Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 03:55 am BST

PDB ID	:	5LFM
Title	:	MamA RS-1 ArsTM double mutant
Authors	:	Zarivach, R.; Cronin, S.L.
Deposited on	:	2016-07-03
Resolution	:	3.28 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 following versions of software and data (see references ⓘ) were used in the production of this report:

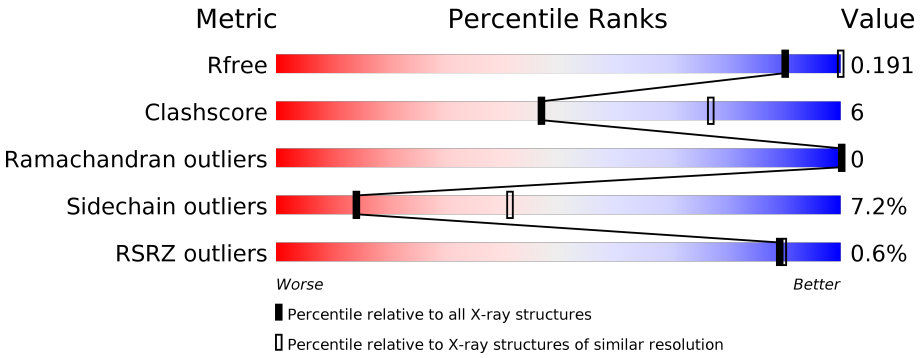
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 3.28 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	1177 (3.32-3.24)
Clashscore	141614	1044 (3.30-3.26)
Ramachandran outliers	138981	1026 (3.30-3.26)
Sidechain outliers	138945	1025 (3.30-3.26)
RSRZ outliers	127900	1141 (3.32-3.24)

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 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% 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	202	74% 11% • 12%
1	B	202	74% 11% • 13%
1	C	202	74% 11% • 13%
1	D	202	77% 8% • 13%
1	E	202	% 72% 13% • 12%
1	F	202	% 77% 7% • 12%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	C	301	-	-	-	X
2	ACT	E	301	-	-	-	X
2	ACT	F	301	-	-	-	X
3	CL	E	302	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8563 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 Magnetosome protein MamA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	177	Total	C	N	O	S	0	0	0
			1423	905	255	256	7			
1	B	176	Total	C	N	O	S	0	0	0
			1417	902	254	254	7			
1	C	176	Total	C	N	O	S	0	0	0
			1417	902	254	254	7			
1	D	176	Total	C	N	O	S	0	0	0
			1417	902	254	254	7			
1	E	177	Total	C	N	O	S	0	0	0
			1423	905	255	256	7			
1	F	178	Total	C	N	O	S	0	0	0
			1431	909	256	259	7			

