



# wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 24, 2024 – 12:51 AM EDT

PDB ID : 2PJQ  
Title : Crystal structure of Q88U62\_LACPL from Lactobacillus plantarum. Northeast Structural Genomics target LpR71  
Authors : Benach, J.; Su, M.; Seetharaman, J.; Forouhar, F.; Chen, C.X.; Cunningham, K.; Ma, L.-C.; Owens, L.; Baran, M.; Acton, T.B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2007-04-16  
Resolution : 2.80 Å(reported)

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We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

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<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>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

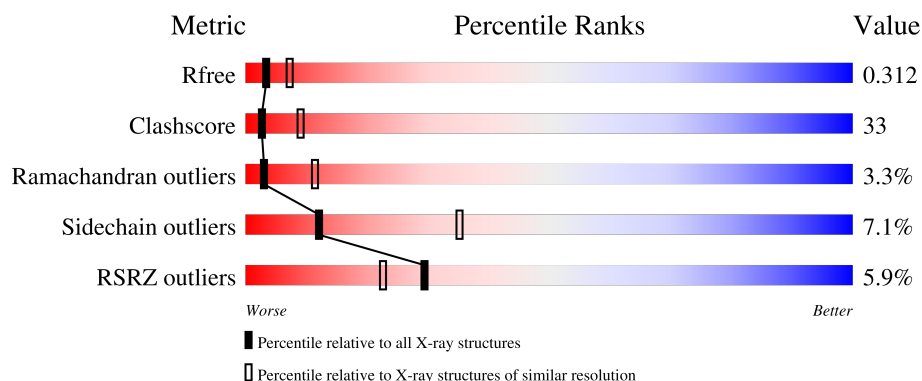
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	164625	3657 (2.80-2.80)
Clashscore	180529	4123 (2.80-2.80)
Ramachandran outliers	177936	4071 (2.80-2.80)
Sidechain outliers	177891	4073 (2.80-2.80)
RSRZ outliers	164620	3659 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	231	<div> <div>2%</div> <div> <div></div> <div>48%</div> <div>33%</div> <div>8%</div> <div>11%</div> </div> </div>
1	B	231	<div> <div>4%</div> <div> <div></div> <div>49%</div> <div>37%</div> <div>• • 8%</div> </div> </div>
1	C	231	<div> <div>8%</div> <div> <div></div> <div>39%</div> <div>45%</div> <div>6%</div> <div>• 8%</div> </div> </div>
1	D	231	<div> <div>6%</div> <div> <div></div> <div>37%</div> <div>48%</div> <div>• • 12%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6575 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 Uncharacterized protein lp\_2664.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	Se	0	0	0
			1625	1019	298	303	5			
1	B	212	Total	C	N	O	Se	0	0	0
			1667	1047	306	308	6			
1	C	212	Total	C	N	O	Se	0	0	0
			1667	1045	306	310	6			
1	D	203	Total	C	N	O	Se	0	0	0
			1604	1008	295	296	5			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	MSE	-	cloning artifact	UNP Q88U62
A	-3	ALA	-	cloning artifact	UNP Q88U62
A	-2	GLY	-	cloning artifact	UNP Q88U62
A	-1	ASP	-	cloning artifact	UNP Q88U62
A	0	PRO	-	cloning artifact	UNP Q88U62
A	1	MSE	MET	modified residue	UNP Q88U62
A	66	MSE	MET	modified residue	UNP Q88U62
A	100	MSE	MET	modified residue	UNP Q88U62
A	158	MSE	MET	modified residue	UNP Q88U62
A	185	MSE	MET	modified residue	UNP Q88U62
A	201	MSE	MET	modified residue	UNP Q88U62
A	219	LEU	-	cloning artifact	UNP Q88U62
A	220	GLU	-	cloning artifact	UNP Q88U62
A	221	HIS	-	cloning artifact	UNP Q88U62
A	222	HIS	-	cloning artifact	UNP Q88U62
A	223	HIS	-	cloning artifact	UNP Q88U62
A	224	HIS	-	cloning artifact	UNP Q88U62
A	225	HIS	-	cloning artifact	UNP Q88U62
A	226	HIS	-	cloning artifact	UNP Q88U62
B	-4	MSE	-	cloning artifact	UNP Q88U62
B	-3	ALA	-	cloning artifact	UNP Q88U62

