



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 28, 2019 – 12:21 PM EDT

PDB ID : 5Y9K
Title : Structure of the belimumab Fab fragment
Authors : Heo, Y.-S.; Shin, W.
Deposited on : 2017-08-25
Resolution : 1.90 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	:	2.4
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.4

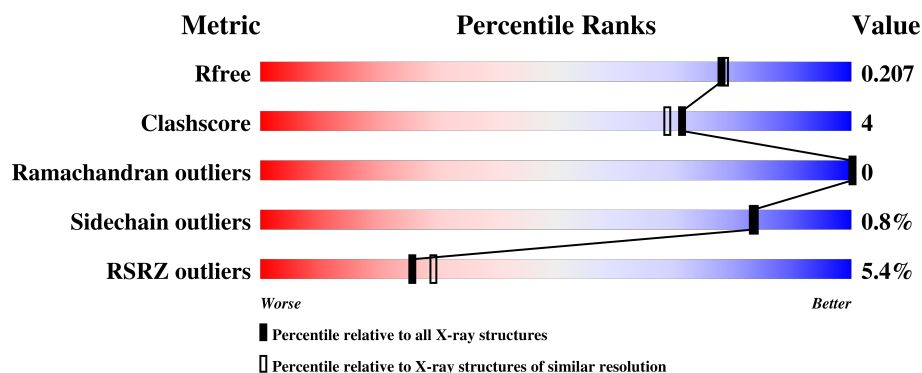
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	111664	5502 (1.90-1.90)
Clashscore	122126	6115 (1.90-1.90)
Ramachandran outliers	120053	6048 (1.90-1.90)
Sidechain outliers	120020	6048 (1.90-1.90)
RSRZ outliers	108989	5379 (1.90-1.90)

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	L	214	<div> <div>5%</div> <div>90%</div> <div>7%</div> <div>.</div> </div>
2	H	235	<div> <div>5%</div> <div>86%</div> <div>6%</div> <div>8%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3564 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 belimumab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	209	Total	C	N	O	S	0	0	0
			1567	975	266	322	4			

- Molecule 2 is a protein called belimumab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	216	Total	C	N	O	S	0	0	0
			1604	1008	278	310	8			

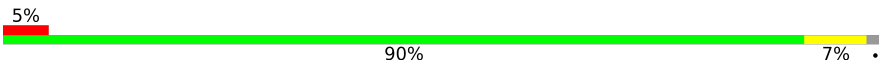
- Molecule 3 is water.

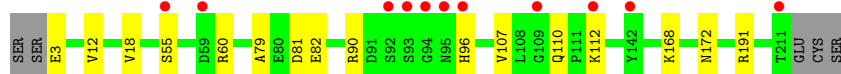
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	196	Total	O	0	0
			196	196		
3	H	197	Total	O	0	0
			197	197		

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.

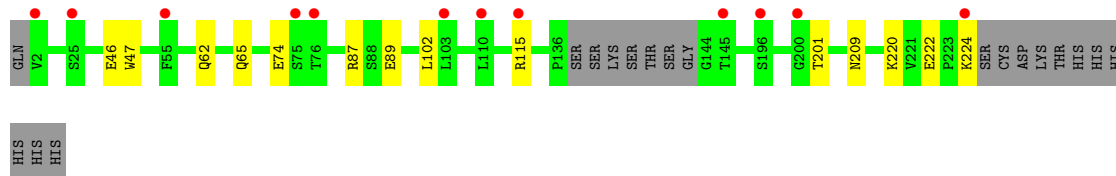
- Molecule 1: belimumab light chain

Chain L: 



- Molecule 2: belimumab heavy chain

Chain H: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.91Å 52.28Å 63.75Å 90.00° 90.72° 90.00°	Depositor
Resolution (Å)	26.31 – 1.90 26.31 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.7 (26.31-1.90) 97.7 (26.31-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.69 (at 1.91Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.167 , 0.206 0.171 , 0.207	Depositor DCC
R_{free} test set	1571 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.094	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 63.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.010 for l,k,-h 0.036 for h,-k,-l 0.024 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3564	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.40% 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, 0.333 respectively for untwinned datasets, and 0.375, 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	L	0.36	0/1605	0.50	0/2193
2	H	0.38	0/1641	0.55	0/2233
All	All	0.37	0/3246	0.53	0/4426

