



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 6, 2021 – 12:08 PM JST

PDB ID : 7CHY  
Title : Crystal Structure Of Human Il-1beta In Complex With Antibody Binding  
Fragment Of IgG26  
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Deposited on : 2020-07-06  
Resolution : 2.65 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](mailto: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.

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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	:	2.16
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.16

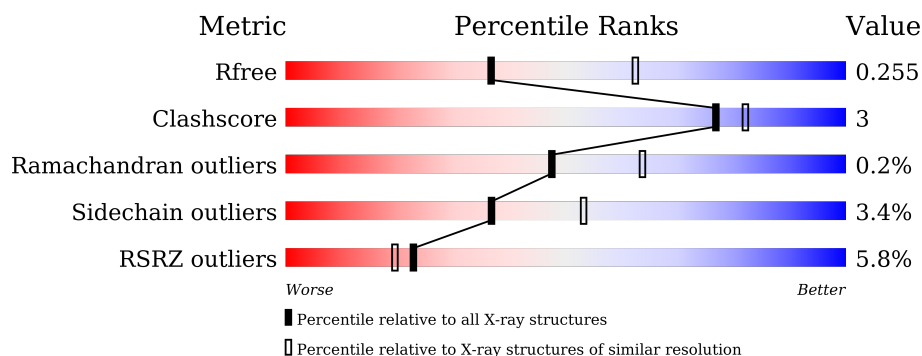
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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))
$R_{free}$	130704	1332 (2.68-2.64)
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

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  $\geq 3$ , 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	L	214	<div> <div>4%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>
2	H	234	<div> <div>%</div> <div>82%</div> <div>9%</div> <div>9%</div> </div>
3	I	157	<div> <div>13%</div> <div>85%</div> <div>10%</div> <div>.</div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4707 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 light chain of antibody binding fragment of IgG26.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	211	Total	C	N	O	S	0	0	0
			1625	1021	270	329	5			

- Molecule 2 is a protein called heavy chain of antibody binding fragment of IgG26.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	212	Total	C	N	O	S	0	0	0
			1603	1024	267	306	6			

- Molecule 3 is a protein called Interleukin-1 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	150	Total	C	N	O	S	0	0	0
			1200	762	198	232	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	113	LEU	-	expression tag	UNP P01584
I	114	GLY	-	expression tag	UNP P01584
I	115	SER	-	expression tag	UNP P01584
I	116	ARG	-	expression tag	UNP P01584

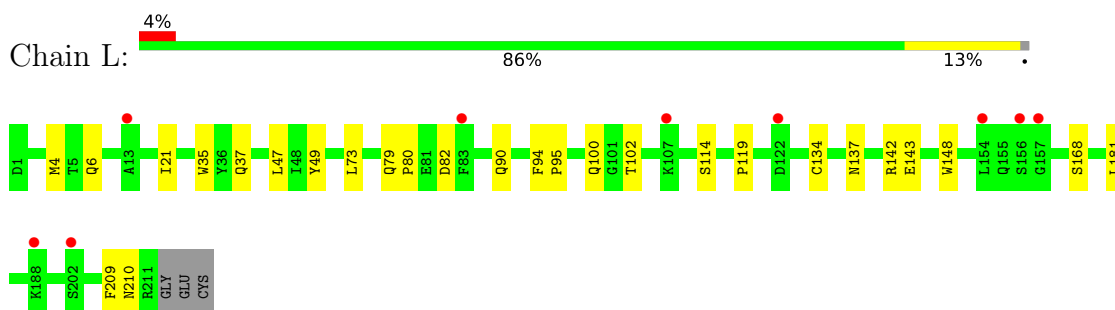
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	63	Total	O	0	0
			63	63		
4	H	172	Total	O	0	0
			172	172		
4	I	44	Total	O	0	0
			44	44		

