



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:36 AM UTC

PDB ID : 2QWW / pdb_00002qww
Title : Crystal structure of multiple antibiotic-resistance repressor (MarR) (YP_013417.1) from *Listeria monocytogenes* 4b F2365 at 2.07 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2007-08-10
Resolution : 2.07 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

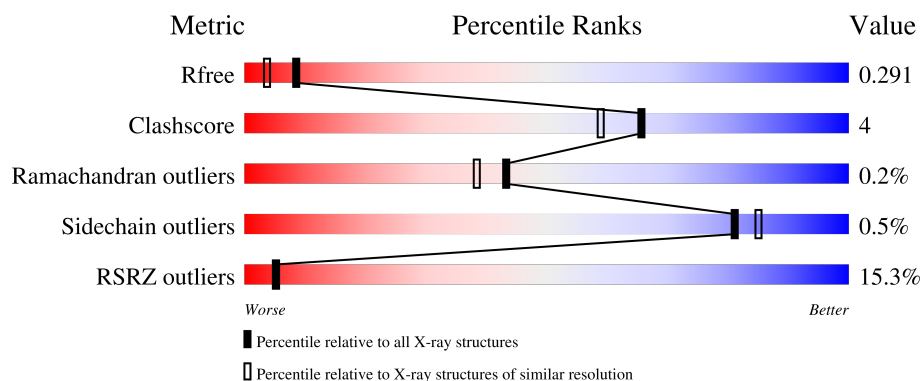
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (2.08-2.04)

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	154	
1	B	154	
1	C	154	
1	D	154	
1	E	154	

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Mol	Chain	Length	Quality of chain
1	F	154	12% 79% 10% • 10%
1	G	154	13% 82% 8% • 9%
1	H	154	16% 86% 7% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8738 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 Transcriptional regulator, MarR family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	Se	0	2	0
			1094	696	181	211	6			
1	B	144	Total	C	N	O	Se	0	2	0
			1075	691	174	205	5			
1	C	138	Total	C	N	O	Se	0	4	0
			1041	665	169	203	4			
1	D	139	Total	C	N	O	Se	0	0	0
			1032	658	169	201	4			
1	E	143	Total	C	N	O	Se	0	1	0
			1047	667	173	203	4			
1	F	139	Total	C	N	O	Se	0	0	0
			1032	657	169	201	5			
1	G	140	Total	C	N	O	Se	0	0	0
			1036	662	174	195	5			
1	H	143	Total	C	N	O	Se	0	3	0
			1048	669	173	202	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q722C0
B	0	GLY	-	expression tag	UNP Q722C0
C	0	GLY	-	expression tag	UNP Q722C0
D	0	GLY	-	expression tag	UNP Q722C0
E	0	GLY	-	expression tag	UNP Q722C0
F	0	GLY	-	expression tag	UNP Q722C0
G	0	GLY	-	expression tag	UNP Q722C0
H	0	GLY	-	expression tag	UNP Q722C0

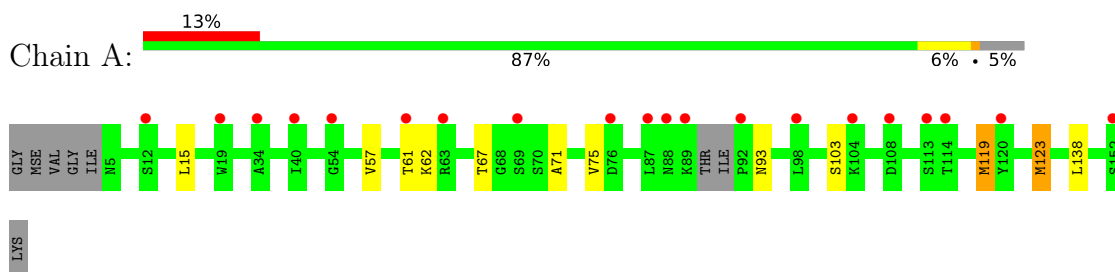
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	49	Total 49	O 49	0	0
2	B	45	Total 45	O 45	0	0
2	C	50	Total 50	O 50	0	0
2	D	37	Total 37	O 37	0	0
2	E	41	Total 41	O 41	0	0
2	F	40	Total 40	O 40	0	0
2	G	36	Total 36	O 36	0	0
2	H	35	Total 35	O 35	0	0

