



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:14 AM UTC

PDB ID : 7KET / pdb_00007ket
Title : Factor H enhancing human antibody fragment (Fab) to meningococcal Factor H binding protein
Authors : Beernink, P.T.; Sands, N.
Deposited on : 2020-10-12
Resolution : 2.00 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

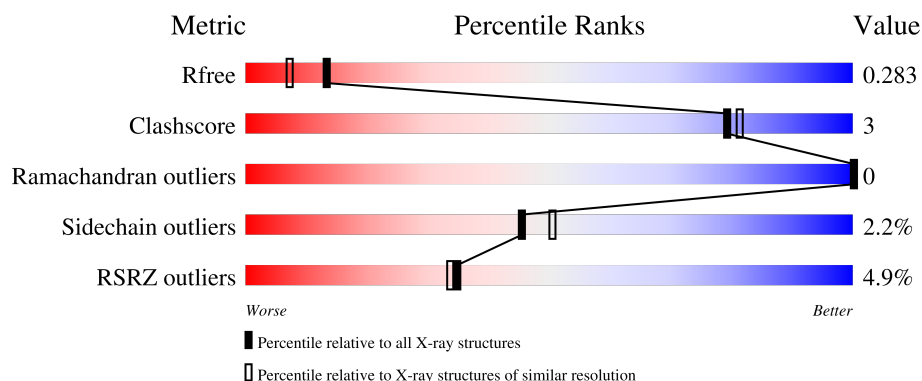
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	B	217	 6% 84% 11% . .
2	A	235	 5% 84% 7% . 8%
3	C	263	 3% 86% . . 8%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10263 atoms, of which 5014 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 Immunoglobulin gamma, kappa light chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	209	Total	C	H	N	O	S	0	1	0
			3160	1008	1551	274	322	5			

- Molecule 2 is a protein called Immunoglobulin gamma, heavy chain Fd fragment.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	A	217	Total	C	H	N	O	S	0	4	0
			3260	1031	1623	276	320	10			

- Molecule 3 is a protein called Factor H binding protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	242	Total	C	H	N	O	S	0	2	0
			3681	1145	1840	330	365	1			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	256	LEU	-	expression tag	UNP Q6VRZ6
C	257	GLU	-	expression tag	UNP Q6VRZ6
C	258	HIS	-	expression tag	UNP Q6VRZ6
C	259	HIS	-	expression tag	UNP Q6VRZ6
C	260	HIS	-	expression tag	UNP Q6VRZ6
C	261	HIS	-	expression tag	UNP Q6VRZ6
C	262	HIS	-	expression tag	UNP Q6VRZ6
C	263	HIS	-	expression tag	UNP Q6VRZ6

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	30	Total	O	0	0
			30	30		

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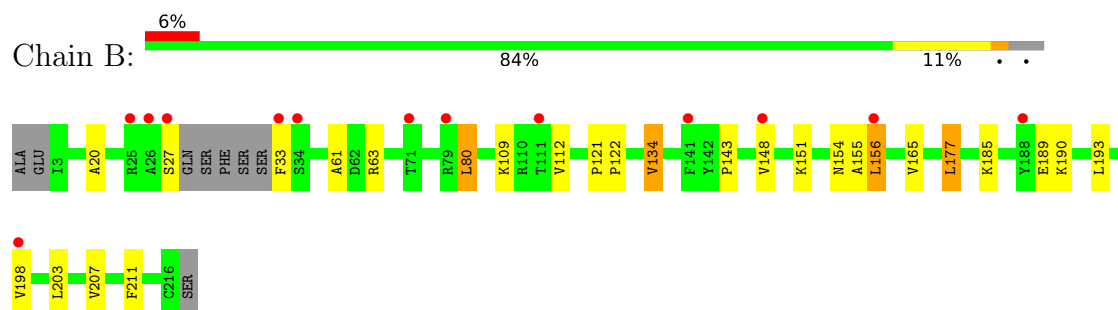
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	67	Total 67	O 67	0	0
4	C	65	Total 65	O 65	0	0

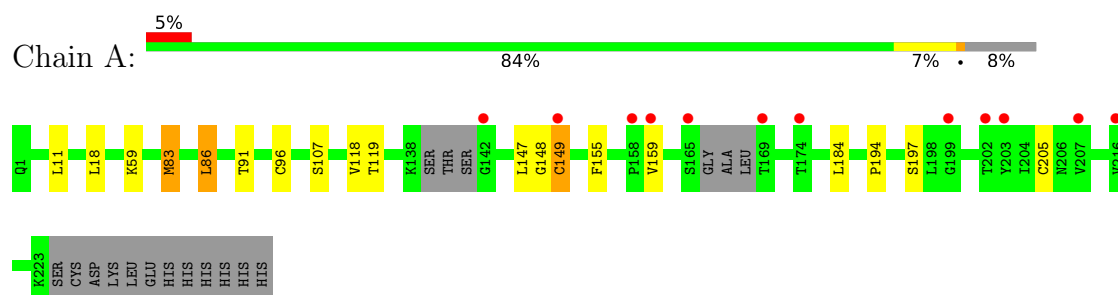
3 Residue-property plots [i](#)

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.

