



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:12 AM GMT

PDB ID : 3HLR
Title : Donor strand complemented FaeG of F4ad fimbriae
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Deposited on : 2009-05-28
Resolution : 2.30 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

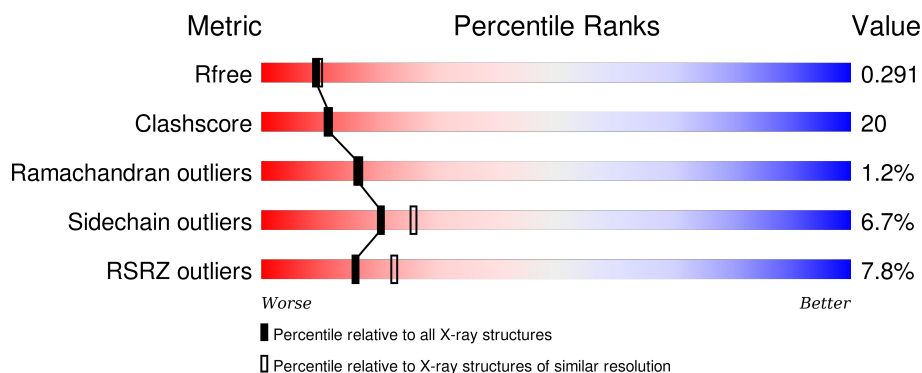
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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. 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	276	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2077 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 K88 fimbrial protein AD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	0	0
			1883	1186	318	375	4			

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	EXPRESSION TAG	UNP P14191
A	1	GLY	-	EXPRESSION TAG	UNP P14191
A	18	SER	ASN	SEE REMARK 999	UNP P14191
A	245	GLY	-	LINKER	UNP P14191
A	246	GLY	-	LINKER	UNP P14191
A	247	GLY	-	LINKER	UNP P14191
A	248	GLY	-	LINKER	UNP P14191
A	249	GLY	-	LINKER	UNP P14191
A	250	GLY	-	LINKER	UNP P14191
A	251	GLY	-	LINKER	UNP P14191
A	252	GLY	-	LINKER	UNP P14191
A	253	GLY	-	LINKER	UNP P14191
A	254	GLY	-	LINKER	UNP P14191

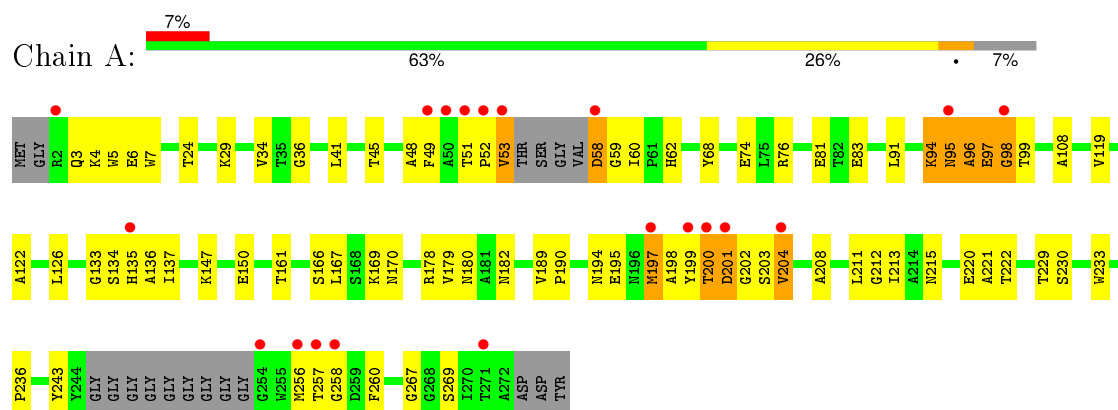
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	194	Total	O	0	0
			194	194		

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 errors 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: K88 fimbrial protein AD



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	37.29 Å 39.73 Å 76.37 Å 90.00° 101.24° 90.00°	Depositor
Resolution (Å)	29.16 – 2.30 30.60 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.9 (29.16-2.30) 95.1 (30.60-2.00)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 2.00 Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.235 , 0.294 0.229 , 0.291	Depositor DCC
R_{free} test set	479 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	19.9	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 38.9	EDS
Estimated twinning fraction	0.057 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 14342 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2077	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.99% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

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	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.28	0/1916	0.50	0/2602

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	94	LYS	Peptide

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	1883	0	1832	76	0
2	A	194	0	0	11	0
All	All	2077	0	1832	76	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 20.

