



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 06:28 PM UTC

PDB ID : 5YUP / pdb_00005yup
Title : Crystal Structure of the Fab fragment of FVIIa antibody mAb4F5
Authors : Jiang, L.G.; Persson, E.; Huang, M.D.
Deposited on : 2017-11-22
Resolution : 1.81 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

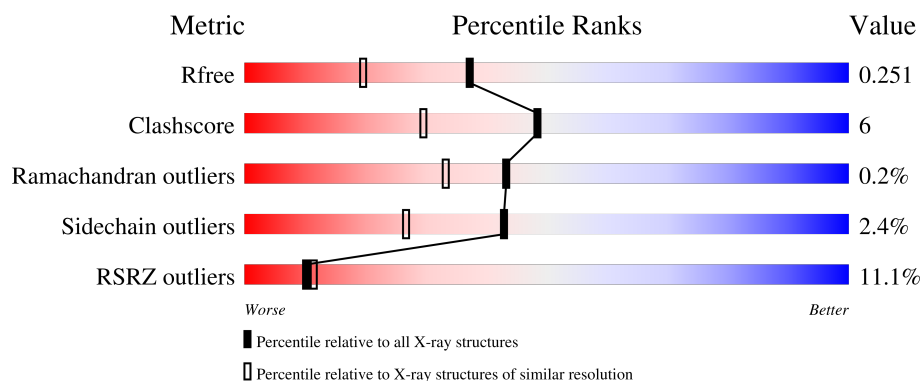
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.81 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}	180053	1112 (1.82-1.82)
Clashscore	190562	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)
RSRZ outliers	180081	1112 (1.82-1.82)

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 $\leq 5\%$. 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	H	219	12% 85% 12% .
2	L	215	10% 87% 11% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3464 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 the heavy chain of the Fab fragment of FVIIa antibody mAb4F5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	219	Total	C	N	O	S	0	0	0
			1658	1047	275	328	8			

- Molecule 2 is a protein called the light chain of the Fab fragment of FVIIa antibody mAb4F5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	215	Total	C	N	O	S	0	0	0
			1675	1042	287	339	7			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	65	Total	O	0	0
			65	65		
3	L	66	Total	O	0	0
			66	66		

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	141.64Å 141.64Å 141.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.82 – 1.81 70.82 – 1.81	Depositor EDS
% Data completeness (in resolution range)	100.0 (70.82-1.81) 100.0 (70.82-1.81)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 1.81Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.207 , 0.249 0.219 , 0.251	Depositor DCC
R_{free} test set	2067 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	32.1	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.032 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3464	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.68% 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

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	$\# Z > 5$	RMSZ	$\# Z > 5$
1	H	0.73	2/1701 (0.1%)	1.05	11/2323 (0.5%)
2	L	0.76	6/1713 (0.4%)	0.96	7/2325 (0.3%)
All	All	0.74	8/3414 (0.2%)	1.01	18/4648 (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	H	0	1
2	L	0	2
All	All	0	3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	31	ILE	CA-C	-7.84	1.41	1.52
2	L	21	ILE	C-O	-5.71	1.18	1.24
1	H	99	ASP	CG-OD2	-5.54	1.14	1.25
2	L	10	PHE	C-O	-5.51	1.17	1.24
1	H	60	ALA	C-O	-5.41	1.17	1.23
2	L	19	VAL	C-O	-5.17	1.18	1.24
2	L	9	LYS	C-O	-5.13	1.18	1.24
2	L	11	LEU	C-O	-5.03	1.18	1.24

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	32	GLY	N-CA-C	-15.94	88.91	110.43
1	H	55	SER	CA-C-N	11.60	141.36	121.64
1	H	55	SER	C-N-CA	11.60	141.36	121.64
2	L	33	LEU	N-CA-C	-8.17	102.89	112.86

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	102	TYR	CB-CA-C	7.69	122.73	111.88
1	H	99	ASP	CA-C-N	6.72	134.57	121.41
1	H	99	ASP	C-N-CA	6.72	134.57	121.41
1	H	54	LEU	CA-C-O	-6.57	111.48	118.77
1	H	99	ASP	CB-CA-C	6.37	123.10	110.42
2	L	214	GLU	N-CA-C	6.36	117.00	108.23
1	H	54	LEU	CA-C-N	6.30	133.57	121.54
1	H	54	LEU	C-N-CA	6.30	133.57	121.54
1	H	164	SER	N-CA-C	-6.27	107.47	114.62
2	L	30	GLY	CA-C-N	5.67	129.60	120.30
2	L	30	GLY	C-N-CA	5.67	129.60	120.30
2	L	31	ILE	O-C-N	5.65	128.19	121.80
2	L	31	ILE	CB-CA-C	-5.65	103.11	112.26
1	H	54	LEU	O-C-N	5.01	127.86	122.34

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	136	THR	Peptide
2	L	30	GLY	Peptide
2	L	31	ILE	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	H	1658	0	1608	23	0
2	L	1675	0	1603	23	0
3	H	65	0	0	0	0
3	L	66	0	0	0	0
All	All	3464	0	3211	36	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 6.