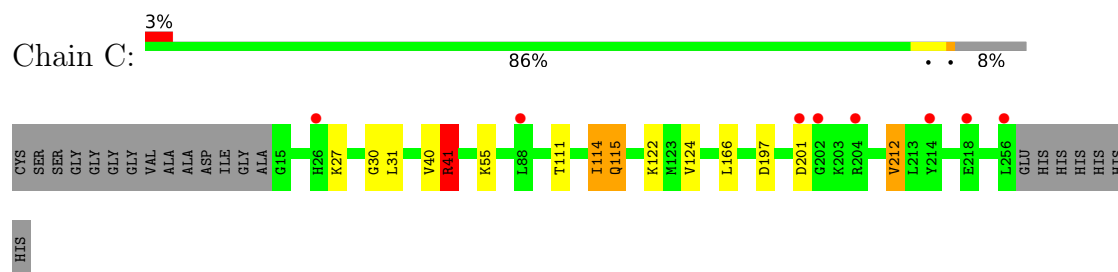
- Molecule 1: Immunoglobulin gamma, kappa light chain



- Molecule 2: Immunoglobulin gamma, heavy chain Fd fragment



- Molecule 3: Factor H binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	38.58Å 86.03Å 107.89Å 90.00° 96.91° 90.00°	Depositor
Resolution (Å)	107.11 – 2.00 107.11 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.0 (107.11-2.00) 95.0 (107.11-2.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.02 (at 2.00Å)	Xtriage
Refinement program	Coot 1.18.2_3874, PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.234 , 0.281 0.242 , 0.283	Depositor DCC
R_{free} test set	1571 reflections (2.68%)	wwPDB-VP
Wilson B-factor (Å ²)	45.3	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10263	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.15% 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	B	0.53	2/1646 (0.1%)	0.75	2/2233 (0.1%)
2	A	0.64	6/1685 (0.4%)	0.71	2/2290 (0.1%)
3	C	0.62	5/1872 (0.3%)	0.77	6/2512 (0.2%)
All	All	0.60	13/5203 (0.2%)	0.74	10/7035 (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	B	0	4
3	C	0	1
All	All	0	5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	115[A]	GLN	C-O	9.56	1.35	1.23
3	C	115[B]	GLN	C-O	9.56	1.35	1.23
1	B	134	VAL	C-O	-8.54	1.14	1.24
2	A	18	LEU	C-O	-7.92	1.14	1.23
2	A	184	LEU	C-O	-7.51	1.14	1.23
3	C	124	VAL	C-O	-7.15	1.17	1.24
2	A	86	LEU	C-O	-7.11	1.15	1.24
2	A	148	GLY	C-O	-6.88	1.15	1.23
1	B	177	LEU	C-O	-6.57	1.15	1.23
3	C	55	LYS	C-O	-5.73	1.16	1.23
3	C	114	ILE	C-O	-5.23	1.18	1.23
2	A	83	MET	C-O	-5.19	1.18	1.24
2	A	107	SER	CA-CB	-5.07	1.46	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	184	LEU	CA-C-O	-6.42	113.79	121.56
3	C	27	LYS	N-CA-C	-5.50	106.57	113.28
1	B	155	ALA	CB-CA-C	-5.44	109.23	117.07
3	C	212	VAL	N-CA-C	-5.34	100.75	108.71
3	C	201	ASP	CA-CB-CG	5.22	117.82	112.60
3	C	41	ARG	CB-CA-C	5.21	118.95	109.83
3	C	115[A]	GLN	O-C-N	-5.07	116.67	122.95
3	C	115[B]	GLN	O-C-N	-5.07	116.67	122.95
2	A	147	LEU	N-CA-C	-5.05	101.73	108.86
1	B	177	LEU	CA-C-O	-5.01	115.45	121.11

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	177	LEU	Mainchain
1	B	33	PHE	Peptide
1	B	80	LEU	Mainchain
3	C	166	LEU	Mainchain

5.2 Too-close contacts ⓘ

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	B	1609	1551	1568	10	1
2	A	1637	1623	1628	11	0
3	C	1841	1840	1840	5	1
4	A	67	0	0	0	0
4	B	30	0	0	1	0
4	C	65	0	0	0	0
All	All	5249	5014	5036	26	1

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 3.