There are 192 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	37	MET	-	initiating methionine	UNP C4XPQ7
A	38	ALA	-	expression tag	UNP C4XPQ7
A	39	MET	-	expression tag	UNP C4XPQ7
A	40	GLY	-	expression tag	UNP C4XPQ7
A	46	PHE	TYR	conflict	UNP C4XPQ7
A	65	PHE	TYR	conflict	UNP C4XPQ7
A	68	LEU	GLU	conflict	UNP C4XPQ7
A	124	ILE	MET	conflict	UNP C4XPQ7
A	140	ALA	GLU	conflict	UNP C4XPQ7
A	141	ALA	LYS	conflict	UNP C4XPQ7
A	143	ALA	GLU	conflict	UNP C4XPQ7
A	218	GLU	-	expression tag	UNP C4XPQ7
A	219	LEU	-	expression tag	UNP C4XPQ7
A	220	ALA	-	expression tag	UNP C4XPQ7
A	221	LEU	-	expression tag	UNP C4XPQ7
A	222	VAL	-	expression tag	UNP C4XPQ7
A	223	PRO	-	expression tag	UNP C4XPQ7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	224	ARG	-	expression tag	UNP C4XPQ7
A	225	GLY	-	expression tag	UNP C4XPQ7
A	226	SER	-	expression tag	UNP C4XPQ7
A	227	SER	-	expression tag	UNP C4XPQ7
A	228	ALA	-	expression tag	UNP C4XPQ7
A	229	HIS	-	expression tag	UNP C4XPQ7
A	230	HIS	-	expression tag	UNP C4XPQ7
A	231	HIS	-	expression tag	UNP C4XPQ7
A	232	HIS	-	expression tag	UNP C4XPQ7
A	233	HIS	-	expression tag	UNP C4XPQ7
A	234	HIS	-	expression tag	UNP C4XPQ7
A	235	HIS	-	expression tag	UNP C4XPQ7
A	236	HIS	-	expression tag	UNP C4XPQ7
A	237	HIS	-	expression tag	UNP C4XPQ7
A	238	HIS	-	expression tag	UNP C4XPQ7
B	37	MET	-	initiating methionine	UNP C4XPQ7
B	38	ALA	-	expression tag	UNP C4XPQ7
B	39	MET	-	expression tag	UNP C4XPQ7
B	40	GLY	-	expression tag	UNP C4XPQ7
B	46	PHE	TYR	conflict	UNP C4XPQ7
B	65	PHE	TYR	conflict	UNP C4XPQ7
B	68	LEU	GLU	conflict	UNP C4XPQ7
B	124	ILE	MET	conflict	UNP C4XPQ7
B	140	ALA	GLU	conflict	UNP C4XPQ7
B	141	ALA	LYS	conflict	UNP C4XPQ7
B	143	ALA	GLU	conflict	UNP C4XPQ7
B	218	GLU	-	expression tag	UNP C4XPQ7
B	219	LEU	-	expression tag	UNP C4XPQ7
B	220	ALA	-	expression tag	UNP C4XPQ7
B	221	LEU	-	expression tag	UNP C4XPQ7
B	222	VAL	-	expression tag	UNP C4XPQ7
B	223	PRO	-	expression tag	UNP C4XPQ7
B	224	ARG	-	expression tag	UNP C4XPQ7
B	225	GLY	-	expression tag	UNP C4XPQ7
B	226	SER	-	expression tag	UNP C4XPQ7
B	227	SER	-	expression tag	UNP C4XPQ7
B	228	ALA	-	expression tag	UNP C4XPQ7
B	229	HIS	-	expression tag	UNP C4XPQ7
B	230	HIS	-	expression tag	UNP C4XPQ7
B	231	HIS	-	expression tag	UNP C4XPQ7
B	232	HIS	-	expression tag	UNP C4XPQ7
B	233	HIS	-	expression tag	UNP C4XPQ7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	234	HIS	-	expression tag	UNP C4XPQ7
B	235	HIS	-	expression tag	UNP C4XPQ7
B	236	HIS	-	expression tag	UNP C4XPQ7
B	237	HIS	-	expression tag	UNP C4XPQ7
B	238	HIS	-	expression tag	UNP C4XPQ7
C	37	MET	-	initiating methionine	UNP C4XPQ7
C	38	ALA	-	expression tag	UNP C4XPQ7
C	39	MET	-	expression tag	UNP C4XPQ7
C	40	GLY	-	expression tag	UNP C4XPQ7
C	46	PHE	TYR	conflict	UNP C4XPQ7
C	65	PHE	TYR	conflict	UNP C4XPQ7
C	68	LEU	GLU	conflict	UNP C4XPQ7
C	124	ILE	MET	conflict	UNP C4XPQ7
C	140	ALA	GLU	conflict	UNP C4XPQ7
C	141	ALA	LYS	conflict	UNP C4XPQ7
C	143	ALA	GLU	conflict	UNP C4XPQ7
C	218	GLU	-	expression tag	UNP C4XPQ7
C	219	LEU	-	expression tag	UNP C4XPQ7
C	220	ALA	-	expression tag	UNP C4XPQ7
C	221	LEU	-	expression tag	UNP C4XPQ7
C	222	VAL	-	expression tag	UNP C4XPQ7
C	223	PRO	-	expression tag	UNP C4XPQ7
C	224	ARG	-	expression tag	UNP C4XPQ7
C	225	GLY	-	expression tag	UNP C4XPQ7
C	226	SER	-	expression tag	UNP C4XPQ7
C	227	SER	-	expression tag	UNP C4XPQ7
C	228	ALA	-	expression tag	UNP C4XPQ7
C	229	HIS	-	expression tag	UNP C4XPQ7
C	230	HIS	-	expression tag	UNP C4XPQ7
C	231	HIS	-	expression tag	UNP C4XPQ7
C	232	HIS	-	expression tag	UNP C4XPQ7
C	233	HIS	-	expression tag	UNP C4XPQ7
C	234	HIS	-	expression tag	UNP C4XPQ7
C	235	HIS	-	expression tag	UNP C4XPQ7
C	236	HIS	-	expression tag	UNP C4XPQ7
C	237	HIS	-	expression tag	UNP C4XPQ7
C	238	HIS	-	expression tag	UNP C4XPQ7
D	37	MET	-	initiating methionine	UNP C4XPQ7
D	38	ALA	-	expression tag	UNP C4XPQ7
D	39	MET	-	expression tag	UNP C4XPQ7
D	40	GLY	-	expression tag	UNP C4XPQ7
D	46	PHE	TYR	conflict	UNP C4XPQ7

