



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 07:00 am BST

PDB ID : 5HO5
Title : MamB
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Deposited on : 2016-01-19
Resolution : 1.99 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
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

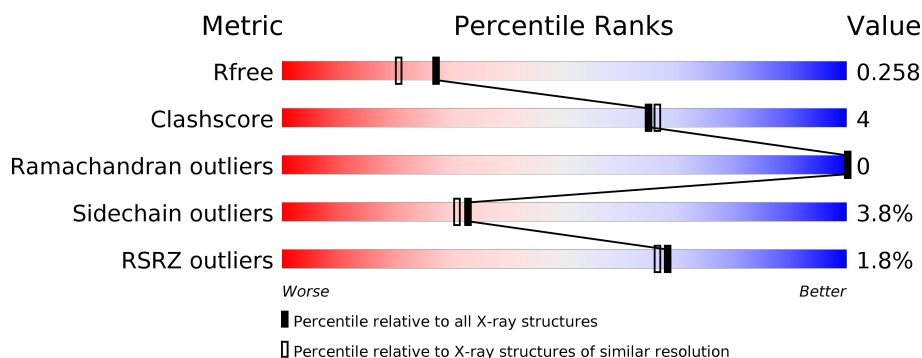
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 $\leq 5\%$. 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	85	<div> <div>2%</div> <div> <div></div> <div>84%</div> <div>9%</div> <div>• • •</div> </div> </div>
1	B	85	<div> <div>4%</div> <div> <div></div> <div>81%</div> <div>16%</div> <div>• •</div> </div> </div>
1	C	85	<div> <div></div> <div> <div></div> <div>85%</div> <div>11%</div> <div>5%</div> </div> </div>
1	D	85	<div> <div>•</div> <div> <div></div> <div>87%</div> <div>8%</div> <div>• •</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2887 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 MamB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	83	Total	C	N	O	S	0	3	0
			680	426	126	127	1			
1	B	84	Total	C	N	O	S	0	2	0
			672	421	124	126	1			
1	C	81	Total	C	N	O	S	0	4	0
			661	412	121	127	1			
1	D	84	Total	C	N	O	S	0	1	0
			667	416	124	126	1			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	209	GLY	-	expression tag	UNP W6KHH6
A	210	SER	-	expression tag	UNP W6KHH6
A	211	HIS	-	expression tag	UNP W6KHH6
A	212	MET	-	expression tag	UNP W6KHH6
A	293	ALA	-	expression tag	UNP W6KHH6
B	209	GLY	-	expression tag	UNP W6KHH6
B	210	SER	-	expression tag	UNP W6KHH6
B	211	HIS	-	expression tag	UNP W6KHH6
B	212	MET	-	expression tag	UNP W6KHH6
B	293	ALA	-	expression tag	UNP W6KHH6
C	209	GLY	-	expression tag	UNP W6KHH6
C	210	SER	-	expression tag	UNP W6KHH6
C	211	HIS	-	expression tag	UNP W6KHH6
C	212	MET	-	expression tag	UNP W6KHH6
C	293	ALA	-	expression tag	UNP W6KHH6
D	209	GLY	-	expression tag	UNP W6KHH6
D	210	SER	-	expression tag	UNP W6KHH6
D	211	HIS	-	expression tag	UNP W6KHH6
D	212	MET	-	expression tag	UNP W6KHH6
D	293	ALA	-	expression tag	UNP W6KHH6

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		

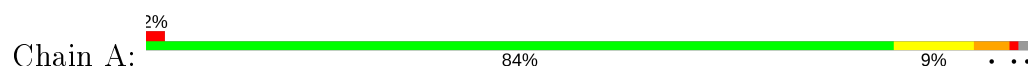
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	52	Total	O	0	0
			52	52		
3	B	55	Total	O	0	0
			55	55		
3	C	53	Total	O	0	0
			53	53		
3	D	46	Total	O	0	0
			46	46		

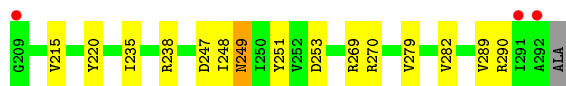
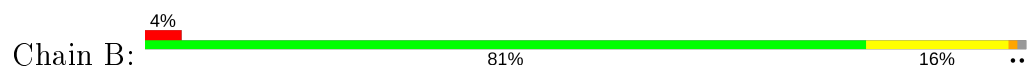
3 Residue-property plots [i](#)

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.