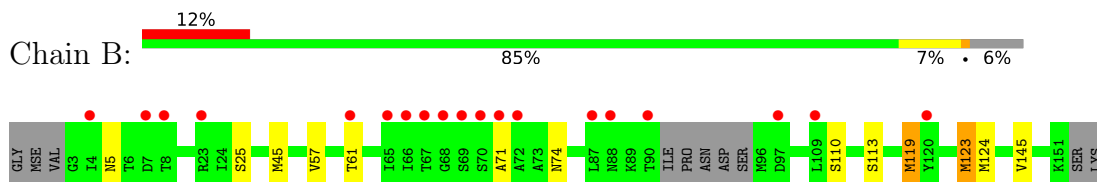
3 Residue-property plots [i](#)

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.

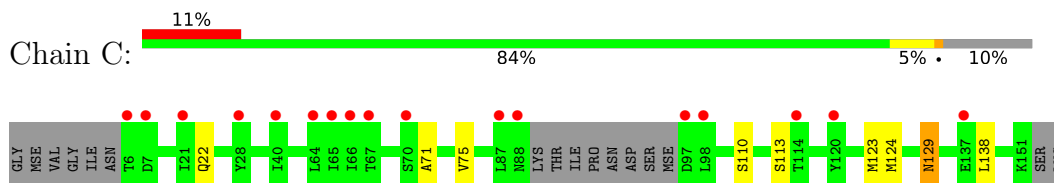
- Molecule 1: Transcriptional regulator, MarR family



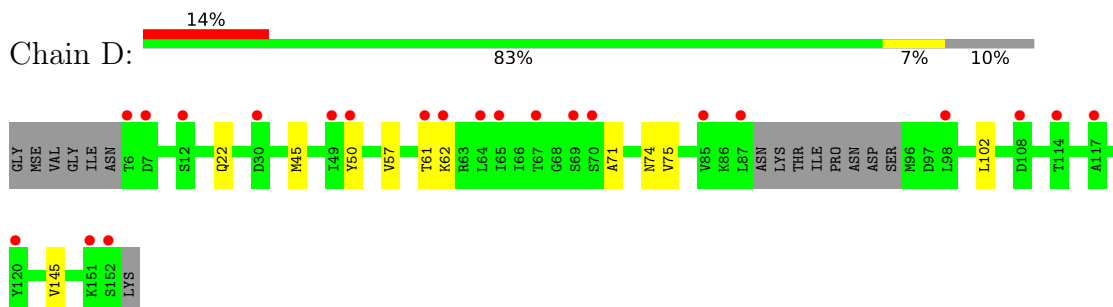
- Molecule 1: Transcriptional regulator, MarR family



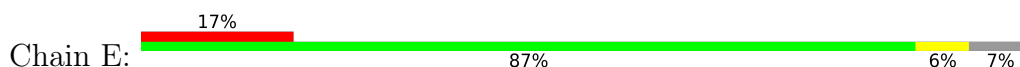
- Molecule 1: Transcriptional regulator, MarR family