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Chain	Residue	Modelled	Actual	Comment	Reference
D	65	PHE	TYR	conflict	UNP C4XPQ7
D	68	LEU	GLU	conflict	UNP C4XPQ7
D	124	ILE	MET	conflict	UNP C4XPQ7
D	140	ALA	GLU	conflict	UNP C4XPQ7
D	141	ALA	LYS	conflict	UNP C4XPQ7
D	143	ALA	GLU	conflict	UNP C4XPQ7
D	218	GLU	-	expression tag	UNP C4XPQ7
D	219	LEU	-	expression tag	UNP C4XPQ7
D	220	ALA	-	expression tag	UNP C4XPQ7
D	221	LEU	-	expression tag	UNP C4XPQ7
D	222	VAL	-	expression tag	UNP C4XPQ7
D	223	PRO	-	expression tag	UNP C4XPQ7
D	224	ARG	-	expression tag	UNP C4XPQ7
D	225	GLY	-	expression tag	UNP C4XPQ7
D	226	SER	-	expression tag	UNP C4XPQ7
D	227	SER	-	expression tag	UNP C4XPQ7
D	228	ALA	-	expression tag	UNP C4XPQ7
D	229	HIS	-	expression tag	UNP C4XPQ7
D	230	HIS	-	expression tag	UNP C4XPQ7
D	231	HIS	-	expression tag	UNP C4XPQ7
D	232	HIS	-	expression tag	UNP C4XPQ7
D	233	HIS	-	expression tag	UNP C4XPQ7
D	234	HIS	-	expression tag	UNP C4XPQ7
D	235	HIS	-	expression tag	UNP C4XPQ7
D	236	HIS	-	expression tag	UNP C4XPQ7
D	237	HIS	-	expression tag	UNP C4XPQ7
D	238	HIS	-	expression tag	UNP C4XPQ7
E	37	MET	-	initiating methionine	UNP C4XPQ7
E	38	ALA	-	expression tag	UNP C4XPQ7
E	39	MET	-	expression tag	UNP C4XPQ7
E	40	GLY	-	expression tag	UNP C4XPQ7
E	46	PHE	TYR	conflict	UNP C4XPQ7
E	65	PHE	TYR	conflict	UNP C4XPQ7
E	68	LEU	GLU	conflict	UNP C4XPQ7
E	124	ILE	MET	conflict	UNP C4XPQ7
E	140	ALA	GLU	conflict	UNP C4XPQ7
E	141	ALA	LYS	conflict	UNP C4XPQ7
E	143	ALA	GLU	conflict	UNP C4XPQ7
E	218	GLU	-	expression tag	UNP C4XPQ7
E	219	LEU	-	expression tag	UNP C4XPQ7
E	220	ALA	-	expression tag	UNP C4XPQ7
E	221	LEU	-	expression tag	UNP C4XPQ7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	222	VAL	-	expression tag	UNP C4XPQ7
E	223	PRO	-	expression tag	UNP C4XPQ7
E	224	ARG	-	expression tag	UNP C4XPQ7
E	225	GLY	-	expression tag	UNP C4XPQ7
E	226	SER	-	expression tag	UNP C4XPQ7
E	227	SER	-	expression tag	UNP C4XPQ7
E	228	ALA	-	expression tag	UNP C4XPQ7
E	229	HIS	-	expression tag	UNP C4XPQ7
E	230	HIS	-	expression tag	UNP C4XPQ7
E	231	HIS	-	expression tag	UNP C4XPQ7
E	232	HIS	-	expression tag	UNP C4XPQ7
E	233	HIS	-	expression tag	UNP C4XPQ7
E	234	HIS	-	expression tag	UNP C4XPQ7
E	235	HIS	-	expression tag	UNP C4XPQ7
E	236	HIS	-	expression tag	UNP C4XPQ7
E	237	HIS	-	expression tag	UNP C4XPQ7
E	238	HIS	-	expression tag	UNP C4XPQ7
F	37	MET	-	initiating methionine	UNP C4XPQ7
F	38	ALA	-	expression tag	UNP C4XPQ7
F	39	MET	-	expression tag	UNP C4XPQ7
F	40	GLY	-	expression tag	UNP C4XPQ7
F	46	PHE	TYR	conflict	UNP C4XPQ7
F	65	PHE	TYR	conflict	UNP C4XPQ7
F	68	LEU	GLU	conflict	UNP C4XPQ7
F	124	ILE	MET	conflict	UNP C4XPQ7
F	140	ALA	GLU	conflict	UNP C4XPQ7
F	141	ALA	LYS	conflict	UNP C4XPQ7
F	143	ALA	GLU	conflict	UNP C4XPQ7
F	218	GLU	-	expression tag	UNP C4XPQ7
F	219	LEU	-	expression tag	UNP C4XPQ7
F	220	ALA	-	expression tag	UNP C4XPQ7
F	221	LEU	-	expression tag	UNP C4XPQ7
F	222	VAL	-	expression tag	UNP C4XPQ7
F	223	PRO	-	expression tag	UNP C4XPQ7
F	224	ARG	-	expression tag	UNP C4XPQ7
F	225	GLY	-	expression tag	UNP C4XPQ7
F	226	SER	-	expression tag	UNP C4XPQ7
F	227	SER	-	expression tag	UNP C4XPQ7
F	228	ALA	-	expression tag	UNP C4XPQ7
F	229	HIS	-	expression tag	UNP C4XPQ7
F	230	HIS	-	expression tag	UNP C4XPQ7
F	231	HIS	-	expression tag	UNP C4XPQ7

