



wwPDB EM Map/Model Validation Report ⓘ

Apr 10, 2016 – 01:39 PM BST

PDB ID : 3J0R
EMDB ID: : EMD-5352
Title : Model of a type III secretion system needle based on a 7 Angstrom resolution cryoEM map
Authors : Fujii, T.; Cheung, M.; Blanco, A.; Kato, T.; Blocker, A.J.; Namba, K.
Deposited on : 2011-11-03
Resolution : 7.70 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

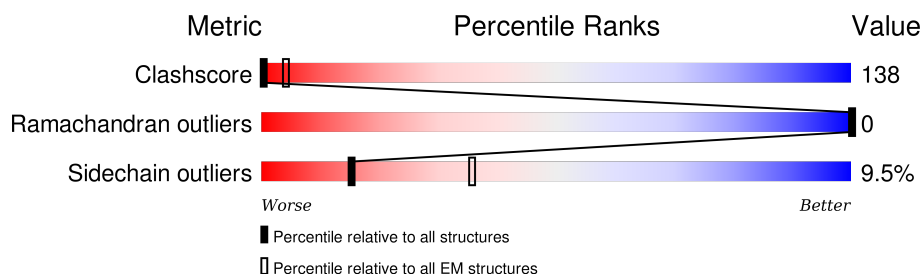
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 7.70 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	82	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 644 atoms, of which 0 are hydrogens and 0 are deuteriums.

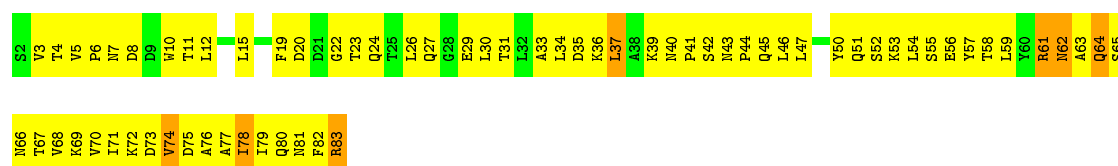
In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Protein mxIH.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
1	A	82	644	402	106	136	0	0

i

- Molecule 1: Protein mxiH



4 Experimental information

Property	Value	Source
Reconstruction method	HELICAL	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	CTFFIND3 Each particle	Depositor
Microscope	JEOL 3200FSC	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	20	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	50000	Depositor
Image detector	TemCam-F415MP	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	1.01	0/653	1.14	0/888

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	644	0	633	176	0
All	All	644	0	633	176	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 138.

