8259	0	8243	79	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 5.

All (79) 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:137:PHE:HA	1:A:142:HIS:CD2	2.00	0.95
1:B:8:ARG:HH11	1:B:15:GLU:HG3	1.29	0.92
1:A:143:LYS:NZ	1:A:146:TYR:HB2	2.04	0.72
1:A:143:LYS:HB3	1:A:146:TYR:HB3	1.71	0.72
1:C:111:LEU:HD13	1:C:151:LYS:HG3	1.74	0.70
1:C:120[B]:GLU:HG2	1:C:121:THR:N	2.07	0.69
1:B:8:ARG:NH1	1:B:15:GLU:HG3	2.07	0.69
1:A:143:LYS:HZ1	1:A:146:TYR:HB2	1.59	0.67
1:B:281:ARG:O	1:B:285:MET:HG3	1.97	0.64
1:C:120[B]:GLU:HG2	1:C:121:THR:H	1.64	0.63
1:A:32:GLN:HG2	1:C:1:MET:HE2	1.79	0.63
1:B:9:VAL:HG22	1:B:78:PHE:HB2	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:VAL:HG12	1:B:76:LEU:HB2	1.82	0.61
1:B:217:TRP:HB2	1:B:228:TYR:HB2	1.85	0.59
1:A:217:TRP:HB2	1:A:228:TYR:HB2	1.85	0.59
1:B:3:LYS:O	1:B:22:PRO:HD3	2.02	0.58
1:B:132:SER:O	1:B:184:ARG:HG3	2.04	0.58
1:A:131:GLN:HG2	1:A:182:MET:HE3	1.85	0.58
1:A:318:MET:O	1:A:322:MET:HG3	2.04	0.57
1:C:217:TRP:HB2	1:C:228:TYR:HB2	1.86	0.57
1:B:111:LEU:HB3	1:B:151:LYS:HE3	1.87	0.57
1:C:273:ARG:HB2	1:C:276:ILE:HD13	1.85	0.57
1:B:276:ILE:O	1:B:280:ILE:HG13	2.05	0.56
1:B:318:MET:O	1:B:322:MET:HG3	2.06	0.56
1:B:256:VAL:HG22	1:B:268:VAL:HG22	1.86	0.55
1:C:86:PRO:HG3	1:C:93:LEU:HD21	1.89	0.55
1:C:9:VAL:HG22	1:C:78:PHE:HB2	1.88	0.55
1:C:256:VAL:HG22	1:C:268:VAL:HG22	1.89	0.54
1:B:8:ARG:HG3	1:B:17:GLU:HG3	1.89	0.54
1:A:273:ARG:HB2	1:A:276:ILE:HD13	1.89	0.54
1:C:296:LYS:HG2	1:C:297:PRO:HD2	1.90	0.54
1:B:319:GLU:HA	1:B:322:MET:HE2	1.89	0.54
1:B:229:GLU:HG2	1:B:237:LYS:HG3	1.91	0.53
1:A:179:HIS:HB3	1:A:182:MET:HE2	1.89	0.53
1:A:86:PRO:HG3	1:A:93:LEU:HD21	1.91	0.52
1:A:256:VAL:HG22	1:A:268:VAL:HG22	1.91	0.51
1:B:164:ASN:HB3	1:B:167:GLN:HG3	1.91	0.51
1:A:153:LEU:HB2	1:A:158:LEU:HD21	1.93	0.51
1:C:10:THR:HG22	1:C:15:GLU:HG3	1.93	0.50
1:C:292:MET:HA	1:C:295:ARG:HG2	1.92	0.50
1:B:273:ARG:HB2	1:B:276:ILE:HD13	1.93	0.50