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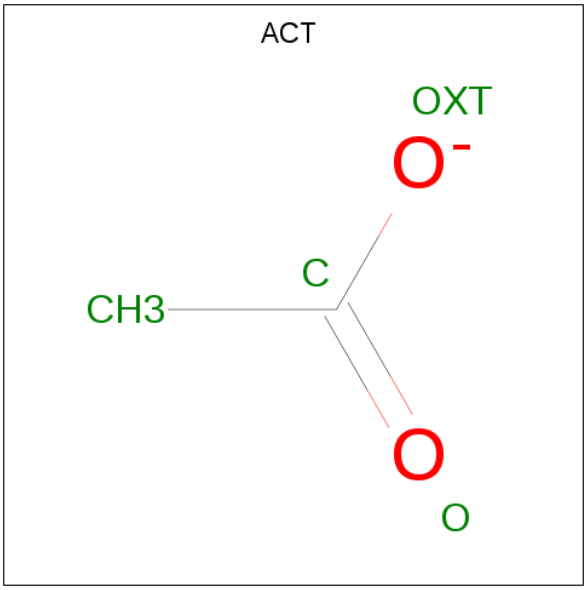
Chain	Residue	Modelled	Actual	Comment	Reference
F	232	HIS	-	expression tag	UNP C4XPQ7
F	233	HIS	-	expression tag	UNP C4XPQ7
F	234	HIS	-	expression tag	UNP C4XPQ7
F	235	HIS	-	expression tag	UNP C4XPQ7
F	236	HIS	-	expression tag	UNP C4XPQ7
F	237	HIS	-	expression tag	UNP C4XPQ7
F	238	HIS	-	expression tag	UNP C4XPQ7

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0
3	A	1	Total Cl 1 1	0	0
3	D	1	Total Cl 1 1	0	0
3	C	2	Total Cl 2 2	0	0
3	E	1	Total Cl 1 1	0	0

- Molecule 4 is water.

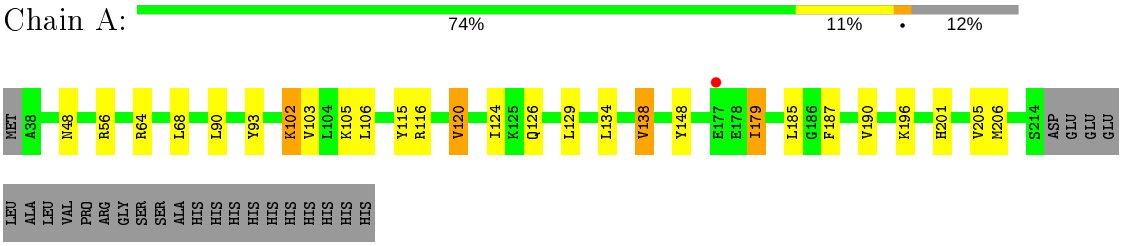
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total O 1 1	0	0
4	C	1	Total O 1 1	0	0
4	D	1	Total O 1 1	0	0
4	E	1	Total O 1 1	0	0
4	F	1	Total O 1 1	0	0

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 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.

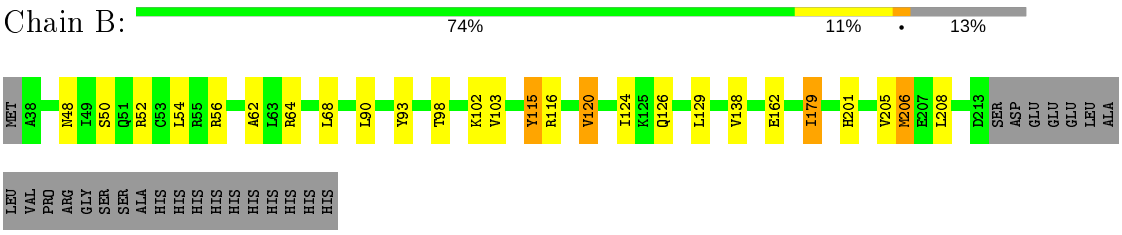
- Molecule 1: Magnetosome protein MamA

Chain A:



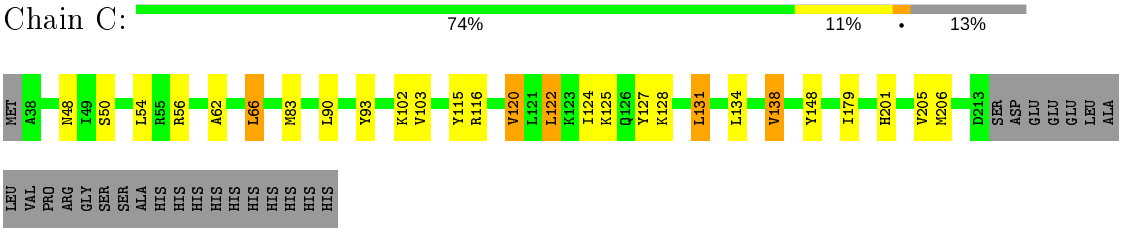
- Molecule 1: Magnetosome protein MamA

Chain B:



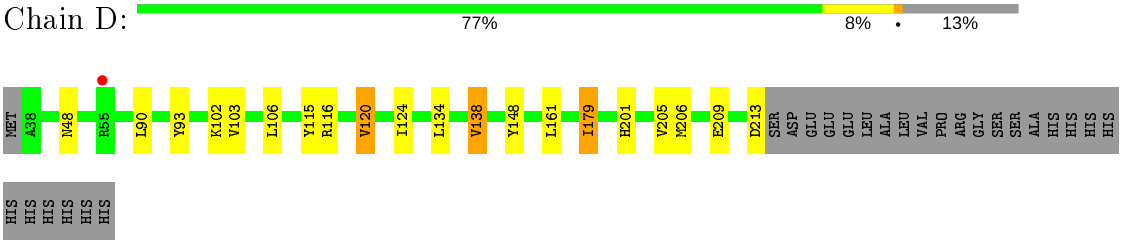
- Molecule 1: Magnetosome protein MamA

Chain C:



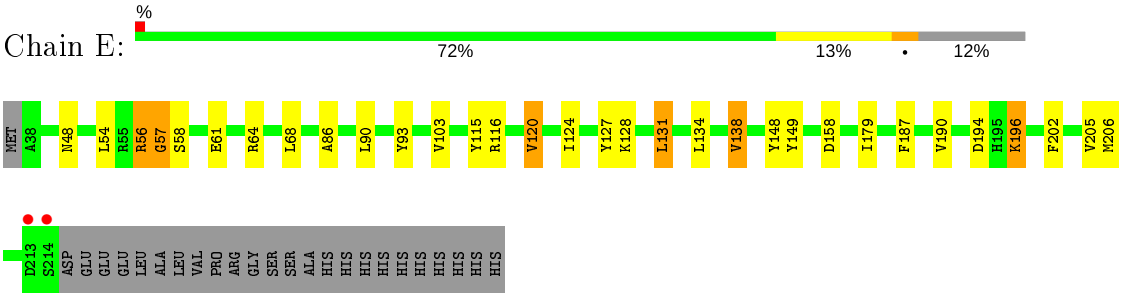
- Molecule 1: Magnetosome protein MamA

Chain D:



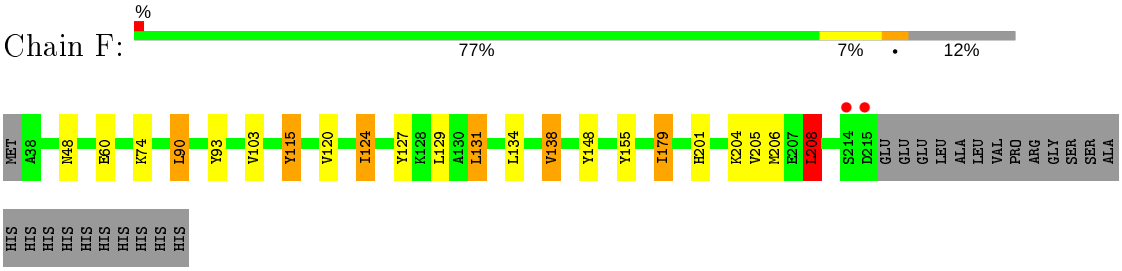
• Molecule 1: Magnetosome protein MamA

Chain E:



• Molecule 1: Magnetosome protein MamA

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α, β, γ	150.12Å 150.12Å 203.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	120.90 – 3.28 48.60 – 3.28	Depositor EDS
% Data completeness (in resolution range)	99.9 (120.90-3.28) 100.0 (48.60-3.28)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.58 (at 3.25Å)	Xtriage
Refinement program	REFMAC 5.8.0151	Depositor
R, R_{free}	0.171 , 0.188 0.175 , 0.191	Depositor DCC
R_{free} test set	1740 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	91.3	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 53.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.008 for -1/2*h+1/2*k-1/2*l,1/2*h-1/2*k-1/2*l,-h-k 0.000 for -1/2*h+1/2*k+1/2*l,1/2*h-1/2*k+1/2*l,h+k 0.000 for -1/2*h-1/2*k+1/2*l,-1/2*h-1/2*k-1/2*l,h-k 0.005 for -1/2*h-1/2*k-1/2*l,-1/2*h-1/2*k+1/2*l,-h+k 0.024 for -h,k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8563	wwPDB-VP
Average B, all atoms (Å^2)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.26% 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

Bond lengths and bond angles in the following residue types are not validated in this section: CL, ACT

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	0.50	0/1451	0.78	1/1951 (0.1%)
1	B	0.50	0/1445	0.80	3/1943 (0.2%)
1	C	0.48	0/1445	0.81	2/1943 (0.1%)
1	D	0.47	0/1445	0.74	0/1943
1	E	0.48	0/1451	0.78	2/1951 (0.1%)
1	F	0.49	0/1459	0.77	4/1962 (0.2%)
All	All	0.49	0/8696	0.78	12/11693 (0.1%)

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	E	0	2

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	185	LEU	CB-CG-CD2	7.62	123.95	111.00
1	F	208	LEU	CB-CG-CD1	6.90	122.72	111.00
1	E	131	LEU	CB-CG-CD2	6.76	122.48	111.00
1	C	122	LEU	CB-CG-CD1	6.70	122.40	111.00
1	F	208	LEU	CB-CG-CD2	-6.44	100.05	111.00
1	F	208	LEU	CA-CB-CG	6.43	130.09	115.30
1	C	131	LEU	CB-CG-CD2	5.91	121.05	111.00
1	F	131	LEU	CB-CG-CD1	5.82	120.89	111.00
1	E	56	ARG	C-N-CA	5.75	134.37	122.30
1	B	206	MET	CG-SD-CE	5.28	108.64	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	138	VAL	CA-CB-CG1	5.06	118.49	110.90
1	B	52	ARG	NE-CZ-NH1	5.02	122.81	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	56	ARG	Peptide
1	E	57	GLY	Mainchain

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	1423	0	1439	18	0
1	B	1417	0	1434	16	0
1	C	1417	0	1434	19	0
1	D	1417	0	1434	13	0
1	E	1423	0	1439	20	0
1	F	1431	0	1443	22	0
2	A	4	0	3	0	0
2	B	4	0	3	1	0
2	C	4	0	3	0	0
2	D	4	0	3	0	0
2	E	4	0	3	0	0
2	F	4	0	3	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
All	All	8563	0	8641	106	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 6.