All (176) 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:37:LEU:HG	1:A:50:TYR:HB2	1.38	1.06
1:A:80:GLN:HA	1:A:83:ARG:HE	1.27	0.97
1:A:51:GLN:HA	1:A:65:SER:HB3	1.50	0.94
1:A:19:PHE:CD2	1:A:77:ALA:HB1	2.03	0.93
1:A:50:TYR:HE1	1:A:64:GLN:HB3	1.36	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:GLU:HB2	1:A:61:ARG:HH21	1.34	0.90
1:A:30:LEU:HD13	1:A:63:ALA:HB3	1.54	0.88
1:A:55:SER:H	1:A:59:LEU:HB2	1.39	0.88
1:A:54:LEU:HD13	1:A:61:ARG:N	1.91	0.84
1:A:5:VAL:HB	1:A:6:PRO:HD3	1.61	0.83
1:A:77:ALA:HA	1:A:80:GLN:OE1	1.80	0.82
1:A:54:LEU:HD22	1:A:61:ARG:CB	2.11	0.81
1:A:37:LEU:CG	1:A:50:TYR:HB2	2.09	0.81
1:A:29:GLU:HB2	1:A:61:ARG:NH2	1.96	0.79
1:A:50:TYR:O	1:A:65:SER:HB3	1.83	0.79
1:A:55:SER:N	1:A:59:LEU:HB2	1.97	0.79
1:A:52:SER:HB3	1:A:62:ASN:HD21	1.48	0.78
1:A:78:ILE:HA	1:A:81:ASN:OD1	1.84	0.78
1:A:31:THR:O	1:A:34:LEU:HG	1.83	0.77
1:A:37:LEU:HG	1:A:50:TYR:CB	2.14	0.77
1:A:20:ASP:O	1:A:24:GLN:HG3	1.83	0.77
1:A:26:LEU:HB3	1:A:61:ARG:HH12	1.50	0.76
1:A:56:GLU:HG3	1:A:57:TYR:H	1.50	0.76
1:A:54:LEU:HD22	1:A:61:ARG:HB3	1.67	0.75
1:A:66:ASN:OD1	1:A:68:VAL:HG12	1.86	0.75
1:A:53:LYS:O	1:A:59:LEU:HD13	1.86	0.75
1:A:54:LEU:HD22	1:A:69:LYS:HE3	1.67	0.75
1:A:51:GLN:HA	1:A:65:SER:CB	2.16	0.75
1:A:50:TYR:OH	1:A:64:GLN:HG2	1.87	0.74
1:A:54:LEU:HD13	1:A:61:ARG:CA	2.18	0.73
1:A:30:LEU:HB2	1:A:63:ALA:CB	2.19	0.73
1:A:61:ARG:HB2	1:A:69:LYS:CD	2.20	0.72
1:A:50:TYR:HE1	1:A:64:GLN:CB	2.02	0.72
1:A:22:GLY:O	1:A:26:LEU:HG	1.90	0.72
1:A:55:SER:N	1:A:59:LEU:HD12	2.04	0.72
1:A:47:LEU:O	1:A:50:TYR:HB3	1.90	0.71
1:A:53:LYS:HE3	1:A:54:LEU:O	1.91	0.71
1:A:42:SER:O	1:A:44:PRO:HD3	1.92	0.70
1:A:56:GLU:HG3	1:A:57:TYR:N	2.07	0.70
1:A:50:TYR:CE1	1:A:64:GLN:HB3	2.22	0.69
1:A:55:SER:H	1:A:59:LEU:HD12	1.55	0.69
1:A:23:THR:HG21	1:A:74:VAL:HA	1.75	0.69
1:A:61:ARG:CB	1:A:69:LYS:HE3	2.21	0.69
1:A:50:TYR:CE1	1:A:64:GLN:HG2	2.27	0.69
1:A:80:GLN:HG3	1:A:83:ARG:HH21	1.58	0.69
1:A:47:LEU:HD12	1:A:50:TYR:HD2	1.57	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:LEU:HD13	1:A:64:GLN:N	2.08	0.68
1:A:55:SER:HB3	1:A:58:THR:OG1	1.94	0.68
1:A:54:LEU:HB2	1:A:61:ARG:HA	1.75	0.67
1:A:26:LEU:HD22	1:A:61:ARG:NH1	2.08	0.67
1:A:31:THR:HA	1:A:34:LEU:CD2	2.25	0.67
1:A:80:GLN:CA	1:A:83:ARG:HE	2.03	0.67
1:A:11:THR:O	1:A:15:LEU:HG	1.94	0.67
1:A:23:THR:CB	1:A:74:VAL:HG13	2.25	0.67
1:A:43:ASN:OD1	1:A:45:GLN:HB3	1.95	0.66
1:A:26:LEU:HA	1:A:61:ARG:HH22	1.60	0.66
1:A:53:LYS:HD2	1:A:54:LEU:H	1.60	0.66
1:A:54:LEU:HD22	1:A:61:ARG:HB2	1.77	0.66
1:A:52:SER:HB3	1:A:62:ASN:ND2	2.10	0.66
1:A:80:GLN:HA	1:A:83:ARG:NE	2.08	0.65
1:A:40:ASN:ND2	1:A:43:ASN:H	1.95	0.65
1:A:62:ASN:HB2	1:A:65:SER:OG	1.98	0.64