All (76) 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:258:GLY:HA2	1:A:260:PHE:H	1.12	1.09
1:A:36:GLY:H	1:A:215:ASN:HD22	1.03	0.98
1:A:258:GLY:HA2	1:A:260:PHE:N	1.81	0.95
1:A:256:MET:O	1:A:257:THR:HG22	1.67	0.93
1:A:201:ASP:CG	1:A:202:GLY:HA2	1.96	0.86
1:A:211:LEU:HB3	2:A:429:HOH:O	1.78	0.83
1:A:6:GLU:HA	1:A:257:THR:HB	1.65	0.78
1:A:97:GLU:O	1:A:99:THR:N	2.18	0.77
1:A:94:LYS:HD2	1:A:98:GLY:HA2	1.70	0.74
1:A:74:GLU:OE1	1:A:76:ARG:HD2	1.88	0.73
1:A:51:THR:CB	1:A:199:TYR:H	2.03	0.71
1:A:83:GLU:HA	1:A:150:GLU:O	1.92	0.69
1:A:3:GLN:CG	1:A:257:THR:HG21	2.28	0.63
1:A:36:GLY:N	1:A:215:ASN:HD22	1.87	0.63
1:A:41:LEU:HB3	2:A:429:HOH:O	1.98	0.63
1:A:256:MET:HE2	1:A:260:PHE:HA	1.81	0.61
1:A:36:GLY:H	1:A:215:ASN:ND2	1.87	0.61
1:A:3:GLN:HG3	1:A:257:THR:HG21	1.82	0.61
1:A:201:ASP:CB	1:A:202:GLY:HA2	2.28	0.60
1:A:48:ALA:HB1	1:A:204:VAL:HG13	1.83	0.60
1:A:236:PRO:HG2	2:A:447:HOH:O	2.02	0.60
1:A:256:MET:HB3	1:A:260:PHE:HD1	1.67	0.60
1:A:94:LYS:O	1:A:95:ASN:HB2	2.02	0.60
1:A:94:LYS:HD2	1:A:98:GLY:CA	2.31	0.58
1:A:136:ALA:HB2	1:A:194:ASN:HD21	1.69	0.57
1:A:74:GLU:CD	1:A:76:ARG:HH11	2.07	0.56
1:A:62:HIS:HE1	1:A:134:SER:HB2	1.70	0.56
1:A:199:TYR:HB3	1:A:203:SER:HB2	1.87	0.56
1:A:3:GLN:O	1:A:4:LYS:HB2	2.06	0.55
1:A:6:GLU:HA	1:A:257:THR:CB	2.37	0.54
1:A:200:THR:HA	1:A:201:ASP:OD1	2.07	0.54
1:A:256:MET:HB3	1:A:260:PHE:CD1	2.43	0.54
1:A:243:TYR:CE2	1:A:258:GLY:HA3	2.44	0.53
1:A:269:SER:HB3	2:A:377:HOH:O	2.08	0.53
1:A:179:VAL:HG11	1:A:204:VAL:HG21	1.90	0.52
1:A:68:TYR:HB2	2:A:404:HOH:O	2.09	0.52
1:A:52:PRO:O	1:A:53:VAL:HG13	2.09	0.52
1:A:48:ALA:HB1	1:A:204:VAL:CG1	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:GLU:HG3	1:A:197:MET:H	1.74	0.51
1:A:126:LEU:HD22	1:A:169:LYS:HG3	1.93	0.50
1:A:170:ASN:HB3	2:A:384:HOH:O	2.10	0.50
1:A:52:PRO:HG3	1:A:198:ALA:CB	2.41	0.50
1:A:133:GLY:HA3	1:A:135:HIS:CE1	2.47	0.50
1:A:7:TRP:HD1	1:A:257:THR:HA	1.77	0.49
1:A:119:VAL:HB	2:A:424:HOH:O	2.12	0.49
1:A:243:TYR:HE2	1:A:258:GLY:HA3	1.77	0.49
1:A:41:LEU:N	2:A:429:HOH:O	2.45	0.49
1:A:95:ASN:O	1:A:96:ALA:HB3	2.12	0.48
1:A:81:GLU:OE1	1:A:81:GLU:HA	2.12	0.48
1:A:51:THR:HA	1:A:52:PRO:HA	1.68	0.47
1:A:179:VAL:HG12	1:A:180:ASN:ND2	2.30	0.47
1:A:122:ALA:HA	1:A:182:ASN:O	2.15	0.46
1:A:34:VAL:HG21	1:A:213:ILE:HG21	1.97	0.46
1:A:220:GLU:O	1:A:220:GLU:HG3	2.15	0.45
1:A:189:VAL:HA	1:A:190:PRO:HD3	1.79	0.45
1:A:60:ILE:HG23	1:A:137:ILE:HG22	1.98	0.44
1:A:45:THR:HG22	1:A:49:PHE:HE2	1.82	0.44
1:A:5:TRP:C	1:A:257:THR:OG1	2.56	0.44
1:A:233:TRP:CZ3	1:A:267:GLY:HA2	2.53	0.43
1:A:200:THR:HA	1:A:201:ASP:CB	2.47	0.43
1:A:5:TRP:O	1:A:257:THR:OG1	2.37	0.43
1:A:161:THR:HB	1:A:166:SER:HB3	2.00	0.43
1:A:167:LEU:HB2	1:A:208:ALA:HB2	2.00	0.42
1:A:51:THR:C	1:A:198:ALA:HA	2.40	0.42
1:A:24:THR:OG1	1:A:29:LYS:HB3	2.19	0.42
1:A:5:TRP:C	1:A:257:THR:HG1	2.21	0.42
1:A:211:LEU:CB	2:A:429:HOH:O	2.51	0.42
1:A:256:MET:HE1	2:A:322:HOH:O	2.19	0.42
1:A:52:PRO:HG3	1:A:198:ALA:HB2	2.01	0.41
1:A:233:TRP:CH2	1:A:267:GLY:HA2	2.55	0.41
1:A:29:LYS:HD2	1:A:221:ALA:O	2.20	0.41
1:A:256:MET:O	1:A:257:THR:CG2	2.54	0.41
1:A:108:ALA:HB1	1:A:212:GLY:O	2.19	0.41
1:A:135:HIS:HD2	2:A:398:HOH:O	2.03	0.41
1:A:94:LYS:O	1:A:95:ASN:CB	2.68	0.40
1:A:58:ASP:HB3	1:A:59:GLY:H	1.58	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	252/276 (91%)	237 (94%)	12 (5%)	3 (1%)	16	16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	95	ASN
1	A	98	GLY
1	A	96	ALA