All (106) 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:E:202:PHE:O	1:E:205:VAL:HG12	1.72	0.88
1:F:127:TYR:O	1:F:131:LEU:HD12	1.83	0.78
1:E:187:PHE:O	1:E:190:VAL:HG22	1.87	0.74
1:A:187:PHE:O	1:A:190:VAL:HG22	1.87	0.74
1:C:127:TYR:O	1:C:131:LEU:HD22	1.90	0.72
1:E:127:TYR:O	1:E:131:LEU:HD22	1.90	0.72
1:E:179:ILE:HG22	1:E:205:VAL:HG23	1.75	0.69
1:D:138:VAL:HG11	1:D:148:TYR:CZ	2.31	0.66
1:F:131:LEU:HD23	1:F:155:TYR:CE1	2.32	0.65
1:A:138:VAL:HG11	1:A:148:TYR:CZ	2.31	0.65
1:F:138:VAL:HG11	1:F:148:TYR:CZ	2.32	0.64
1:E:138:VAL:HG11	1:E:148:TYR:CZ	2.32	0.64
1:C:138:VAL:HG11	1:C:148:TYR:CZ	2.32	0.63
1:F:138:VAL:HG11	1:F:148:TYR:CE1	2.37	0.60
1:F:131:LEU:CD2	1:F:155:TYR:CD1	2.84	0.60
1:D:138:VAL:HG11	1:D:148:TYR:CE1	2.37	0.59
1:E:138:VAL:HG11	1:E:148:TYR:CE1	2.38	0.59
1:E:179:ILE:CG2	1:E:205:VAL:HG23	2.33	0.59
1:F:131:LEU:HD21	1:F:155:TYR:CD1	2.37	0.59
1:A:138:VAL:HG11	1:A:148:TYR:CE1	2.37	0.58
1:E:58:SER:OG	1:E:61:GLU:OE1	2.19	0.58
1:C:138:VAL:HG11	1:C:148:TYR:CE1	2.38	0.58
1:B:54:LEU:HD22	1:B:62:ALA:CB	2.34	0.57
1:B:179:ILE:HD13	1:B:208:LEU:HB3	1.87	0.55
1:C:54:LEU:HD22	1:C:62:ALA:CB	2.35	0.55
1:F:138:VAL:CG1	1:F:148:TYR:CZ	2.91	0.54
1:B:179:ILE:CD1	1:B:208:LEU:HB3	2.37	0.54
1:F:204:LYS:O	1:F:208:LEU:HD22	2.08	0.53
1:D:138:VAL:CG1	1:D:148:TYR:CZ	2.91	0.53
1:E:138:VAL:CG1	1:E:148:TYR:CZ	2.91	0.53
1:F:124:ILE:N	1:F:124:ILE:HD13	2.23	0.53
1:A:138:VAL:CG1	1:A:148:TYR:CZ	2.91	0.53
1:C:138:VAL:CG1	1:C:148:TYR:CZ	2.91	0.53
1:E:86:ALA:O	1:E:90:LEU:HD22	2.10	0.52
1:E:134:LEU:O	1:E:138:VAL:HG23	2.11	0.51
1:F:131:LEU:CD2	1:F:155:TYR:CE1	2.93	0.51
1:F:134:LEU:O	1:F:138:VAL:HG23	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:134:LEU:O	1:D:138:VAL:HG23	2.12	0.50
1:C:134:LEU:O	1:C:138:VAL:HG23	2.11	0.50
1:C:125:LYS:HD2	1:C:127:TYR:OH	2.11	0.50
1:A:56:ARG:NH1	1:D:209:GLU:OE2	2.45	0.50
1:E:179:ILE:H	1:E:179:ILE:HD12	1.76	0.50
1:A:134:LEU:O	1:A:138:VAL:HG23	2.11	0.49
1:A:138:VAL:CG1	1:A:148:TYR:CE2	2.96	0.49
1:D:138:VAL:CG1	1:D:148:TYR:CE2	2.96	0.48
1:E:138:VAL:CG1	1:E:148:TYR:CE2	2.97	0.48
1:C:138:VAL:CG1	1:C:148:TYR:CE2	2.97	0.48
1:C:179:ILE:H	1:C:179:ILE:HD12	1.77	0.48
1:F:60:GLU:OE1	1:F:90:LEU:CD1	2.61	0.47
1:A:179:ILE:HD12	1:A:179:ILE:H	1.78	0.47
1:A:201:HIS:O	1:A:205:VAL:HG23	2.14	0.47
1:D:179:ILE:H	1:D:179:ILE:HD12	1.78	0.47
1:F:138:VAL:CG1	1:F:148:TYR:CE2	2.97	0.47
1:F:201:HIS:O	1:F:205:VAL:HG23	2.15	0.47
1:D:201:HIS:O	1:D:205:VAL:HG23	2.15	0.47
1:C:50:SER:CB	1:C:66:LEU:CD1	2.93	0.47
1:C:201:HIS:O	1:C:205:VAL:HG23	2.14	0.47
1:C:50:SER:HB2	1:C:66:LEU:CD1	2.45	0.47
1:B:64:ARG:O	1:B:68:LEU:HD23	2.15	0.47
1:B:201:HIS:O	1:B:205:VAL:HG23	2.15	0.46
1:B:179:ILE:HD13	1:B:208:LEU:CB	2.46	0.46
