



wwPDB EM Map/Model Validation Report ⓘ

Apr 10, 2016 – 01:34 PM BST

PDB ID : 2BGZ
EMDB ID: : EMD-1132
Title : ATOMIC MODEL OF THE BACTERIAL FLAGELLAR BASED ON DOCK-
ING AN X-RAY DERIVED HOOK STRUCTURE INTO AN EM MAP.
Authors : Shaikh, T.R.; Thomas, D.R.; Chen, J.Z.; Samatey, F.A.; Matsunami, H.;
Imada, K.; Namba, K.; Derosier, D.J.
Deposited on : 2005-01-06
Resolution : 12.00 Å(reported)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could
stem from errors in the original structure(s) used in the fitting.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/EMValidationReportHelp>

MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : trunk27241

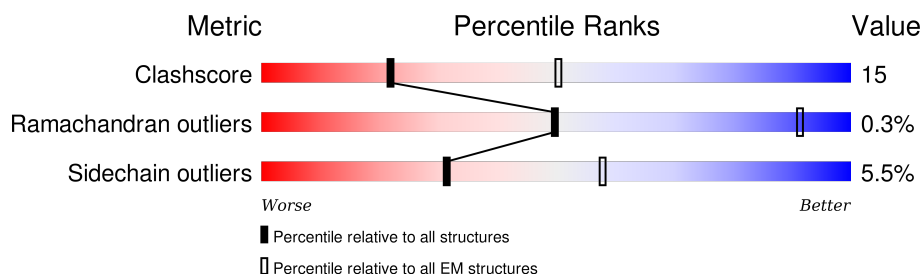
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 12.00 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)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	299	

2 Entry composition [i](#)

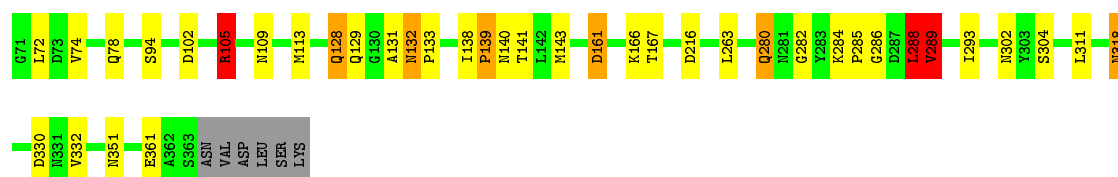
There is only 1 type of molecule in this entry. The entry contains 2156 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 FLAGELLAR HOOK PROTEIN FLGE.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	293	Total	C	N	O	S	0	0
			2156	1327	369	454	6		

- Molecule 1: FLAGELLAR HOOK PROTEIN FLGE



4 Experimental information

Property	Value	Source
Reconstruction method	CRYSTALLOGRAPHY	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	Not provided	Depositor
Microscope	FEI CM200 FEG	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	Not provided	Depositor
Minimum defocus (nm)	1300	Depositor
Maximum defocus (nm)	2700	Depositor
Magnification	66000	Depositor
Image detector	KODAK SO-163 FILM	Depositor

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	$\# Z > 2$	RMSZ	$\# Z > 2$
1	A	0.66	1/2193 (0.0%)	1.21	12/2993 (0.4%)

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	A	139	PRO	C-N	-5.20	1.22	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	139	PRO	CA-C-N	-30.50	50.09	117.20
1	A	139	PRO	C-N-CA	-24.51	60.42	121.70
1	A	139	PRO	O-C-N	21.78	157.55	122.70
1	A	105	ARG	NE-CZ-NH1	8.41	124.50	120.30
1	A	288	LEU	CB-CA-C	-7.22	96.47	110.20
1	A	289	VAL	N-CA-C	6.43	128.37	111.00
1	A	105	ARG	NE-CZ-NH2	-6.38	117.11	120.30
1	A	143	MET	CG-SD-CE	6.02	109.83	100.20
1	A	330	ASP	CB-CG-OD2	6.01	123.71	118.30
1	A	216	ASP	CB-CG-OD2	5.25	123.03	118.30
1	A	102	ASP	CB-CG-OD1	5.19	122.97	118.30
1	A	161	ASP	CB-CG-OD1	5.13	122.92	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	139	PRO	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	2156	0	2054	64	0
All	All	2156	0	2054	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 15.