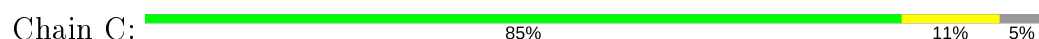
- Molecule 1: Magnetosome protein MamB



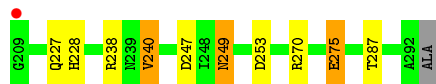
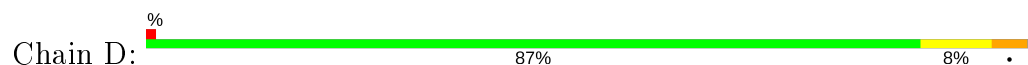
- Molecule 1: Magnetosome protein MamB



- Molecule 1: Magnetosome protein MamB



- Molecule 1: Magnetosome protein MamB



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	37.16Å 96.66Å 96.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.44 – 1.99 24.38 – 1.99	Depositor EDS
% Data completeness (in resolution range)	92.2 (68.44-1.99) 92.4 (24.38-1.99)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 1.99Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.207 , 0.249 0.215 , 0.258	Depositor DCC
R_{free} test set	1168 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	21.8	Xtriage
Anisotropy	0.246	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2887	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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 44.78 % 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 1.4535e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.97	0/701	1.11	7/951 (0.7%)
1	B	0.89	0/689	1.15	8/936 (0.9%)
1	C	0.93	0/684	1.12	8/929 (0.9%)
1	D	0.93	1/681 (0.1%)	1.14	6/925 (0.6%)
All	All	0.93	1/2755 (0.0%)	1.13	29/3741 (0.8%)

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	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	275	GLU	CG-CD	5.58	1.60	1.51

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	253	ASP	CB-CG-OD1	9.38	126.74	118.30
1	D	253	ASP	CB-CG-OD1	9.02	126.42	118.30
1	B	269	ARG	NE-CZ-NH1	8.39	124.50	120.30
1	B	253	ASP	CB-CG-OD1	7.71	125.23	118.30
1	C	280	ARG	NE-CZ-NH1	-7.17	116.71	120.30
1	C	270	ARG	NE-CZ-NH1	7.02	123.81	120.30
1	B	238	ARG	NE-CZ-NH2	-6.87	116.86	120.30
1	B	269	ARG	NE-CZ-NH2	-6.78	116.91	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	270	ARG	NE-CZ-NH1	6.54	123.57	120.30
1	D	253	ASP	CB-CG-OD2	-6.21	112.72	118.30
1	B	238	ARG	NE-CZ-NH1	6.20	123.40	120.30
1	B	290	ARG	NE-CZ-NH2	6.18	123.39	120.30
1	D	270	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	C	280	ARG	NE-CZ-NH2	6.04	123.32	120.30
1	B	270	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	D	247	ASP	CB-CG-OD1	5.83	123.54	118.30
1	C	270	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	D	238	ARG	NE-CZ-NH2	-5.55	117.52	120.30
1	C	247[A]	ASP	CB-CG-OD1	5.55	123.30	118.30
1	C	247[B]	ASP	CB-CG-OD1	5.55	123.30	118.30
1	D	270	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	C	238	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	A	269[A]	ARG	NE-CZ-NH1	5.37	122.99	120.30
1	A	269[B]	ARG	NE-CZ-NH1	5.37	122.99	120.30
1	A	270	ARG	NE-CZ-NH2	-5.13	117.73	120.30
1	A	269[A]	ARG	NE-CZ-NH2	-5.12	117.74	120.30
1	A	269[B]	ARG	NE-CZ-NH2	-5.12	117.74	120.30
1	C	238	ARG	NE-CZ-NH2	-5.05	117.77	120.30
1	B	253	ASP	CB-CG-OD2	-5.05	113.75	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	210	SER	Peptide