All (26) 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:A:149[A]:CYS:HG	2:A:205:CYS:HG	1.41	0.68
2:A:11:LEU:CD1	2:A:155:PHE:HE2	2.11	0.64
2:A:91:THR:HG23	2:A:119:THR:HA	1.82	0.60
2:A:11:LEU:HD11	2:A:155:PHE:HE2	1.71	0.55
3:C:111[B]:THR:HG21	3:C:114:ILE:CG2	2.40	0.52
3:C:111[B]:THR:HG21	3:C:114:ILE:HG23	1.92	0.51
2:A:149[B]:CYS:HB3	2:A:205:CYS:HG	1.76	0.51
3:C:115[A]:GLN:NE2	3:C:122:LYS:O	2.45	0.50
2:A:83:MET:HB3	2:A:86:LEU:HD21	1.94	0.49
2:A:194:PRO:HG2	2:A:197:SER:OG	2.13	0.49
1:B:20:ALA:HB2	1:B:80:LEU:HD11	1.96	0.48
1:B:122:PRO:HD3	1:B:134:VAL:HG22	1.98	0.45
3:C:30:GLY:HA2	3:C:31:LEU:C	2.42	0.45
3:C:41:ARG:HA	3:C:41:ARG:HD2	1.52	0.43
2:A:149[B]:CYS:HB3	2:A:205:CYS:SG	2.59	0.43
2:A:11:LEU:CD1	2:A:155:PHE:CE2	2.98	0.42
1:B:112:VAL:HG22	1:B:143:PRO:HD3	2.01	0.41
1:B:203:LEU:HD13	1:B:207:VAL:HG23	2.02	0.41
1:B:109:LYS:NZ	4:B:306:HOH:O	2.53	0.41
1:B:148:VAL:HG22	1:B:198:VAL:HG12	2.01	0.41
1:B:121:PRO:HB3	1:B:211:PHE:CE1	2.56	0.41
2:A:11:LEU:HD13	2:A:155:PHE:CE2	2.56	0.41
1:B:185:LYS:HE3	1:B:189:GLU:OE1	2.21	0.41
2:A:83:MET:HE1	2:A:118:VAL:HG21	2.03	0.40
1:B:61:ALA:HB3	1:B:63:ARG:HG3	2.04	0.40
1:B:151:LYS:HE2	1:B:156:LEU:HG	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:LYS:HZ1	3:C:197:ASP:OD1[1_656]	1.50	0.10

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	B	206/217 (95%)	194 (94%)	12 (6%)	0	100	100
2	A	215/235 (92%)	209 (97%)	6 (3%)	0	100	100
3	C	242/263 (92%)	229 (95%)	13 (5%)	0	100	100
All	All	663/715 (93%)	632 (95%)	31 (5%)	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	B	181/187 (97%)	176 (97%)	5 (3%)	38	41
2	A	188/200 (94%)	182 (97%)	6 (3%)	34	35
3	C	191/202 (95%)	188 (98%)	3 (2%)	55	62
All	All	560/589 (95%)	546 (98%)	14 (2%)	45	45

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	27	SER
1	B	154	ASN
1	B	156	LEU
1	B	165	VAL
1	B	193	LEU
2	A	59	LYS
2	A	96[A]	CYS
2	A	96[B]	CYS
2	A	149[A]	CYS
2	A	149[B]	CYS
2	A	159	VAL
3	C	40	VAL
3	C	41	ARG

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Mol	Chain	Res	Type
3	C	212	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	173	HIS
3	C	119	HIS
3	C	138	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 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	B	209/217 (96%)	0.86	13 (6%)	26 25	45, 75, 109, 139	1 (0%)
2	A	217/235 (92%)	0.59	12 (5%)	30 29	30, 61, 100, 138	4 (1%)
3	C	242/263 (92%)	0.43	8 (3%)	49 48	28, 59, 97, 133	2 (0%)
All	All	668/715 (93%)	0.62	33 (4%)	35 34	28, 64, 104, 139	7 (1%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	27	SER	5.3
1	B	33	PHE	3.8
2	A	158	PRO	3.6
3	C	214	TYR	3.4
1	B	26	ALA	3.4
3	C	88	LEU	3.3
3	C	256	LEU	3.3
2	A	149[A]	CYS	3.1
2	A	216	VAL	3.0
3	C	218	GLU	2.8
1	B	156	LEU	2.8
1	B	34[A]	SER	2.8
1	B	148	VAL	2.7
1	B	79	ARG	2.6
2	A	174	THR	2.6
2	A	202	THR	2.6
2	A	199	GLY	2.5
2	A	159	VAL	2.5
1	B	198	VAL	2.4
2	A	169	THR	2.4
2	A	207	VAL	2.3
3	C	201	ASP	2.2
2	A	203	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
3	C	204	ARG	2.1
2	A	165	SER	2.1
1	B	25	ARG	2.1
1	B	111	THR	2.1
1	B	188	TYR	2.1
1	B	71	THR	2.0
3	C	202	GLY	2.0
1	B	141	PHE	2.0
3	C	26	HIS	2.0
2	A	142	GLY	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.