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	L	1567	0	1499	12	0
2	H	1604	0	1584	16	0
3	H	197	0	0	10	0
3	L	196	0	0	3	0
All	All	3564	0	3083	24	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 (24) 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)
2:H:115:ARG:NH1	3:H:303:HOH:O	2.19	0.74

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:191:ARG:NE	3:L:302:HOH:O	2.27	0.66
2:H:87:ARG:NH1	3:H:307:HOH:O	2.31	0.64
2:H:46:GLU:OE1	3:H:302:HOH:O	2.16	0.60
2:H:62:GLN:NE2	2:H:65:GLN:OE1	2.36	0.59
2:H:115:ARG:NH1	3:H:301:HOH:O	2.09	0.57
1:L:79:ALA:O	1:L:172:ASN:ND2	2.36	0.57
2:H:89:GLU:HB3	3:H:386:HOH:O	2.07	0.55
1:L:90:ARG:NH1	2:H:102:LEU:HD21	2.22	0.55
1:L:96:HIS:HB2	2:H:47:TRP:CZ3	2.42	0.54
2:H:74:GLU:HG3	3:H:402:HOH:O	2.07	0.53
1:L:90:ARG:CZ	2:H:102:LEU:HD21	2.40	0.51
2:H:209:ASN:OD1	3:H:304:HOH:O	2.19	0.50
1:L:82:GLU:OE2	1:L:110:GLN:NE2	2.44	0.50
1:L:168:LYS:NZ	3:L:305:HOH:O	2.34	0.48
1:L:3:GLU:HG2	3:L:460:HOH:O	2.13	0.48
1:L:60:ARG:NH1	1:L:81:ASP:OD2	2.40	0.47
2:H:220:LYS:HE3	2:H:222:GLU:HG2	1.98	0.45
2:H:65:GLN:NE2	3:H:313:HOH:O	2.49	0.45
1:L:12:VAL:HB	1:L:18:VAL:HB	2.00	0.44
2:H:201:THR:OG1	3:H:305:HOH:O	2.22	0.43
1:L:107:VAL:O	1:L:110:GLN:NE2	2.49	0.42
1:L:96:HIS:CB	2:H:47:TRP:CZ3	3.03	0.40
2:H:89:GLU:CG	3:H:344:HOH:O	2.69	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	L	207/214 (97%)	203 (98%)	4 (2%)	0	100	100
2	H	212/235 (90%)	206 (97%)	6 (3%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	419/449 (93%)	409 (98%)	10 (2%)	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	L	174/179 (97%)	172 (99%)	2 (1%)	76	75
2	H	179/197 (91%)	178 (99%)	1 (1%)	87	88
All	All	353/376 (94%)	350 (99%)	3 (1%)	83	83

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	55	SER
1	L	112	LYS
2	H	224	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	96	HIS

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.

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	L	209/214 (97%)	0.18	11 (5%) 26 30	16, 27, 51, 73	0
2	H	216/235 (91%)	0.08	12 (5%) 24 28	16, 24, 46, 72	0
All	All	425/449 (94%)	0.13	23 (5%) 26 29	16, 25, 48, 73	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	211	THR	6.0
2	H	200	GLY	5.0
1	L	95	ASN	3.8
1	L	92	SER	3.5
2	H	115	ARG	3.4
2	H	75	SER	3.3
2	H	25	SER	3.2
1	L	96	HIS	3.1
1	L	94	GLY	3.0
2	H	55	PHE	2.9
2	H	224	LYS	2.9
2	H	76	THR	2.9
2	H	110	LEU	2.7
1	L	109	GLY	2.6
1	L	93	SER	2.3
2	H	145	THR	2.3
1	L	59	ASP	2.2
1	L	55	SER	2.2
1	L	112	LYS	2.1
2	H	103	LEU	2.1
2	H	196	SER	2.1
1	L	142	TYR	2.1
2	H	2	VAL	2.0

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.