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	GLY	-	cloning artifact	UNP Q88U62
B	-1	ASP	-	cloning artifact	UNP Q88U62
B	0	PRO	-	cloning artifact	UNP Q88U62
B	1	MSE	MET	modified residue	UNP Q88U62
B	66	MSE	MET	modified residue	UNP Q88U62
B	100	MSE	MET	modified residue	UNP Q88U62
B	158	MSE	MET	modified residue	UNP Q88U62
B	185	MSE	MET	modified residue	UNP Q88U62
B	201	MSE	MET	modified residue	UNP Q88U62
B	219	LEU	-	cloning artifact	UNP Q88U62
B	220	GLU	-	cloning artifact	UNP Q88U62
B	221	HIS	-	cloning artifact	UNP Q88U62
B	222	HIS	-	cloning artifact	UNP Q88U62
B	223	HIS	-	cloning artifact	UNP Q88U62
B	224	HIS	-	cloning artifact	UNP Q88U62
B	225	HIS	-	cloning artifact	UNP Q88U62
B	226	HIS	-	cloning artifact	UNP Q88U62
C	-4	MSE	-	cloning artifact	UNP Q88U62
C	-3	ALA	-	cloning artifact	UNP Q88U62
C	-2	GLY	-	cloning artifact	UNP Q88U62
C	-1	ASP	-	cloning artifact	UNP Q88U62
C	0	PRO	-	cloning artifact	UNP Q88U62
C	1	MSE	MET	modified residue	UNP Q88U62
C	66	MSE	MET	modified residue	UNP Q88U62
C	100	MSE	MET	modified residue	UNP Q88U62
C	158	MSE	MET	modified residue	UNP Q88U62
C	185	MSE	MET	modified residue	UNP Q88U62
C	201	MSE	MET	modified residue	UNP Q88U62
C	219	LEU	-	cloning artifact	UNP Q88U62
C	220	GLU	-	cloning artifact	UNP Q88U62
C	221	HIS	-	cloning artifact	UNP Q88U62
C	222	HIS	-	cloning artifact	UNP Q88U62
C	223	HIS	-	cloning artifact	UNP Q88U62
C	224	HIS	-	cloning artifact	UNP Q88U62
C	225	HIS	-	cloning artifact	UNP Q88U62
C	226	HIS	-	cloning artifact	UNP Q88U62
D	-4	MSE	-	cloning artifact	UNP Q88U62
D	-3	ALA	-	cloning artifact	UNP Q88U62
D	-2	GLY	-	cloning artifact	UNP Q88U62
D	-1	ASP	-	cloning artifact	UNP Q88U62
D	0	PRO	-	cloning artifact	UNP Q88U62
D	1	MSE	MET	modified residue	UNP Q88U62

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Chain	Residue	Modelled	Actual	Comment	Reference
D	66	MSE	MET	modified residue	UNP Q88U62
D	100	MSE	MET	modified residue	UNP Q88U62
D	158	MSE	MET	modified residue	UNP Q88U62
D	185	MSE	MET	modified residue	UNP Q88U62
D	201	MSE	MET	modified residue	UNP Q88U62
D	219	LEU	-	cloning artifact	UNP Q88U62
D	220	GLU	-	cloning artifact	UNP Q88U62
D	221	HIS	-	cloning artifact	UNP Q88U62
D	222	HIS	-	cloning artifact	UNP Q88U62
D	223	HIS	-	cloning artifact	UNP Q88U62
D	224	HIS	-	cloning artifact	UNP Q88U62
D	225	HIS	-	cloning artifact	UNP Q88U62
D	226	HIS	-	cloning artifact	UNP Q88U62