1:F:179:ILE:HD13	1:F:208:LEU:HB3	1.97	0.46
1:C:66:LEU:HD23	1:C:83:MET:HA	1.98	0.46
1:E:64:ARG:O	1:E:68:LEU:HD23	2.15	0.46
1:A:64:ARG:O	1:A:68:LEU:HD23	2.16	0.46
1:B:124:ILE:HG21	1:B:126:GLN:HE21	1.82	0.45
1:B:179:ILE:HG22	1:B:205:VAL:HG13	1.98	0.45
1:F:179:ILE:CD1	1:F:208:LEU:HB3	2.46	0.45
1:D:93:TYR:HB3	1:D:124:ILE:HD11	1.97	0.45
1:A:124:ILE:HG22	1:A:124:ILE:O	2.17	0.44
1:F:179:ILE:HG22	1:F:205:VAL:HG13	1.99	0.44
1:C:50:SER:O	1:C:54:LEU:HD23	2.17	0.44
1:B:124:ILE:O	1:B:124:ILE:HG22	2.18	0.44
1:C:179:ILE:HG22	1:C:205:VAL:HG13	1.99	0.44
1:D:179:ILE:HG22	1:D:205:VAL:HG13	1.99	0.44
1:A:179:ILE:HG22	1:A:205:VAL:HG13	1.99	0.44
1:E:93:TYR:HB2	1:E:124:ILE:HD11	2.00	0.43
1:F:60:GLU:OE1	1:F:90:LEU:HD11	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:TYR:HB3	1:A:124:ILE:HD11	2.00	0.43
1:B:50:SER:O	1:B:54:LEU:HD23	2.18	0.43
1:E:194:ASP:OD1	1:E:196:LYS:HD2	2.18	0.42
1:C:124:ILE:HG22	1:C:124:ILE:O	2.19	0.42
1:C:93:TYR:HB3	1:C:124:ILE:HD11	2.01	0.42
1:E:124:ILE:O	1:E:124:ILE:HG22	2.20	0.42
1:F:179:ILE:HD13	1:F:208:LEU:CB	2.49	0.42
1:A:105:LYS:NZ	1:B:98:THR:HG22	2.35	0.42
1:E:116:ARG:O	1:E:120:VAL:HG13	2.20	0.42
1:F:115:TYR:CD1	2:F:301:ACT:H3	2.55	0.41
1:F:93:TYR:CB	1:F:124:ILE:HD11	2.50	0.41
1:F:179:ILE:CG2	1:F:205:VAL:HG13	2.51	0.41
1:A:102:LYS:O	1:A:106:LEU:HD13	2.20	0.41
1:B:179:ILE:CG2	1:B:205:VAL:HG13	2.50	0.41
1:D:179:ILE:CG2	1:D:205:VAL:HG13	2.50	0.41
1:C:116:ARG:O	1:C:120:VAL:HG13	2.20	0.41
1:C:179:ILE:CG2	1:C:205:VAL:HG13	2.51	0.41
1:A:179:ILE:CG2	1:A:205:VAL:HG13	2.51	0.41
1:D:116:ARG:O	1:D:120:VAL:HG13	2.20	0.41
1:A:124:ILE:HG21	1:A:126:GLN:HE21	1.85	0.41
1:E:54:LEU:O	1:E:57:GLY:CA	2.69	0.41
1:B:116:ARG:O	1:B:120:VAL:HG13	2.20	0.41
1:D:102:LYS:O	1:D:106:LEU:HD13	2.20	0.41
1:B:115:TYR:CD1	2:B:301:ACT:H2	2.56	0.41
1:E:128:LYS:HD2	1:E:158:ASP:OD1	2.21	0.40
1:A:116:ARG:O	1:A:120:VAL:HG13	2.20	0.40
1:B:93:TYR:HB3	1:B:124:ILE:HD11	2.02	0.40
1:B:162:GLU:H	1:B:162:GLU:CD	2.25	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	175/202 (87%)	168 (96%)	7 (4%)	0	100	100
1	B	174/202 (86%)	168 (97%)	6 (3%)	0	100	100
1	C	174/202 (86%)	168 (97%)	6 (3%)	0	100	100
1	D	174/202 (86%)	168 (97%)	6 (3%)	0	100	100
1	E	175/202 (87%)	167 (95%)	8 (5%)	0	100	100
1	F	176/202 (87%)	169 (96%)	7 (4%)	0	100	100
All	All	1048/1212 (86%)	1008 (96%)	40 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	146/168 (87%)	135 (92%)	11 (8%)	13	39
1	B	145/168 (86%)	135 (93%)	10 (7%)	15	44
1	C	145/168 (86%)	133 (92%)	12 (8%)	11	35
1	D	145/168 (86%)	135 (93%)	10 (7%)	15	44
1	E	146/168 (87%)	138 (94%)	8 (6%)	21	52
1	F	147/168 (88%)	135 (92%)	12 (8%)	11	36
All	All	874/1008 (87%)	811 (93%)	63 (7%)	14	41