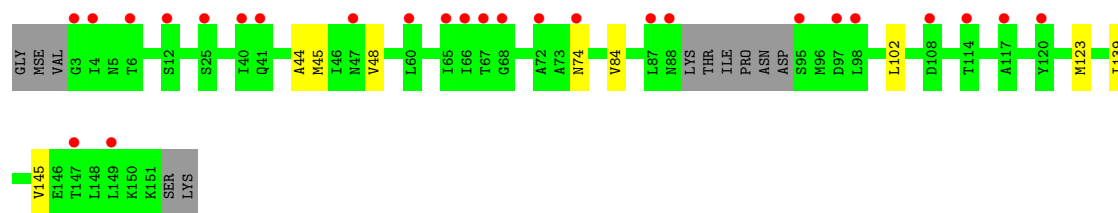


- Molecule 1: Transcriptional regulator, MarR family

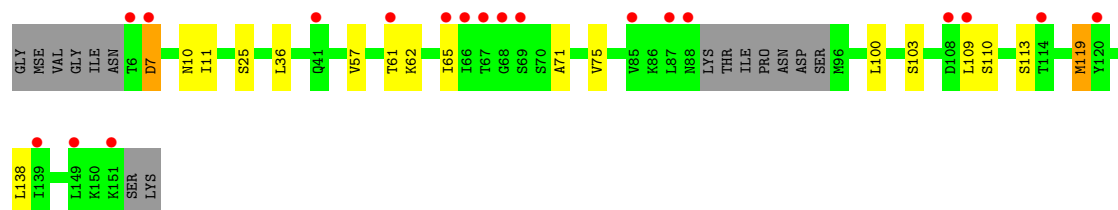
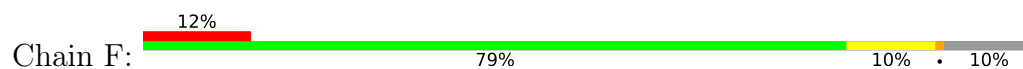


- Molecule 1: Transcriptional regulator, MarR family

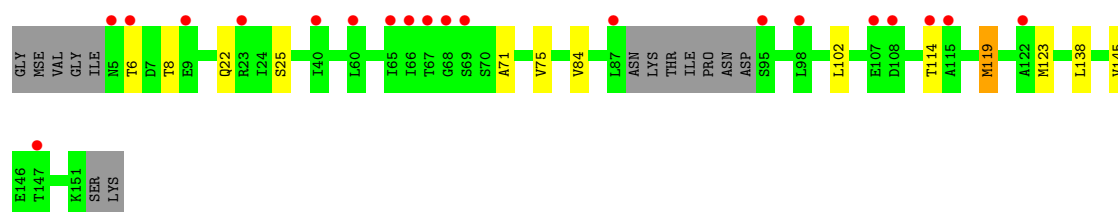
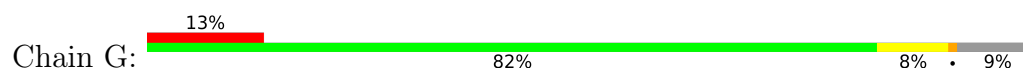




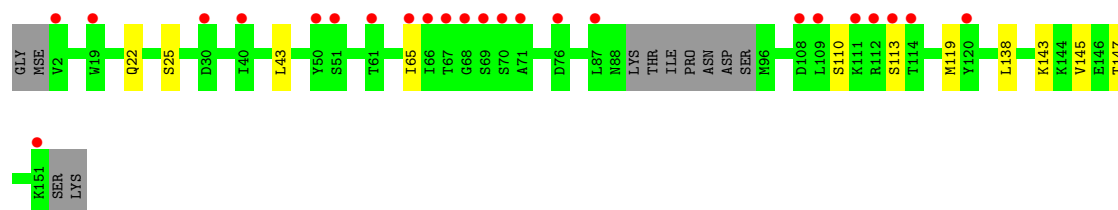
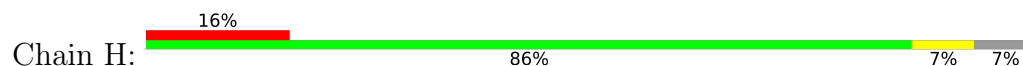
- Molecule 1: Transcriptional regulator, MarR family



