

wwPDB X-ray Structure Validation Summary Report (i)

Aug 30, 2022 – 05:06 pm BST

PDB ID	:	7PHS
Title	:	highly potent IL6 antagonistic antibody selected from a camelid immune phage
		display repertoire
Authors	:	Cambillau, C.; Pannecoucke, E.; De Haard, H.
Deposited on		
Resolution	:	2.77 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

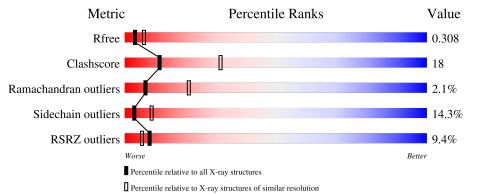
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.30
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.30

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.77 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	4107 (2.80-2.76)
Clashscore	141614	4575 (2.80-2.76)
Ramachandran outliers	138981	4487 (2.80-2.76)
Sidechain outliers	138945	4489 (2.80-2.76)
RSRZ outliers	127900	4027 (2.80-2.76)

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 of 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 <=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	Н	223	65%	27%	6% •
2	L	212	7%	20%	6%



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3325 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 Heavy chain of Fab129D3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Н	223	Total 1667	C 1059	N 269	0 334	${ m S}{ m 5}$	0	0	0

• Molecule 2 is a protein called Light chain of Fab129D3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	211	Total 1569	C 984	N 261	O 319	${ m S}{ m 5}$	0	0	0

• Molecule 3 is water.

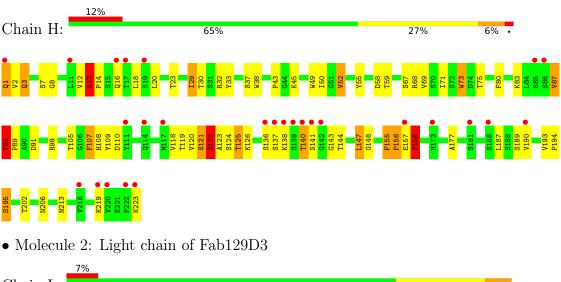
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	Н	51	$\begin{array}{cc} \text{Total} & \text{O} \\ 51 & 51 \end{array}$	0	0
3	L	38	Total O 38 38	0	0



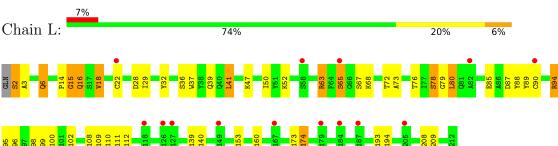


3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Heavy chain of Fab129D3





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants	58.99Å 58.99Å 285.83Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.05 - 2.77	Depositor
Resolution (A)	36.02 - 2.77	EDS
% Data completeness	98.4 (41.05 - 2.77)	Depositor
(in resolution range)	98.8(36.02-2.77)	EDS
R _{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$3.29 (at 2.77 \text{\AA})$	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.273 , 0.303	Depositor
II, II, <i>free</i>	0.286 , 0.308	DCC
R_{free} test set	1344 reflections $(9.91%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	51.9	Xtriage
Anisotropy	0.761	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	(Not available), (Not available)	EDS
L-test for twinning ²	$ < L > = 0.48, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	3325	wwPDB-VP
Average B, all atoms $(Å^2)$	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.37% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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	Bo	nd lengths	Bond angles		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	Н	0.69	3/1714~(0.2%)	0.92	4/2349~(0.2%)	
2	L	0.52	0/1607	0.75	1/2197~(0.0%)	
All	All	0.61	3/3321~(0.1%)	0.84	5/4546~(0.1%)	

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	Н	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	Η	88	THR	C-N	8.76	1.50	1.34
1	Н	13	LYS	C-N	8.26	1.50	1.34
1	Н	155	PHE	C-N	8.20	1.49	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
1	Н	122	SER	C-N-CA	9.71	145.98	121.70
1	Н	58	ASP	CB-CA-C	6.22	122.84	110.40
1	Н	158	PRO	N-CA-CB	-5.24	96.83	102.60
1	Н	107	PHE	C-N-CA	5.21	134.71	121.70
2	L	110	LEU	N-CA-C	5.16	124.92	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	Н	122	SER	Mainchain,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	Н	1667	0	1605	65	0
2	L	1569	0	1524	59	0
3	Н	51	0	0	1	0
3	L	38	0	0	1	0
All	All	3325	0	3129	112	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 18.

The worst 5 of 112 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:H:105:THR:HG22	2:L:52:LYS:NZ	1.60	1.15
1:H:138:LYS:NZ	2:L:209:THR:O	1.79	1.13
2:L:14:PRO:HG3	2:L:111:GLY:HA3	1.24	1.06
1:H:105:THR:HG22	2:L:52:LYS:HZ1	1.16	0.99
2:L:2:SER:N	2:L:94:ARG:HH22	1.69	0.91

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	Н	221/223~(99%)	198 (90%)	17 (8%)	6 (3%)	5 15
2	L	209/212~(99%)	194 (93%)	12~(6%)	3~(1%)	11 31
All	All	430/435~(99%)	392 (91%)	29~(7%)	9~(2%)	7 21

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	Н	43	PRO
1	Н	108	HIS
1	Н	32	ARG
1	Н	195	SER
2	L	174	ASN

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	Н	190/194~(98%)	160 (84%)	30 (16%)	2 7
2	L	174/175~(99%)	152 (87%)	22 (13%)	4 12
All	All	364/369~(99%)	312~(86%)	52 (14%)	3 9

 $5~{\rm of}~52$ residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	Н	213	ASN
2	L	41	LEU
2	L	160	LYS
1	Н	223	LYS
2	L	16	GLN

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such side chains are listed below:

	Chain	\mathbf{Res}	Type
2	L	6	GLN

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Mol	Chain	Res	Type
2	L	16	GLN
2	L	95	ASN
2	L	112	GLN
2	L	192	HIS

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 monosaccharides 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 (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, 95^{th} 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	$\langle RSRZ \rangle$	#RSRZ>2	$\mathbf{OWAB}(\mathbf{A}^2)$	Q<0.9
1	Н	223/223~(100%)	0.83	27 (12%) 4 3	34, 56, 76, 99	10 (4%)
2	L	211/212 (99%)	0.37	14 (6%) 18 13	34, 52, 69, 80	10 (4%)
All	All	434/435~(99%)	0.61	41 (9%) 8 6	34, 53, 74, 99	20 (4%)

The worst 5 of 41 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	Н	137	SER	15.8
1	Н	141	SER	8.4
1	Н	140	THR	7.9
1	Н	181	SER	7.9
1	Н	138	LYS	5.4

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 monosaccharides 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.