1:B:86:PRO:HG3	1:B:93:LEU:HD21	1.95	0.48
1:A:137:PHE:CD2	1:A:142:HIS:HB3	2.48	0.48
1:A:137:PHE:HA	1:A:142:HIS:NE2	2.29	0.48
1:C:111:LEU:HB3	1:C:151:LYS:HE3	1.95	0.48
1:C:156:ARG:HG3	1:C:160:GLN:HE22	1.78	0.48
1:B:15:GLU:O	1:B:15:GLU:HG2	2.14	0.47
1:B:41:GLU:HG2	1:B:305:MET:HE1	1.95	0.47
1:B:111:LEU:HD13	1:B:151:LYS:HG2	1.96	0.47
1:A:137:PHE:HB2	1:A:176:HIS:CE1	2.50	0.47
1:B:212:LYS:HE3	1:B:212:LYS:HB3	1.74	0.47
1:A:132:SER:O	1:A:184:ARG:HG3	2.15	0.47
1:A:161:HIS:HB2	1:A:168:TRP:CZ2	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:LYS:HD3	1:A:134:TYR:CE1	2.51	0.45
1:C:110:ILE:HD13	1:C:115:ILE:HB	1.99	0.45
1:C:281:ARG:O	1:C:285:MET:HG3	2.17	0.45
1:B:110:ILE:HD13	1:B:115:ILE:HB	1.99	0.45
1:B:16:LEU:HD11	1:B:37:ILE:HG23	1.98	0.45
1:A:335:LYS:HA	1:A:338:GLU:HG2	1.97	0.44
1:C:88:ASP:HB3	1:C:91:GLU:HB2	1.99	0.44
1:A:143:LYS:HZ3	1:A:146:TYR:HB2	1.79	0.44
1:A:255:PHE:CZ	1:A:269:PHE:HB2	2.53	0.44
1:A:161:HIS:HB2	1:A:168:TRP:HZ2	1.83	0.43
1:B:255:PHE:CZ	1:B:269:PHE:HB2	2.53	0.43
1:A:131:GLN:HG2	1:A:182:MET:CE	2.48	0.43
1:A:164:ASN:OD1	1:A:167:GLN:HG3	2.18	0.42
1:A:275:ARG:O	1:A:279:ARG:HG3	2.18	0.42
1:A:110:ILE:HD13	1:A:115:ILE:HB	2.00	0.42
1:A:18:PHE:HB3	1:C:1:MET:HE1	2.01	0.42
1:A:226:ASN:HB3	1:A:236:PRO:HB3	2.02	0.42
1:B:241:PRO:HG2	1:B:244:GLU:HG3	2.01	0.42
1:A:242:TRP:HB3	1:A:288:HIS:HD2	1.84	0.41
1:A:335:LYS:O	1:A:339:LYS:HG3	2.21	0.41
1:A:323:LEU:HD23	1:A:323:LEU:HA	1.98	0.41
1:A:249:SER:O	1:A:255:PHE:HA	2.21	0.41
1:A:318:MET:HE2	1:A:318:MET:HB3	1.95	0.40
1:C:225:LEU:HD11	1:C:284:CYS:SG	2.61	0.40
1:C:255:PHE:CZ	1:C:269:PHE:HB2	2.56	0.40
1:A:93:LEU:HD22	1:A:98:THR:HG22	2.04	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	332/347 (96%)	320 (96%)	11 (3%)	1 (0%)	36	45
1	B	339/347 (98%)	323 (95%)	15 (4%)	1 (0%)	36	45
1	C	328/347 (94%)	318 (97%)	10 (3%)	0	100	100
All	All	999/1041 (96%)	961 (96%)	36 (4%)	2 (0%)	43	56