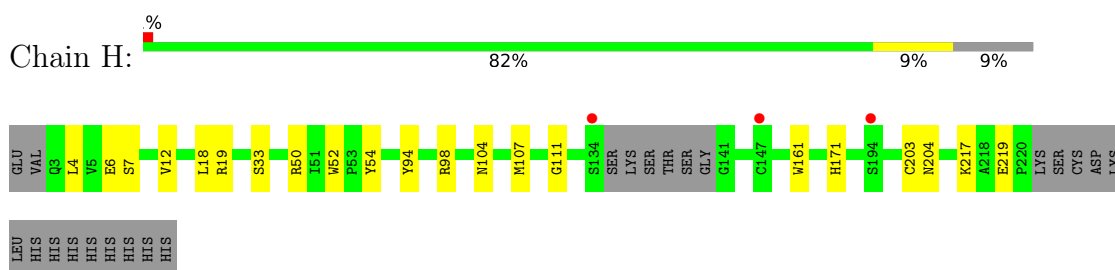
### 3 Residue-property plots [i](#)

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.

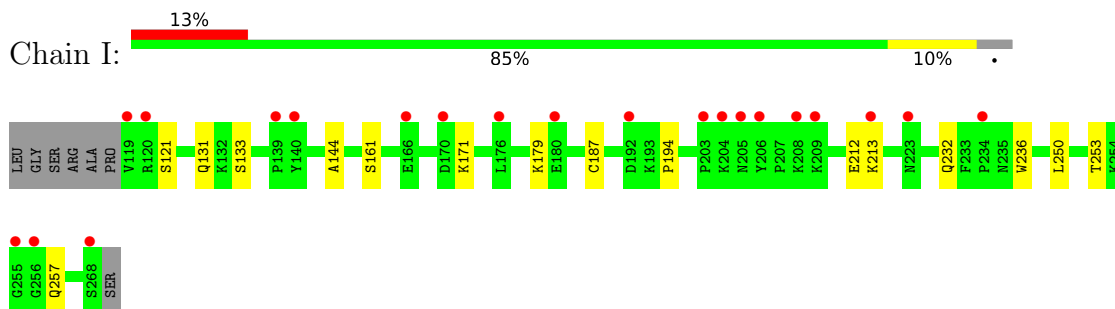
- Molecule 1: light chain of antibody binding fragment of IgG26



- Molecule 2: heavy chain of antibody binding fragment of IgG26



- Molecule 3: Interleukin-1 beta



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	156.34Å 112.24Å 38.58Å 90.00° 94.85° 90.00°	Depositor
Resolution (Å)	19.70 – 2.65 19.75 – 2.65	Depositor EDS
% Data completeness (in resolution range)	93.7 (19.70-2.65) 93.9 (19.75-2.65)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.34 (at 2.67Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.200 , 0.254 0.201 , 0.255	Depositor DCC
$R_{free}$ test set	963 reflections (5.32%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.7	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 42.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4707	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	L	0.64	0/1662	0.74	0/2259
2	H	0.63	0/1648	0.76	0/2250
3	I	0.65	0/1222	0.74	0/1643
All	All	0.64	0/4532	0.75	0/6152