- Molecule 1: Transcriptional regulator, MarR family



- Molecule 1: Transcriptional regulator, MarR family



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.25Å 159.70Å 90.27Å 90.00° 103.26° 90.00°	Depositor
Resolution (Å)	46.98 – 2.07 46.98 – 2.07	Depositor EDS
% Data completeness (in resolution range)	98.1 (46.98-2.07) 98.1 (46.98-2.07)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 2.07Å)	Xtriage
Refinement program	REFMAC 5.2.0005, PHENIX	Depositor
R, R_{free}	0.239 , 0.293 0.242 , 0.291	Depositor DCC
R_{free} test set	3967 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	26.4	Xtriage
Anisotropy	0.472	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.037 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8738	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.51 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.2438e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.77	2/1103 (0.2%)	0.96	2/1482 (0.1%)
1	B	0.81	2/1081 (0.2%)	0.94	0/1455
1	C	0.72	1/1056 (0.1%)	0.89	1/1424 (0.1%)
1	D	0.71	0/1035	0.88	0/1393
1	E	0.76	1/1052 (0.1%)	0.94	0/1418
1	F	0.77	1/1035 (0.1%)	0.99	3/1393 (0.2%)
1	G	0.76	2/1038 (0.2%)	0.88	1/1394 (0.1%)
1	H	0.71	0/1060	0.92	0/1430
All	All	0.75	9/8460 (0.1%)	0.92	7/11389 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	123	MSE	SE-CE	-10.92	1.62	1.95
1	B	123	MSE	SE-CE	-7.84	1.72	1.95
1	G	123	MSE	SE-CE	-7.56	1.72	1.95
1	C	123	MSE	SE-CE	-7.33	1.73	1.95
1	G	119	MSE	SE-CE	-6.33	1.76	1.95
1	E	123	MSE	SE-CE	-6.08	1.77	1.95
1	F	119	MSE	SE-CE	-5.73	1.78	1.95
1	A	119	MSE	SE-CE	-5.60	1.78	1.95
1	B	119	MSE	SE-CE	-5.47	1.79	1.95

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	119	MSE	CG-SE-CE	-5.75	86.28	98.92
1	F	103	SER	N-CA-C	-5.74	102.27	110.59
1	C	124	MSE	N-CA-C	-5.44	105.43	111.36
1	A	103	SER	N-CA-C	-5.41	102.88	110.35
1	F	100	LEU	N-CA-C	5.27	118.02	109.06
1	F	7	ASP	N-CA-C	-5.12	105.59	111.07
1	A	93	ASN	N-CA-C	5.09	117.57	111.71

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	1094	0	1120	8	0
1	B	1075	0	1102	12	0
1	C	1041	0	1067	5	0
1	D	1032	0	1052	9	0
1	E	1047	0	1045	5	0
1	F	1032	0	1053	14	0
1	G	1036	0	1072	11	0
1	H	1048	0	1056	9	0
2	A	49	0	0	1	0
2	B	45	0	0	0	0
2	C	50	0	0	1	0
2	D	37	0	0	0	0
2	E	41	0	0	0	0
2	F	40	0	0	0	0
2	G	36	0	0	1	0
2	H	35	0	0	0	0
All	All	8738	0	8567	64	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 4.