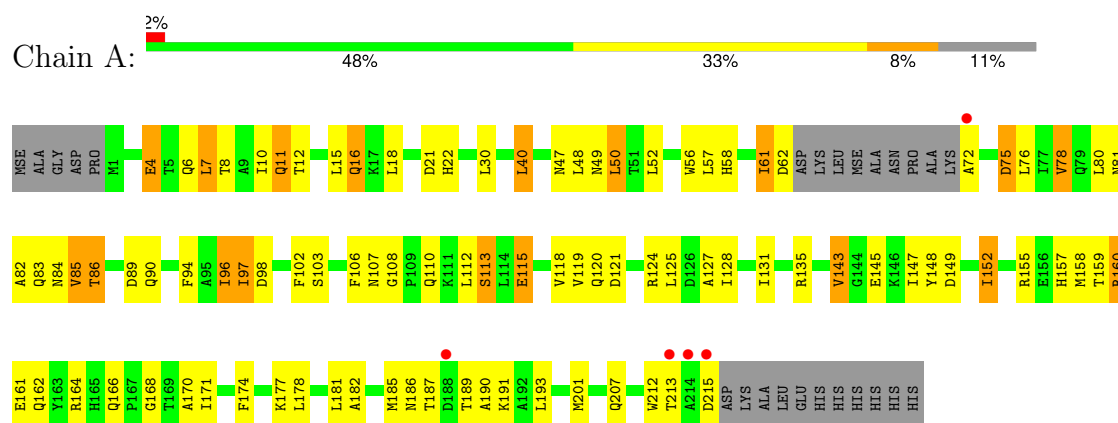
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total O 2 2	0	0
2	B	6	Total O 6 6	0	0
2	C	3	Total O 3 3	0	0
2	D	1	Total O 1 1	0	0

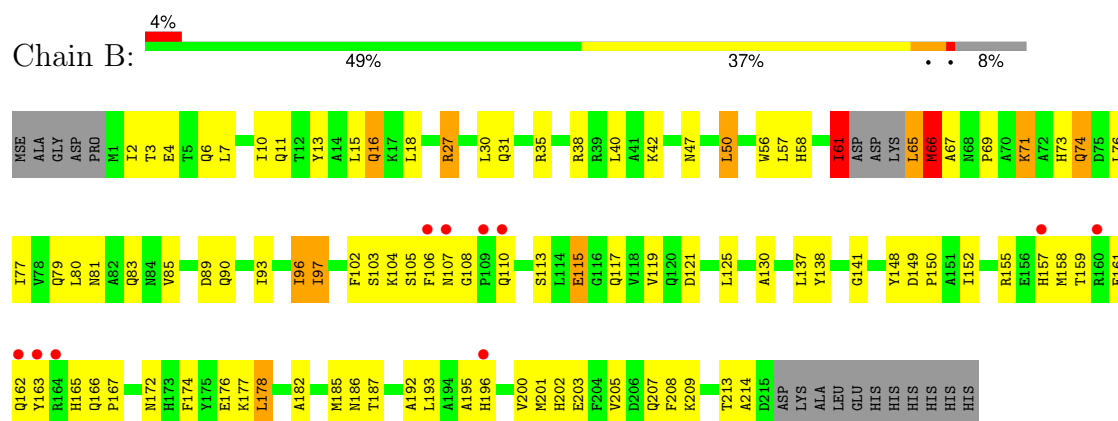
### 3 Residue-property plots

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

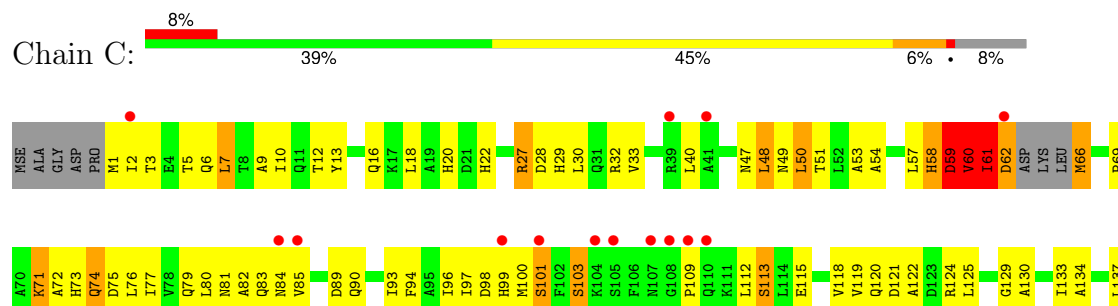
#### • Molecule 1: Uncharacterized protein lp\_2664