1:A:19:PHE:HB3	1:A:77:ALA:HB1	1.79	0.64
1:A:79:ILE:HG22	1:A:83:ARG:NE	2.12	0.64
1:A:34:LEU:HD12	1:A:35:ASP:N	2.11	0.64
1:A:23:THR:HG22	1:A:73:ASP:OD1	1.98	0.64
1:A:19:PHE:HD2	1:A:77:ALA:HB1	1.61	0.64
1:A:23:THR:HB	1:A:74:VAL:HG13	1.78	0.64
1:A:63:ALA:HB1	1:A:70:VAL:CG2	2.28	0.64
1:A:37:LEU:N	1:A:37:LEU:HD13	2.12	0.63
1:A:30:LEU:CD1	1:A:63:ALA:HB3	2.27	0.63
1:A:55:SER:H	1:A:59:LEU:CB	2.11	0.63
1:A:31:THR:HA	1:A:34:LEU:HD21	1.81	0.62
1:A:75:ASP:O	1:A:78:ILE:HD12	2.00	0.62
1:A:53:LYS:C	1:A:59:LEU:HD13	2.20	0.61
1:A:61:ARG:HB2	1:A:69:LYS:HD2	1.82	0.61
1:A:61:ARG:HG3	1:A:69:LYS:HE3	1.83	0.61
1:A:79:ILE:HG22	1:A:83:ARG:CZ	2.30	0.61
1:A:19:PHE:CG	1:A:77:ALA:HB1	2.36	0.61
1:A:30:LEU:O	1:A:30:LEU:HD12	2.00	0.61
1:A:19:PHE:HB3	1:A:77:ALA:CB	2.31	0.61
1:A:80:GLN:CG	1:A:83:ARG:HH21	2.13	0.60
1:A:47:LEU:HD12	1:A:50:TYR:CD2	2.37	0.60
1:A:3:VAL:O	1:A:6:PRO:HD2	2.00	0.60
1:A:47:LEU:CD1	1:A:50:TYR:HD2	2.13	0.60
1:A:61:ARG:HB2	1:A:69:LYS:HE3	1.83	0.60
1:A:62:ASN:HB2	1:A:65:SER:HG	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:LEU:HD13	1:A:63:ALA:CB	2.31	0.59
1:A:19:PHE:O	1:A:23:THR:HG23	2.03	0.59
1:A:26:LEU:CA	1:A:61:ARG:HH22	2.15	0.59
1:A:50:TYR:CZ	1:A:64:GLN:HG2	2.38	0.58
1:A:67:THR:HG23	1:A:68:VAL:N	2.18	0.58
1:A:39:LYS:O	1:A:41:PRO:HD3	2.03	0.58
1:A:52:SER:H	1:A:65:SER:CB	2.17	0.57
1:A:26:LEU:CB	1:A:61:ARG:HH12	2.15	0.57
1:A:61:ARG:HG3	1:A:69:LYS:CE	2.35	0.57
1:A:54:LEU:HD22	1:A:69:LYS:CE	2.32	0.57
1:A:61:ARG:HB2	1:A:69:LYS:CE	2.35	0.57
1:A:61:ARG:HG3	1:A:69:LYS:NZ	2.20	0.56
1:A:68:VAL:HG13	1:A:69:LYS:N	2.21	0.56
1:A:50:TYR:HE1	1:A:64:GLN:CG	2.17	0.56
1:A:66:ASN:CG	1:A:68:VAL:HG12	2.26	0.56
1:A:63:ALA:HB1	1:A:70:VAL:HG21	1.88	0.55
1:A:63:ALA:HA	1:A:70:VAL:HG22	1.88	0.55
1:A:77:ALA:O	1:A:80:GLN:HB2	2.07	0.55
1:A:19:PHE:CB	1:A:77:ALA:HB1	2.37	0.55
1:A:50:TYR:HE1	1:A:64:GLN:HG2	1.70	0.55
1:A:37:LEU:HD12	1:A:46:LEU:HB3	1.89	0.54
1:A:55:SER:H	1:A:59:LEU:CD1	2.19	0.53
1:A:30:LEU:HB2	1:A:63:ALA:HB3	1.88	0.53
1:A:8:ASP:O	1:A:12:LEU:HG	2.08	0.53
1:A:51:GLN:HG3	1:A:51:GLN:O	2.09	0.53
1:A:27:GLN:O	1:A:30:LEU:HB3	2.07	0.52
1:A:30:LEU:CB	1:A:63:ALA:HB3	2.38	0.52
1:A:61:ARG:CG	1:A:69:LYS:HE3	2.39	0.52
1:A:53:LYS:CD	1:A:54:LEU:H	2.20	0.52
1:A:37:LEU:HD12	1:A:46:LEU:C	2.30	0.52
1:A:72:LYS:HZ2	1:A:76:ALA:HB2	1.74	0.52
1:A:78:ILE:HD12	1:A:78:ILE:H	1.75	0.52
1:A:10:TRP:CE3	1:A:10:TRP:HA	2.45	0.52
1:A:63:ALA:CB	1:A:70:VAL:HG22	2.40	0.51
1:A:61:ARG:HG3	1:A:61:ARG:HH11	1.75	0.51
1:A:63:ALA:HB1	1:A:70:VAL:HG22	1.91	0.51
1:A:72:LYS:NZ	1:A:76:ALA:HB2	2.25	0.51
1:A:34:LEU:HD12	1:A:35:ASP:CA	2.42	0.51
1:A:68:VAL:O	1:A:72:LYS:HB2	2.11	0.50
1:A:5:VAL:CB	1:A:6:PRO:HD3	2.37	0.49