All (64) 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:G:25:SER:HA	1:G:119:MSE:HE1	1.61	0.81
1:B:61:THR:HG22	1:B:71:ALA:CB	2.19	0.72
1:D:61:THR:HG22	1:D:71:ALA:CB	2.22	0.70
1:F:61:THR:HG22	1:F:71:ALA:CB	2.23	0.69
1:G:25:SER:CA	1:G:119:MSE:HE1	2.23	0.67
1:F:25:SER:CB	1:F:119:MSE:HE1	2.23	0.67
1:A:61:THR:HG22	1:A:71:ALA:CB	2.25	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:36:LEU:HB3	1:F:109:LEU:HD21	1.79	0.65
1:H:43:LEU:HD21	1:H:113[A]:SER:OG	1.98	0.64
1:D:71:ALA:O	1:D:75:VAL:HG23	2.01	0.61
1:G:25:SER:CB	1:G:119:MSE:HE1	2.32	0.59
1:H:110:SER:O	1:H:113[B]:SER:OG	2.20	0.58
1:A:119:MSE:HE3	1:A:123:MSE:HE2	1.86	0.58
1:F:25:SER:HB3	1:F:119:MSE:HE1	1.85	0.58
1:F:57:VAL:O	1:F:61:THR:HG23	2.04	0.57
1:B:25:SER:HB3	1:B:119:MSE:HE1	1.87	0.55
1:C:110:SER:O	1:C:113[B]:SER:OG	2.26	0.53
1:A:57:VAL:O	1:A:61:THR:HG23	2.08	0.53
1:B:25:SER:CB	1:B:119:MSE:HE1	2.39	0.52
1:B:57:VAL:O	1:B:61:THR:HG23	2.11	0.51
1:C:138:LEU:HD21	1:D:145:VAL:HG21	1.93	0.51
1:D:57:VAL:O	1:D:61:THR:HG23	2.10	0.51
2:A:167:HOH:O	1:B:124:MSE:HG3	2.09	0.51
1:D:45:MSE:HE1	1:D:74:ASN:HB2	1.93	0.50
1:E:84:VAL:HG12	1:E:102:LEU:HD23	1.92	0.50
1:H:25:SER:CB	1:H:119:MSE:HE1	2.42	0.49
1:C:22[B]:GLN:OE1	1:D:22:GLN:NE2	2.46	0.48
1:H:143:LYS:O	1:H:147:THR:HG23	2.13	0.48
1:B:25:SER:CA	1:B:119:MSE:HE1	2.44	0.48
1:C:71:ALA:O	1:C:75:VAL:HG23	2.13	0.48
1:H:25:SER:CA	1:H:119:MSE:HE1	2.44	0.48
1:C:129:ASN:N	1:C:129:ASN:HD22	2.11	0.48
1:A:61:THR:OG1	1:A:62:LYS:N	2.47	0.47
1:G:145:VAL:HG21	1:H:138:LEU:HD21	1.96	0.47
2:C:169:HOH:O	1:G:6:THR:HG21	2.16	0.46
1:G:138:LEU:HD21	1:H:145:VAL:HG21	1.96	0.46
1:E:139:ILE:HD11	1:F:10:ASN:HB3	1.98	0.46
1:F:25:SER:CA	1:F:119:MSE:HE1	2.46	0.45
1:B:61:THR:HG22	1:B:71:ALA:HB3	1.97	0.45
1:G:114:THR:OG1	2:G:171:HOH:O	2.20	0.44
1:B:25:SER:HA	1:B:119:MSE:HE1	1.99	0.44
1:F:71:ALA:O	1:F:75:VAL:HG23	2.17	0.43
1:E:145:VAL:HG21	1:F:138:LEU:HD21	1.99	0.43
1:A:119:MSE:HE3	1:A:123:MSE:CE	2.48	0.43
1:F:61:THR:OG1	1:F:62:LYS:N	2.51	0.43
1:H:25:SER:HA	1:H:119:MSE:HE1	2.00	0.43
1:E:44:ALA:O	1:E:48:VAL:HG23	2.19	0.43
1:G:6:THR:HG22	1:G:8:THR:H	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:45:MSE:HE1	1:D:74:ASN:CB	2.48	0.43
1:A:15:LEU:HD21	1:B:123:MSE:HE3	2.01	0.43
1:E:45:MSE:HE1	1:E:74:ASN:HB2	2.01	0.42
1:G:22:GLN:HG2	1:H:22:GLN:HG2	2.00	0.42
1:G:84:VAL:HG12	1:G:102:LEU:HD23	2.00	0.42
1:A:138:LEU:HD21	1:B:145:VAL:HG21	2.01	0.42
1:B:110:SER:O	1:B:113:SER:OG	2.32	0.42
1:B:45:MSE:HE1	1:B:74:ASN:HB2	2.02	0.41
1:G:71:ALA:O	1:G:75:VAL:HG23	2.21	0.41
1:F:61:THR:HG22	1:F:71:ALA:HB3	1.97	0.41
1:A:71:ALA:O	1:A:75:VAL:HG23	2.21	0.41
1:D:50:TYR:HB2	1:D:102:LEU:HD11	2.02	0.41
1:D:61:THR:HG22	1:D:71:ALA:HB3	2.01	0.40
1:F:7:ASP:O	1:F:11:ILE:HG13	2.21	0.40
1:F:25:SER:HA	1:F:119:MSE:HE1	2.03	0.40
1:F:110:SER:O	1:F:113:SER:OG	2.24	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	144/154 (94%)	142 (99%)	2 (1%)	0	100	100
1	B	142/154 (92%)	140 (99%)	2 (1%)	0	100	100
1	C	138/154 (90%)	136 (99%)	2 (1%)	0	100	100
1	D	135/154 (88%)	134 (99%)	1 (1%)	0	100	100
1	E	140/154 (91%)	138 (99%)	2 (1%)	0	100	100
1	F	135/154 (88%)	132 (98%)	2 (2%)	1 (1%)	18	10
1	G	136/154 (88%)	135 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	142/154 (92%)	141 (99%)	0	1 (1%)	18	10
All	All	1112/1232 (90%)	1098 (99%)	12 (1%)	2 (0%)	43	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	65	ILE
1	F	65	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	116/127 (91%)	115 (99%)	1 (1%)	70	74
1	B	110/127 (87%)	109 (99%)	1 (1%)	70	74
1	C	111/127 (87%)	110 (99%)	1 (1%)	70	74
1	D	108/127 (85%)	107 (99%)	1 (1%)	70	74
1	E	105/127 (83%)	105 (100%)	0	100	100
1	F	109/127 (86%)	109 (100%)	0	100	100
1	G	108/127 (85%)	108 (100%)	0	100	100
1	H	107/127 (84%)	107 (100%)	0	100	100
All	All	874/1016 (86%)	870 (100%)	4 (0%)	81	85