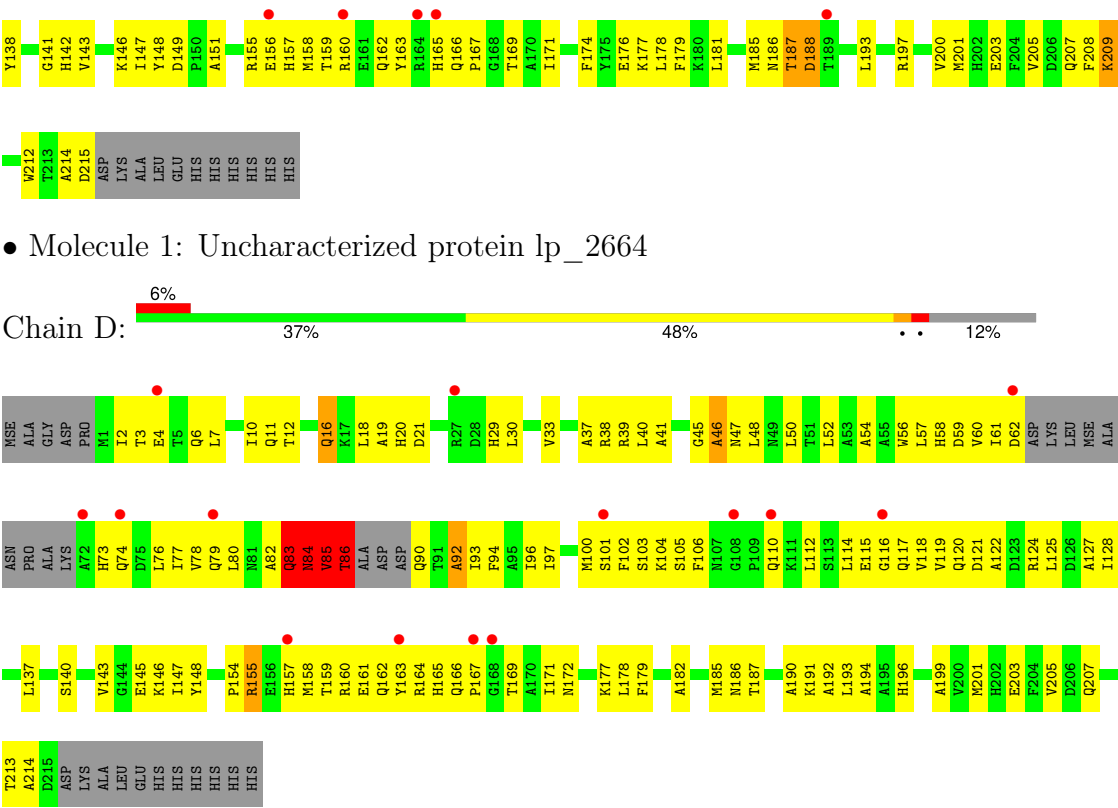


#### • Molecule 1: Uncharacterized protein lp\_2664



#### • Molecule 1: Uncharacterized protein lp\_2664





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.93Å 167.44Å 63.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 20.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	87.1 (20.00-2.80) 93.0 (20.00-2.80)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.85 (at 2.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.236 , 0.307 0.242 , 0.312	Depositor DCC
$R_{free}$ test set	2125 reflections (9.84%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.7	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 29.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	6575	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.43	0/1653	0.69	0/2232
1	B	0.50	0/1695	0.74	6/2287 (0.3%)
1	C	0.70	5/1696 (0.3%)	0.78	5/2290 (0.2%)
1	D	0.62	5/1631 (0.3%)	0.86	11/2200 (0.5%)
All	All	0.57	10/6675 (0.1%)	0.77	22/9009 (0.2%)