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

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	1625	0	1577	13	0
2	H	1603	0	1549	9	0
3	I	1200	0	1198	6	0
4	H	172	0	0	1	0
4	I	44	0	0	0	0
4	L	63	0	0	0	0
All	All	4707	0	4324	26	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 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)
3:I:187:CYS:SG	3:I:250:LEU:HD11	2.40	0.61
2:H:161:TRP:CH2	2:H:203:CYS:HB3	2.44	0.53
1:L:4:MET:HE3	1:L:90:GLN:HG2	1.92	0.51
1:L:35:TRP:CD2	1:L:73:LEU:HB2	2.46	0.51
3:I:213:LYS:HB3	3:I:232:GLN:HE21	1.74	0.51
3:I:133:SER:O	3:I:144:ALA:HA	2.14	0.48
2:H:12:VAL:HG11	2:H:18:LEU:HD13	1.95	0.48
1:L:37:GLN:HB2	1:L:47:LEU:HD11	1.95	0.48
3:I:212:GLU:OE1	3:I:212:GLU:N	2.48	0.47
1:L:137:ASN:HD21	2:H:171:HIS:CD2	2.32	0.47
2:H:4:LEU:O	2:H:111:GLY:HA2	2.14	0.47
3:I:194:PRO:HG2	3:I:236:TRP:CD2	2.51	0.46
1:L:6:GLN:NE2	1:L:102:THR:HG23	2.31	0.45
1:L:6:GLN:O	1:L:100:GLN:NE2	2.50	0.45
2:H:19:ARG:HD2	4:H:325:HOH:O	2.16	0.45
1:L:21:ILE:HG23	1:L:102:THR:HG21	1.98	0.45
2:H:6:GLU:OE2	2:H:94:TYR:O	2.34	0.45
1:L:49:TYR:CZ	2:H:104:ASN:HB3	2.53	0.44
3:I:121:SER:HB3	3:I:161:SER:HA	1.99	0.44
1:L:94:PHE:HA	1:L:95:PRO:C	2.37	0.43
1:L:134:CYS:HB2	1:L:148:TRP:CH2	2.54	0.42
1:L:21:ILE:HD12	1:L:73:LEU:HD23	2.02	0.42
2:H:98:ARG:O	2:H:107:MET:HA	2.19	0.41
1:L:119:PRO:HB3	1:L:209:PHE:CE2	2.55	0.41
1:L:79:GLN:HB3	1:L:80:PRO:HD2	2.02	0.41
2:H:33:SER:HB3	2:H:52:TRP:CE2	2.55	0.41

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	209/214 (98%)	201 (96%)	8 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	208/234 (89%)	199 (96%)	8 (4%)	1 (0%)	29	43
3	I	148/157 (94%)	138 (93%)	10 (7%)	0	100	100
All	All	565/605 (93%)	538 (95%)	26 (5%)	1 (0%)	47	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	54	TYR

### 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	L	186/188 (99%)	179 (96%)	7 (4%)	33	49
2	H	173/194 (89%)	168 (97%)	5 (3%)	42	60
3	I	138/143 (96%)	133 (96%)	5 (4%)	35	51
All	All	497/525 (95%)	480 (97%)	17 (3%)	37	53

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	82	ASP
1	L	114	SER
1	L	142	ARG
1	L	143	GLU
1	L	168	SER
1	L	181	LEU
1	L	210	ASN
2	H	7	SER
2	H	50	ARG
2	H	204	ASN
2	H	217	LYS
2	H	219	GLU
3	I	131	GLN

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Mol	Chain	Res	Type
3	I	171	LYS
3	I	179	LYS
3	I	253	THR
3	I	257	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	124	GLN
1	L	137	ASN
1	L	138	ASN
1	L	210	ASN
3	I	232	GLN
3	I	257	GLN

### 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

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	L	211/214 (98%)	0.38	9 (4%) 35 31	36, 86, 127, 142	0
2	H	212/234 (90%)	-0.31	3 (1%) 75 73	25, 37, 91, 107	0
3	I	150/157 (95%)	0.64	21 (14%) 2 2	36, 83, 138, 151	0
All	All	573/605 (94%)	0.19	33 (5%) 23 19	25, 69, 127, 151	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	83	PHE	6.9
3	I	170	ASP	4.9
2	H	134	SER	4.2
3	I	176	LEU	3.6
3	I	166	GLU	3.5
3	I	204	LYS	3.4
3	I	268	SER	3.4
3	I	206	TYR	3.3
3	I	139	PRO	3.1
3	I	180	GLU	2.9
3	I	255	GLY	2.8
3	I	223	ASN	2.8
1	L	122	ASP	2.7
3	I	203	PRO	2.6
1	L	156	SER	2.6
1	L	202	SER	2.6
1	L	188	LYS	2.6
1	L	13	ALA	2.5
3	I	208	LYS	2.5
3	I	209	LYS	2.4
1	L	107	LYS	2.3
3	I	119	VAL	2.3
3	I	234	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
2	H	147	CYS	2.3
3	I	213	LYS	2.2
1	L	157	GLY	2.2
3	I	256	GLY	2.2
1	L	154	LEU	2.2
3	I	140	TYR	2.2
3	I	120	ARG	2.1
3	I	205	ASN	2.1
2	H	194	SER	2.1
3	I	192	ASP	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 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.