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	THR
1	B	5	ASN
1	C	129	ASN
1	D	62	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such

sidechains are listed below:

Mol	Chain	Res	Type
1	B	31	GLN
1	C	129	ASN
1	D	22	GLN
1	E	22	GLN
1	F	10	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 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 [i](#)

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	141/154 (91%)	1.22	20 (14%) 6 6	20, 41, 56, 71	1 (0%)
1	B	139/154 (90%)	1.17	19 (13%) 6 6	17, 40, 55, 64	2 (1%)
1	C	134/154 (87%)	1.16	17 (12%) 8 7	21, 40, 54, 57	4 (2%)
1	D	134/154 (87%)	1.20	22 (16%) 4 4	31, 41, 54, 57	0
1	E	138/154 (89%)	1.27	26 (18%) 3 3	27, 40, 54, 61	1 (0%)
1	F	134/154 (87%)	1.25	19 (14%) 6 6	29, 40, 54, 57	0
1	G	135/154 (87%)	1.16	20 (14%) 5 5	30, 40, 53, 57	0
1	H	138/154 (89%)	1.33	24 (17%) 4 3	20, 40, 53, 59	3 (2%)
All	All	1093/1232 (88%)	1.22	167 (15%) 5 5	17, 40, 54, 71	11 (1%)

All (167) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	88	ASN	7.4
1	H	68	GLY	5.9
1	D	87	LEU	5.3
1	B	87	LEU	5.2
1	A	92	PRO	5.0
1	F	65	ILE	4.9
1	C	7	ASP	4.9
1	C	88	ASN	4.8
1	B	68	GLY	4.8
1	F	88	ASN	4.8
1	C	65	ILE	4.5
1	F	7	ASP	4.5
1	F	67	THR	4.4
1	D	6	THR	4.3
1	A	152	SER	4.3
1	H	65	ILE	4.3