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

Mol	Chain	Res	Type
1	A	48	ASN
1	A	90	LEU
1	A	102	LYS
1	A	103	VAL
1	A	115	TYR
1	A	120	VAL
1	A	129	LEU

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Mol	Chain	Res	Type
1	A	138	VAL
1	A	179	ILE
1	A	196	LYS
1	A	206	MET
1	B	48	ASN
1	B	56	ARG
1	B	90	LEU
1	B	102	LYS
1	B	103	VAL
1	B	115	TYR
1	B	120	VAL
1	B	129	LEU
1	B	179	ILE
1	B	206	MET
1	C	48	ASN
1	C	56	ARG
1	C	66	LEU
1	C	90	LEU
1	C	102	LYS
1	C	103	VAL
1	C	115	TYR
1	C	120	VAL
1	C	122	LEU
1	C	128	LYS
1	C	138	VAL
1	C	206	MET
1	D	48	ASN
1	D	90	LEU
1	D	103	VAL
1	D	115	TYR
1	D	120	VAL
1	D	138	VAL
1	D	161	LEU
1	D	179	ILE
1	D	206	MET
1	D	213	ASP
1	E	48	ASN
1	E	103	VAL
1	E	115	TYR
1	E	120	VAL
1	E	138	VAL
1	E	149	TYR

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Mol	Chain	Res	Type
1	E	196	LYS
1	E	206	MET
1	F	48	ASN
1	F	74	LYS
1	F	90	LEU
1	F	103	VAL
1	F	115	TYR
1	F	120	VAL
1	F	124	ILE
1	F	129	LEU
1	F	138	VAL
1	F	179	ILE
1	F	206	MET
1	F	208	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	GLN
1	A	126	GLN
1	A	182	HIS
1	A	189	ASN
1	B	126	GLN
1	D	182	HIS
1	E	182	HIS
1	F	89	ASN
1	F	126	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 carbohydrates in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	ACT	C	301	-	1,3,3	2.02	1 (100%)	0,3,3	0.00	-
2	ACT	A	301	-	1,3,3	1.73	0	0,3,3	0.00	-
2	ACT	F	301	-	1,3,3	1.61	0	0,3,3	0.00	-
2	ACT	D	301	-	1,3,3	1.55	0	0,3,3	0.00	-
2	ACT	E	301	-	1,3,3	1.58	0	0,3,3	0.00	-
2	ACT	B	301	-	1,3,3	1.66	0	0,3,3	0.00	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	ACT	CH3-C	2.02	1.51	1.48

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	301	ACT	1	0
2	B	301	ACT	1	0

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

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^{th} 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	177/202 (87%)	0.14	1 (0%) 89 90	67, 89, 137, 180	1 (0%)
1	B	176/202 (87%)	0.19	0 100 100	67, 88, 134, 168	0
1	C	176/202 (87%)	0.24	0 100 100	82, 107, 145, 167	0
1	D	176/202 (87%)	0.27	1 (0%) 89 90	74, 97, 133, 163	0
1	E	177/202 (87%)	0.17	2 (1%) 80 80	84, 105, 140, 176	0
1	F	178/202 (88%)	0.26	2 (1%) 80 80	79, 98, 148, 200	0
All	All	1060/1212 (87%)	0.21	6 (0%) 89 90	67, 100, 140, 200	1 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	215	ASP	3.7
1	E	214	SER	2.8
1	F	214	SER	2.5
1	D	55	ARG	2.3
1	E	213	ASP	2.1
1	A	177	GLU	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	CL	E	302	1/1	0.52	0.53	122,122,122,122	0
2	ACT	E	301	4/4	0.72	0.76	119,130,134,137	0
2	ACT	C	301	4/4	0.73	0.64	113,121,125,127	0
2	ACT	F	301	4/4	0.75	0.91	108,118,125,126	0
3	CL	D	302	1/1	0.80	0.46	108,108,108,108	0
2	ACT	D	301	4/4	0.80	0.66	104,117,121,122	0
2	ACT	A	301	4/4	0.81	0.45	104,111,113,115	0
3	CL	C	302	1/1	0.84	0.30	117,117,117,117	0
2	ACT	B	301	4/4	0.84	0.56	100,112,116,117	0
3	CL	B	302	1/1	0.90	0.47	112,112,112,112	0
3	CL	C	303	1/1	0.92	0.40	106,106,106,106	0
3	CL	A	302	1/1	0.94	0.31	105,105,105,105	0

6.5 Other polymers (i)

There are no such residues in this entry.