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	680	0	668	6	0
1	B	672	0	665	9	0
1	C	661	0	649	3	0
1	D	667	0	654	5	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	52	0	0	0	0
3	B	55	0	0	0	0
3	C	53	0	0	1	0
3	D	46	0	0	1	0
All	All	2887	0	2636	22	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 (22) 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:B:235:ILE:HG13	1:B:248[B]:ILE:CD1	2.17	0.74
1:D:227:GLN:NE2	3:D:302:HOH:O	2.27	0.68
1:A:210:SER:HA	1:A:212:MET:N	2.19	0.58
1:D:227:GLN:HG2	1:D:228:HIS:CD2	2.40	0.57
1:A:280:ARG:CG	1:A:280:ARG:HH21	2.17	0.56
1:D:240:VAL:CG2	1:D:240:VAL:O	2.56	0.53
1:B:251:TYR:CD1	1:B:289[A]:VAL:CG1	2.94	0.51
1:C:227:GLN:HG3	1:C:228:HIS:CD2	2.46	0.51
1:D:240:VAL:O	1:D:240:VAL:HG23	2.14	0.48
1:B:248[A]:ILE:HG22	1:B:249:ASN:N	2.31	0.46
1:B:220:TYR:HH	1:C:210:SER:N	2.15	0.45
1:B:215:VAL:HG13	1:B:279:VAL:CG2	2.47	0.45
1:B:247:ASP:O	1:B:248[B]:ILE:HD13	2.18	0.44
1:B:279:VAL:CG1	1:B:282:VAL:CG2	2.96	0.43
1:D:249:ASN:ND2	1:D:287:THR:OG1	2.49	0.43
1:A:242:GLU:O	1:A:242:GLU:HG3	2.19	0.43
1:C:211:HIS:HB2	3:C:334:HOH:O	2.18	0.43
1:A:280:ARG:CG	1:A:280:ARG:NH2	2.82	0.42
1:A:210:SER:HA	1:A:211:HIS:C	2.39	0.42
1:B:279:VAL:HG11	1:B:282:VAL:CG2	2.50	0.41
1:A:269[A]:ARG:NH2	1:A:284:VAL:H	2.19	0.41
1:B:279:VAL:HG11	1:B:282:VAL:HG22	2.03	0.40

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	84/85 (99%)	83 (99%)	1 (1%)	0	100	100
1	B	84/85 (99%)	84 (100%)	0	0	100	100
1	C	83/85 (98%)	82 (99%)	1 (1%)	0	100	100
1	D	83/85 (98%)	82 (99%)	1 (1%)	0	100	100
All	All	334/340 (98%)	331 (99%)	3 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	75/72 (104%)	69 (92%)	6 (8%)	12	7
1	B	74/72 (103%)	73 (99%)	1 (1%)	67	72
1	C	75/72 (104%)	74 (99%)	1 (1%)	69	74
1	D	73/72 (101%)	70 (96%)	3 (4%)	30	28
All	All	297/288 (103%)	286 (96%)	11 (4%)	33	32

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	210	SER
1	A	227	GLN

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Mol	Chain	Res	Type
1	A	242	GLU
1	A	249	ASN
1	A	280	ARG
1	A	290	ARG
1	B	249	ASN
1	C	290	ARG
1	D	240	VAL
1	D	249	ASN
1	D	275	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	249	ASN
1	B	249	ASN
1	D	249	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [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, 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	83/85 (97%)	0.01	2 (2%) 59 57	14, 23, 51, 75	0
1	B	84/85 (98%)	-0.01	3 (3%) 42 42	16, 25, 46, 75	0
1	C	81/85 (95%)	-0.03	0 100 100	16, 25, 52, 63	0
1	D	84/85 (98%)	0.06	1 (1%) 79 78	17, 28, 44, 72	0
All	All	332/340 (97%)	0.01	6 (1%) 68 66	14, 26, 52, 75	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	292	ALA	4.7
1	B	292	ALA	3.9
1	A	290	ARG	3.1
1	B	209	GLY	2.8
1	B	291	ILE	2.3
1	D	209	GLY	2.1

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 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, 95th 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(\AA^2)	Q<0.9
2	MG	B	301	1/1	0.90	0.13	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.