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	D	0	4

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	60	VAL	CA-CB	-12.41	1.28	1.54
1	C	60	VAL	CB-CG1	-10.32	1.31	1.52
1	D	84	ASN	CA-C	-9.27	1.28	1.52
1	C	60	VAL	CB-CG2	-8.10	1.35	1.52
1	D	82	ALA	CA-CB	-7.50	1.36	1.52

The worst 5 of 22 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	83	GLN	CA-C-N	-9.64	95.99	117.20
1	C	61	ILE	CA-C-N	8.79	136.54	117.20
1	C	66	MSE	CG-SE-CE	-8.73	79.68	98.90
1	D	84	ASN	CA-C-N	-8.42	98.67	117.20
1	D	85	VAL	CA-CB-CG1	-8.40	98.31	110.90

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	83	GLN	Peptide,Mainchain
1	D	84	ASN	Mainchain
1	D	85	VAL	Mainchain

## 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	1625	0	1590	90	0
1	B	1667	0	1642	98	0
1	C	1667	0	1635	136	0
1	D	1604	0	1576	121	0
2	A	2	0	0	1	0
2	B	6	0	0	1	0
2	C	3	0	0	0	0
2	D	1	0	0	0	0
All	All	6575	0	6443	427	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 33.

The worst 5 of 427 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:C:60:VAL:O	1:C:61:ILE:HG22	1.45	1.12
1:B:18:LEU:HD12	1:B:30:LEU:HD11	1.42	0.99
1:D:38:ARG:HH12	1:D:48:LEU:HD21	1.27	0.99
1:C:187:THR:HG22	1:C:188:ASP:H	1.26	0.98
1:C:174:PHE:HA	1:C:178:LEU:HB2	1.44	0.96

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	A	202/231 (87%)	171 (85%)	23 (11%)	8 (4%)	2	8
1	B	208/231 (90%)	179 (86%)	24 (12%)	5 (2%)	5	18
1	C	208/231 (90%)	172 (83%)	24 (12%)	12 (6%)	1	4
1	D	197/231 (85%)	163 (83%)	32 (16%)	2 (1%)	13	39
All	All	815/924 (88%)	685 (84%)	103 (13%)	27 (3%)	3	11

5 of 27 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	85	VAL
1	A	97	ILE
1	A	113	SER
1	C	82	ALA
1	C	187	THR

### 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	165/178 (93%)	151 (92%)	14 (8%)	8	27
1	B	169/178 (95%)	156 (92%)	13 (8%)	10	31
1	C	169/178 (95%)	156 (92%)	13 (8%)	10	31
1	D	163/178 (92%)	156 (96%)	7 (4%)	25	57
All	All	666/712 (94%)	619 (93%)	47 (7%)	12	35

5 of 47 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	27	ARG
1	C	113	SER
1	C	50	LEU
1	C	61	ILE
1	C	157	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 36 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	186	ASN
1	D	162	GLN
1	D	47	ASN
1	D	99	HIS
1	B	74	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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	201/231 (87%)	-0.10	5 (2%) 58 49	2, 17, 44, 63	0
1	B	206/231 (89%)	0.03	10 (4%) 36 28	1, 18, 82, 97	0
1	C	206/231 (89%)	0.51	19 (9%) 16 12	3, 36, 77, 101	0
1	D	198/231 (85%)	0.58	14 (7%) 23 18	3, 38, 83, 89	0
All	All	811/924 (87%)	0.25	48 (5%) 29 22	1, 26, 78, 101	0

The worst 5 of 48 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	215	ASP	8.3
1	A	214	ALA	6.9
1	A	213	THR	5.4
1	C	110	GLN	4.6
1	D	167	PRO	4.1

### 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 monosaccharides 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.