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	252	ASP
1	B	71	ARG

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	288/314 (92%)	269 (93%)	19 (7%)	15	21
1	B	303/314 (96%)	283 (93%)	20 (7%)	15	21
1	C	282/314 (90%)	268 (95%)	14 (5%)	22	33
All	All	873/942 (93%)	820 (94%)	53 (6%)	17	25

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	15	GLU
1	A	40	ARG
1	A	73	GLU
1	A	97	ILE
1	A	143	LYS
1	A	160	GLN
1	A	163	LEU
1	A	164	ASN
1	A	165	LYS
1	A	171	ARG

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Mol	Chain	Res	Type
1	A	184	ARG
1	A	237	LYS
1	A	238	ILE
1	A	249	SER
1	A	260	ILE
1	A	262	LYS
1	A	296	LYS
1	A	326	GLU
1	B	3	LYS
1	B	16	LEU
1	B	60	LYS
1	B	91	GLU
1	B	97	ILE
1	B	136	ASP
1	B	140	GLU
1	B	158	LEU
1	B	159	GLU
1	B	165	LYS
1	B	171	ARG
1	B	180	ARG
1	B	212	LYS
1	B	237	LYS
1	B	252	ASP
1	B	260	ILE
1	B	289	GLU
1	B	292	MET
1	B	313	LYS
1	B	323	LEU
1	C	17	GLU
1	C	53	LYS
1	C	97	ILE
1	C	123	VAL
1	C	136	ASP
1	C	139	LYS
1	C	140	GLU
1	C	143	LYS
1	C	160	GLN
1	C	162	LYS
1	C	171	ARG
1	C	179	HIS
1	C	237	LYS
1	C	260	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	50	GLN
1	A	142	HIS
1	A	161	HIS
1	A	226	ASN
1	A	288	HIS
1	A	317	GLN
1	A	325	ASN
1	B	50	GLN
1	B	62	ASN
1	B	304	GLN
1	B	325	ASN
1	C	50	GLN
1	C	138	ASN
1	C	160	GLN
1	C	226	ASN
1	C	230	GLN

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 ⓘ

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	336/347 (96%)	0.59	18 (5%) 31 31	10, 29, 67, 83	0
1	B	341/347 (98%)	0.83	26 (7%) 20 19	15, 42, 63, 83	0
1	C	328/347 (94%)	0.85	37 (11%) 10 9	14, 38, 71, 93	2 (0%)
All	All	1005/1041 (96%)	0.75	81 (8%) 18 17	10, 35, 67, 93	2 (0%)

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	141	VAL	5.4
1	C	327	LYS	4.8
1	B	161	HIS	4.3
1	C	319	GLU	3.6
1	B	141	VAL	3.6
1	B	143	LYS	3.5
1	C	324	GLU	3.5
1	C	267	PHE	3.4
1	C	260	ILE	3.4
1	B	67	ALA	3.4
1	B	148	ALA	3.4
1	B	163	LEU	3.3
1	C	285	MET	3.3
1	C	239	GLY	3.2
1	B	58	TRP	3.1
1	A	340	ILE	3.0
1	C	264	ALA	3.0
1	A	339	LYS	3.0
1	A	145	GLY	3.0
1	A	144	SER	2.9
1	B	165	LYS	2.9
1	C	325	ASN	2.8
1	B	147	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	149	GLY	2.7
1	C	238	ILE	2.7
1	C	318	MET	2.7
1	A	167	GLN	2.7
1	C	265	PRO	2.7
1	B	156	ARG	2.7
1	C	301	GLU	2.7
1	C	317	GLN	2.7
1	A	137	PHE	2.7
1	C	320	ARG	2.7
1	B	300	ILE	2.6
1	A	142	HIS	2.6
1	C	326	GLU	2.6
1	A	161	HIS	2.6
1	C	263	LYS	2.6
1	C	323	LEU	2.5
1	C	212	LYS	2.5
1	A	175	TRP	2.5
1	C	322	MET	2.5
1	A	171	ARG	2.4
1	C	207	SER	2.4
1	A	143	LYS	2.4
1	A	182	MET	2.4
1	B	15	GLU	2.4
1	B	157	VAL	2.4
1	C	311	GLU	2.4
1	B	154	PRO	2.3
1	C	300	ILE	2.3
1	B	140	GLU	2.3
1	B	138	ASN	2.3
1	C	282	ALA	2.3
1	C	295	ARG	2.3
1	C	299	THR	2.3
1	B	5	ILE	2.3
1	B	183	LEU	2.3
1	A	252	ASP	2.3
1	C	257	ILE	2.2
1	C	141	VAL	2.2
1	A	155	GLN	2.2
1	A	147	LEU	2.2
1	B	56	SER	2.2
1	B	137	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	160	GLN	2.1
1	C	135	GLY	2.1
1	C	213	GLY	2.1
1	B	116	TYR	2.1
1	C	162	LYS	2.1
1	B	302	VAL	2.0
1	C	321	ALA	2.0
1	C	112	ASN	2.0
1	C	274	LEU	2.0
1	A	159	GLU	2.0
1	B	3	LYS	2.0
1	B	72	LYS	2.0
1	C	252	ASP	2.0
1	B	155	GLN	2.0
1	C	62	ASN	2.0
1	B	146	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.