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	193/204 (95%)	180 (93%)	13 (7%)	20	26

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

Mol	Chain	Res	Type
1	A	53	VAL
1	A	58	ASP
1	A	91	LEU
1	A	97	GLU
1	A	147	LYS
1	A	178	ARG
1	A	197	MET
1	A	200	THR

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Mol	Chain	Res	Type
1	A	201	ASP
1	A	204	VAL
1	A	222	THR
1	A	229	THR
1	A	230	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	148	ASN
1	A	180	ASN
1	A	194	ASN
1	A	215	ASN
1	A	238	ASN
1	A	261	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

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	258/276 (93%)	0.46	20 (7%) 16 22	12, 22, 41, 58	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	257	THR	8.9
1	A	256	MET	8.5
1	A	200	THR	6.9
1	A	258	GLY	6.5
1	A	197	MET	5.3
1	A	51	THR	4.7
1	A	199	TYR	4.7
1	A	50	ALA	4.1
1	A	53	VAL	3.9
1	A	254	GLY	3.7
1	A	95	ASN	3.6
1	A	49	PHE	3.5
1	A	135	HIS	3.4
1	A	98	GLY	3.2
1	A	58	ASP	3.0
1	A	52	PRO	2.9
1	A	2	ARG	2.6
1	A	271	THR	2.5
1	A	204	VAL	2.3
1	A	201	ASP	2.2

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 [i](#)

There are no carbohydrates 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.