All (36) 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:103:ARG:NH1	2:L:51:TRP:NE1	1.83	1.23
1:H:103:ARG:NH1	2:L:51:TRP:HE1	1.39	1.15
1:H:103:ARG:HH12	2:L:51:TRP:NE1	1.41	1.08
1:H:103:ARG:NH1	2:L:51:TRP:CD1	2.46	0.83
2:L:31:ILE:H	2:L:31:ILE:HD12	1.53	0.72
1:H:103:ARG:HH11	2:L:51:TRP:NE1	1.86	0.70
2:L:31:ILE:HD12	2:L:31:ILE:N	2.08	0.67
2:L:31:ILE:H	2:L:31:ILE:CD1	2.13	0.61
1:H:103:ARG:HH12	2:L:51:TRP:HE1	0.67	0.61
1:H:12:VAL:HG11	1:H:85:LEU:HD13	1.83	0.60
2:L:84:LEU:HD13	2:L:169:SER:HA	1.86	0.58
2:L:146:ASN:HB3	2:L:198:THR:HG23	1.86	0.58
1:H:104:VAL:HG23	1:H:107:TRP:NE1	2.21	0.56
1:H:11:LEU:HB2	1:H:151:PRO:HG3	1.89	0.55
1:H:103:ARG:HD2	2:L:50:HIS:HB2	1.89	0.53
1:H:3:LYS:HB2	1:H:25:SER:HB2	1.90	0.53
2:L:32:GLY:C	2:L:34:VAL:N	2.65	0.52
1:H:62:THR:HG23	1:H:63:VAL:HG13	1.92	0.51
1:H:102:TYR:CD2	2:L:95:TYR:OH	2.63	0.51
2:L:34:VAL:HG11	2:L:72:PHE:CD2	2.46	0.51
1:H:175:GLN:HG2	1:H:176:SER:OG	2.11	0.50
1:H:200:ASN:ND2	1:H:211:ASP:OD1	2.35	0.49
1:H:123:PRO:HB3	1:H:149:TYR:HB3	1.95	0.48
2:L:32:GLY:C	2:L:34:VAL:H	2.22	0.48
2:L:187:TYR:CZ	2:L:212:ARG:HD3	2.49	0.48
1:H:104:VAL:HG22	2:L:37:TYR:OH	2.13	0.47
2:L:31:ILE:N	2:L:31:ILE:CD1	2.76	0.47
1:H:6:GLU:HG3	1:H:95:CYS:SG	2.55	0.47
2:L:146:ASN:HB3	2:L:198:THR:CG2	2.44	0.46
1:H:104:VAL:HG23	1:H:107:TRP:HE1	1.81	0.45
1:H:30:SER:O	1:H:53:ASN:HB2	2.19	0.42
1:H:60:ALA:O	1:H:64:THR:HG23	2.20	0.42
1:H:18:ARG:HB3	1:H:82:MET:HE3	2.02	0.41
1:H:103:ARG:HD2	2:L:50:HIS:CB	2.50	0.41
2:L:29:VAL:O	2:L:33:LEU:HB2	2.19	0.41
2:L:13:THR:OG1	2:L:17:ASN:HB2	2.20	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 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	H	217/219 (99%)	213 (98%)	3 (1%)	1 (0%)	24	14
2	L	213/215 (99%)	208 (98%)	5 (2%)	0	100	100
All	All	430/434 (99%)	421 (98%)	8 (2%)	1 (0%)	43	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	55	SER

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	H	186/186 (100%)	182 (98%)	4 (2%)	45	30
2	L	193/193 (100%)	188 (97%)	5 (3%)	40	23
All	All	379/379 (100%)	370 (98%)	9 (2%)	43	27

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	55	SER
1	H	136	THR
1	H	154	VAL
1	H	176	SER
2	L	31	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	L	157	GLN
2	L	161	LEU
2	L	198	THR
2	L	215	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	L	8	HIS
2	L	78	ASN
2	L	138	ASN
2	L	211	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 oligosaccharides 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	H	219/219 (100%)	0.86	27 (12%) 8 9	23, 35, 50, 98	0
2	L	215/215 (100%)	0.70	21 (9%) 13 15	23, 33, 46, 86	0
All	All	434/434 (100%)	0.79	48 (11%) 10 11	23, 34, 49, 98	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	31	ILE	8.5
2	L	33	LEU	7.9
1	H	136	THR	4.4
1	H	102	TYR	4.3
1	H	100	GLY	4.2
1	H	103	ARG	4.2
1	H	55	SER	4.0
1	H	219	CYS	4.0
2	L	15	VAL	3.8
2	L	16	GLY	3.6
2	L	164	TRP	3.5
1	H	209	LYS	3.3
2	L	77	ASN	3.3
2	L	215	CYS	3.1
2	L	203	THR	3.1
2	L	214	GLU	2.9
2	L	42	GLY	2.7
1	H	177	ASP	2.7
2	L	10	PHE	2.5
1	H	217	ARG	2.5
1	H	218	ASP	2.5
2	L	78	ASN	2.5
1	H	41	PRO	2.5
1	H	32	TYR	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	H	99	ASP	2.4
1	H	215	VAL	2.3
2	L	195	CYS	2.3
1	H	106	TYR	2.3
1	H	31	ASP	2.3
1	H	210	VAL	2.3
2	L	198	THR	2.3
1	H	135	GLN	2.3
1	H	137	ASN	2.1
2	L	34	VAL	2.1
2	L	57	THR	2.1
1	H	7	SER	2.1
1	H	133	ALA	2.1
1	H	104	VAL	2.1
2	L	160	VAL	2.1
2	L	32	GLY	2.1
1	H	204	PRO	2.1
1	H	174	LEU	2.1
1	H	162	SER	2.0
2	L	169	SER	2.0
2	L	204	SER	2.0
1	H	196	THR	2.0
1	H	165	SER	2.0
2	L	79	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 oligosaccharides 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.