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Mol	Chain	Res	Type	RSRZ
1	H	151	LYS	4.3
1	A	87	LEU	4.2
1	A	108	ASP	4.1
1	D	108	ASP	4.1
1	E	98	LEU	4.1
1	F	120	TYR	4.1
1	H	108	ASP	4.1
1	B	67	THR	4.1
1	F	151	LYS	4.0
1	F	66	ILE	4.0
1	C	120	TYR	3.9
1	H	87	LEU	3.9
1	G	108	ASP	3.9
1	H	69	SER	3.9
1	C	67	THR	3.9
1	G	114	THR	3.8
1	E	41	GLN	3.8
1	G	67	THR	3.8
1	B	70	SER	3.7
1	G	95	SER	3.7
1	E	4	ILE	3.7
1	G	115	ALA	3.7
1	E	108	ASP	3.6
1	B	120	TYR	3.6
1	G	87	LEU	3.6
1	A	114	THR	3.5
1	A	89	LYS	3.5
1	D	114	THR	3.5
1	H	109	LEU	3.5
1	B	90	THR	3.5
1	F	6	THR	3.5
1	A	88	ASN	3.3
1	D	65	ILE	3.3
1	F	68	GLY	3.3
1	G	6	THR	3.3
1	F	69	SER	3.3
1	C	114	THR	3.2
1	E	95	SER	3.2
1	A	61	THR	3.1
1	H	71	ALA	3.1
1	A	12	SER	3.1
1	G	122	ALA	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	87	LEU	3.1
1	F	61	THR	3.1
1	F	87	LEU	3.1
1	B	61	THR	3.0
1	G	98	LEU	3.0
1	B	7	ASP	3.0
1	E	120	TYR	2.9
1	D	12	SER	2.9
1	D	152	SER	2.9
1	D	117	ALA	2.9
1	C	98	LEU	2.8
1	H	67	THR	2.8
1	B	69	SER	2.8
1	H	70	SER	2.8
1	H	76	ASP	2.8
1	D	50	TYR	2.8
1	H	50	TYR	2.8
1	C	6	THR	2.8
1	D	61	THR	2.8
1	E	67	THR	2.8
1	G	5	ASN	2.8
1	E	3	GLY	2.7
1	B	72	ALA	2.7
1	D	67	THR	2.7
1	H	61	THR	2.7
1	H	2	VAL	2.7
1	G	68	GLY	2.7
1	D	30	ASP	2.7
1	G	40	ILE	2.7
1	E	87	LEU	2.6
1	B	8	THR	2.6
1	B	109	LEU	2.6
1	A	40	ILE	2.6
1	B	71	ALA	2.6
1	B	88	ASN	2.5
1	E	72	ALA	2.5
1	A	19	TRP	2.5
1	H	111	LYS	2.5
1	H	120	TYR	2.5
1	C	70	SER	2.5
1	D	62	LYS	2.5
1	E	12	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	104	LYS	2.5
1	C	137	GLU	2.4
1	D	7	ASP	2.4
1	H	30	ASP	2.4
1	E	117	ALA	2.4
1	F	149	LEU	2.4
1	E	6	THR	2.4
1	D	151	LYS	2.4
1	E	68	GLY	2.4
1	G	65	ILE	2.4
1	C	28	TYR	2.3
1	A	34	ALA	2.3
1	F	108	ASP	2.3
1	H	66	ILE	2.3
1	G	147	THR	2.3
1	G	107	GLU	2.3
1	E	47	ASN	2.3
1	G	23	ARG	2.3
1	A	113	SER	2.3
1	D	70	SER	2.3
1	A	76	ASP	2.3
1	A	69	SER	2.3
1	C	66	ILE	2.3
1	B	97	ASP	2.3
1	E	65	ILE	2.2
1	D	85	VAL	2.2
1	D	98	LEU	2.2
1	F	109	LEU	2.2
1	G	60	LEU	2.2
1	F	41	GLN	2.2
1	D	69	SER	2.2
1	H	113[A]	SER	2.2
1	F	85	VAL	2.2
1	A	54	GLY	2.2
1	G	69	SER	2.2
1	H	112	ARG	2.2
1	B	65	ILE	2.2
1	E	114	THR	2.2
1	E	147	THR	2.2
1	G	66	ILE	2.2
1	H	51	SER	2.2
1	F	114	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	63	ARG	2.1
1	A	98	LEU	2.1
1	D	64	LEU	2.1
1	E	40	ILE	2.1
1	E	66	ILE	2.1
1	F	139	ILE	2.1
1	H	40	ILE	2.1
1	E	60	LEU	2.1
1	C	97	ASP	2.1
1	A	120	TYR	2.1
1	D	120	TYR	2.1
1	H	114	THR	2.1
1	B	66	ILE	2.1
1	H	19	TRP	2.1
1	E	149	LEU	2.1
1	B	4	ILE	2.1
1	C	40	ILE	2.1
1	C	21	ILE	2.0
1	D	49	ILE	2.0
1	G	9	GLU	2.0
1	E	25	SER	2.0
1	E	74	ASN	2.0
1	C	64	LEU	2.0
1	E	97	ASP	2.0
1	B	23	ARG	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.