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:A:105:ARG:HD2	1:A:140:ASN:CB	1.46	1.43
1:A:288:LEU:CG	1:A:289:VAL:N	1.77	1.42
1:A:288:LEU:CD2	1:A:289:VAL:H	1.37	1.36
1:A:105:ARG:CD	1:A:140:ASN:HB3	1.55	1.35
1:A:288:LEU:CG	1:A:289:VAL:HG23	1.60	1.31
1:A:288:LEU:CD2	1:A:302:ASN:O	1.81	1.27
1:A:105:ARG:HD2	1:A:140:ASN:CG	1.54	1.27
1:A:288:LEU:HD21	1:A:302:ASN:O	1.35	1.24
1:A:288:LEU:CD1	1:A:289:VAL:HG23	1.73	1.17
1:A:288:LEU:HG	1:A:289:VAL:CG2	1.75	1.16
1:A:288:LEU:CD2	1:A:289:VAL:N	2.01	1.16
1:A:105:ARG:CD	1:A:140:ASN:ND2	2.17	1.08
1:A:288:LEU:HD23	1:A:289:VAL:H	0.96	1.08
1:A:105:ARG:HD2	1:A:140:ASN:HB3	1.09	1.06
1:A:105:ARG:HD2	1:A:140:ASN:ND2	1.71	1.05
1:A:288:LEU:HG	1:A:289:VAL:HG23	1.11	1.05
1:A:288:LEU:HD23	1:A:289:VAL:N	1.67	1.03
1:A:105:ARG:CG	1:A:140:ASN:ND2	2.21	1.03
1:A:105:ARG:CD	1:A:140:ASN:HD22	1.71	1.02
1:A:288:LEU:HD22	1:A:302:ASN:O	1.63	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:ARG:CG	1:A:140:ASN:HD22	1.84	0.88
1:A:288:LEU:HG	1:A:289:VAL:N	1.32	0.88
1:A:105:ARG:HG3	1:A:140:ASN:ND2	1.90	0.84
1:A:288:LEU:CG	1:A:289:VAL:CG2	2.44	0.84
1:A:105:ARG:CD	1:A:140:ASN:CB	2.28	0.81
1:A:105:ARG:NE	1:A:140:ASN:HB3	1.95	0.81
1:A:288:LEU:CD1	1:A:289:VAL:CG2	2.57	0.80
1:A:288:LEU:CD2	1:A:289:VAL:CA	2.63	0.76
1:A:288:LEU:HG	1:A:289:VAL:CA	2.17	0.74
1:A:138:ILE:HG23	1:A:311:LEU:HD13	1.71	0.72
1:A:288:LEU:HD21	1:A:289:VAL:HB	1.73	0.71
1:A:288:LEU:HG	1:A:289:VAL:CB	2.21	0.70
1:A:105:ARG:CZ	1:A:140:ASN:HB3	2.23	0.69
1:A:128:GLN:HE21	1:A:128:GLN:H	1.42	0.68
1:A:105:ARG:HD3	1:A:138:ILE:O	1.93	0.68
1:A:318:ASN:OD1	1:A:351:ASN:ND2	2.32	0.63
1:A:288:LEU:HA	1:A:304:SER:HB3	1.80	0.62
1:A:105:ARG:HD3	1:A:140:ASN:HB3	1.75	0.61
1:A:105:ARG:NE	1:A:140:ASN:HD22	1.97	0.61
1:A:288:LEU:HD12	1:A:289:VAL:HG23	1.76	0.60
1:A:72:LEU:HG	1:A:361:GLU:HG3	1.83	0.60
1:A:288:LEU:HD23	1:A:289:VAL:CA	2.28	0.60
1:A:288:LEU:CG	1:A:289:VAL:H	1.56	0.59
1:A:288:LEU:CG	1:A:289:VAL:CB	2.81	0.57
1:A:94:SER:HB2	1:A:332:VAL:HG22	1.86	0.57
1:A:105:ARG:NH1	1:A:140:ASN:HB3	2.21	0.56
1:A:288:LEU:HD11	1:A:289:VAL:CG2	2.33	0.55
1:A:128:GLN:CG	1:A:131:ALA:HB2	2.36	0.55
1:A:132:ASN:HD22	1:A:133:PRO:HD2	1.73	0.54
1:A:288:LEU:CD2	1:A:289:VAL:CB	2.87	0.53
1:A:288:LEU:HD21	1:A:289:VAL:CB	2.40	0.52
1:A:284:LYS:O	1:A:286:GLY:N	2.43	0.52
1:A:128:GLN:HG2	1:A:128:GLN:O	2.11	0.51
1:A:132:ASN:HD22	1:A:133:PRO:CD	2.24	0.51
1:A:109:ASN:HD21	1:A:113:MET:CE	2.26	0.49
1:A:105:ARG:CD	1:A:140:ASN:CG	2.44	0.46
1:A:288:LEU:C	1:A:289:VAL:CG2	2.83	0.45
1:A:288:LEU:CG	1:A:289:VAL:CA	2.79	0.44
1:A:284:LYS:O	1:A:285:PRO:C	2.56	0.44
1:A:74:VAL:HG11	1:A:293:ILE:HD12	2.00	0.44
1:A:132:ASN:HD22	1:A:133:PRO:N	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:GLN:NE2	1:A:282:GLY:H	2.17	0.42
1:A:109:ASN:HD21	1:A:113:MET:HE3	1.85	0.41
1:A:105:ARG:CZ	1:A:140:ASN:CB	2.96	0.41

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 PDB entries followed by that with respect to all EM entries.

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	291/299 (97%)	281 (97%)	9 (3%)	1 (0%)	46 83

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	289	VAL

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 PDB entries followed by that with respect to all EM entries.

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	235/241 (98%)	222 (94%)	13 (6%)	27 63

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	105	ARG
1	A	128	GLN
1	A	129	GLN
1	A	132	ASN
1	A	141	THR
1	A	161	ASP
1	A	166	LYS
1	A	167	THR
1	A	263	LEU
1	A	280	GLN
1	A	288	LEU
1	A	318	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	88	ASN
1	A	98	GLN
1	A	109	ASN
1	A	111	GLN
1	A	128	GLN
1	A	132	ASN
1	A	140	ASN
1	A	196	ASN
1	A	244	ASN
1	A	280	GLN
1	A	318	ASN
1	A	322	ASN
1	A	351	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

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.