1:A:54:LEU:HB3	1:A:69:LYS:HD2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:LYS:HA	1:A:75:ASP:OD1	2.13	0.49
1:A:26:LEU:HA	1:A:61:ARG:NH2	2.27	0.48
1:A:54:LEU:CB	1:A:69:LYS:HD2	2.43	0.48
1:A:79:ILE:O	1:A:82:PHE:HB2	2.13	0.48
1:A:63:ALA:CA	1:A:70:VAL:HG22	2.42	0.48
1:A:40:ASN:HD21	1:A:43:ASN:H	1.60	0.48
1:A:34:LEU:HD12	1:A:35:ASP:HA	1.96	0.47
1:A:26:LEU:HD22	1:A:61:ARG:HH12	1.78	0.47
1:A:33:ALA:O	1:A:37:LEU:HD22	2.13	0.47
1:A:34:LEU:HB2	1:A:64:GLN:OE1	2.15	0.47
1:A:30:LEU:HD11	1:A:64:GLN:HB2	1.97	0.46
1:A:53:LYS:HD2	1:A:54:LEU:N	2.28	0.46
1:A:54:LEU:CD2	1:A:61:ARG:HB3	2.43	0.46
1:A:78:ILE:O	1:A:81:ASN:HB2	2.16	0.46
1:A:19:PHE:N	1:A:19:PHE:CD1	2.83	0.46
1:A:54:LEU:HD13	1:A:61:ARG:CB	2.46	0.45
1:A:61:ARG:HG3	1:A:61:ARG:NH1	2.32	0.44
1:A:23:THR:HA	1:A:26:LEU:HG	1.99	0.44
1:A:54:LEU:HA	1:A:59:LEU:CB	2.48	0.44
1:A:63:ALA:CA	1:A:70:VAL:CG2	2.95	0.43
1:A:24:GLN:O	1:A:27:GLN:HB3	2.18	0.43
1:A:61:ARG:O	1:A:61:ARG:HD3	2.18	0.43
1:A:54:LEU:HA	1:A:59:LEU:HB2	2.01	0.43
1:A:23:THR:HG21	1:A:74:VAL:CG1	2.48	0.43
1:A:52:SER:CB	1:A:62:ASN:HD21	2.26	0.43
1:A:55:SER:H	1:A:59:LEU:CG	2.32	0.42
1:A:79:ILE:CG2	1:A:83:ARG:CZ	2.95	0.42
1:A:29:GLU:CB	1:A:61:ARG:HH21	2.19	0.42
1:A:19:PHE:CE2	1:A:81:ASN:ND2	2.87	0.42
1:A:7:ASN:O	1:A:11:THR:HG23	2.19	0.42
1:A:31:THR:HA	1:A:34:LEU:HD23	2.00	0.42
1:A:75:ASP:HA	1:A:78:ILE:CD1	2.49	0.42
1:A:68:VAL:CG1	1:A:69:LYS:N	2.83	0.42
1:A:4:THR:O	1:A:7:ASN:HB2	2.19	0.42
1:A:53:LYS:CG	1:A:54:LEU:N	2.83	0.41
1:A:67:THR:CG2	1:A:68:VAL:N	2.83	0.41
1:A:71:ILE:O	1:A:74:VAL:HG23	2.20	0.41
1:A:37:LEU:CD1	1:A:46:LEU:HB3	2.50	0.41
1:A:50:TYR:HH	1:A:64:GLN:HG2	1.83	0.41
1:A:36:LYS:HB3	1:A:46:LEU:HD13	2.03	0.41
1:A:54:LEU:CD2	1:A:69:LYS:HD2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:LYS:HZ2	1:A:76:ALA:CB	2.34	0.41
1:A:79:ILE:HG22	1:A:83:ARG:CD	2.51	0.41
1:A:53:LYS:CD	1:A:54:LEU:N	2.83	0.41
1:A:46:LEU:HD23	1:A:46:LEU:HA	1.94	0.41
1:A:72:LYS:O	1:A:76:ALA:HB2	2.21	0.41
1:A:66:ASN:HB3	1:A:69:LYS:HB2	2.03	0.40
1:A:23:THR:O	1:A:26:LEU:HB2	2.22	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 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	80/82 (98%)	80 (100%)	0	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 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	74/74 (100%)	67 (90%)	7 (10%)	11	41

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	LEU
1	A	61	ARG
1	A	62	ASN
1	A	64	GLN
1	A	74	VAL
1	A	78	ILE
1	A	83	ARG

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	40	ASN
1	A	62	ASN
1	A	66	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](#)

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.