



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2023 – 12:17 PM EDT

PDB ID : 8F6O  
Title : anti-BTLA monoclonal antibody h22B3 in complex with BTLA  
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Deposited on : 2022-11-16  
Resolution : 2.31 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.

The types of validation reports are described at

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

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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  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtrriage (Phenix) : 1.13  
EDS : 2.35  
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) : 2.35

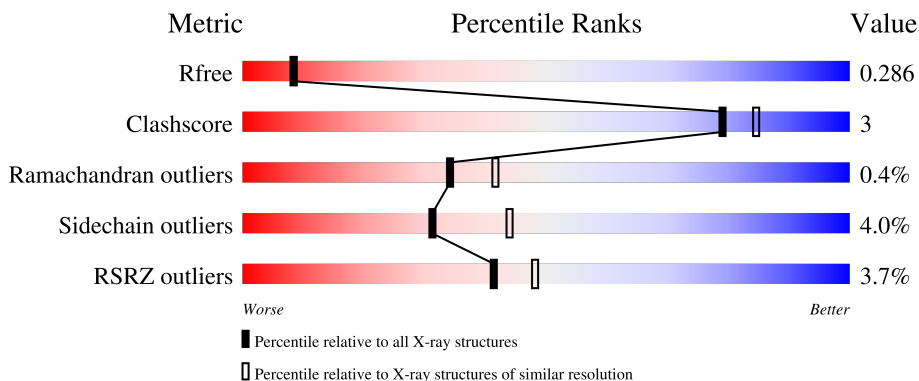
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


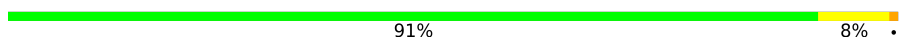

The reported resolution of this entry is 2.31 Å.

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	5974 (2.34-2.30)
Clashscore	141614	6604 (2.34-2.30)
Ramachandran outliers	138981	6523 (2.34-2.30)
Sidechain outliers	138945	6523 (2.34-2.30)
RSRZ outliers	127900	5855 (2.34-2.30)

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	A	232	 87% 9% . .
2	B	217	 91% 8% .
3	C	130	 13% 68% 9% 22%

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 4258 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 h22B3 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	224	1687	1064	274	341	8	0	1	0

- Molecule 2 is a protein called h22B3 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	217	1644	1027	274	338	5	0	1	0

- Molecule 3 is a protein called B- and T-lymphocyte attenuator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	101	769	486	131	146	6	0	1	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	151	GLU	-	expression tag	UNP Q7Z6A9
C	152	PHE	-	expression tag	UNP Q7Z6A9
C	153	HIS	-	expression tag	UNP Q7Z6A9
C	154	HIS	-	expression tag	UNP Q7Z6A9
C	155	HIS	-	expression tag	UNP Q7Z6A9
C	156	HIS	-	expression tag	UNP Q7Z6A9
C	157	HIS	-	expression tag	UNP Q7Z6A9
C	158	HIS	-	expression tag	UNP Q7Z6A9
C	159	HIS	-	expression tag	UNP Q7Z6A9
C	160	HIS	-	expression tag	UNP Q7Z6A9

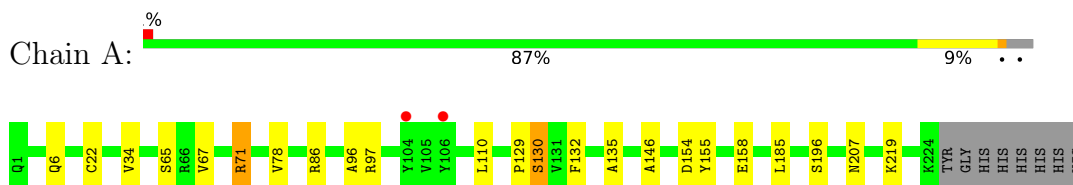
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	92	Total O 93 93	0	1
4	B	62	Total O 63 63	0	1
4	C	2	Total O 2 2	0	0

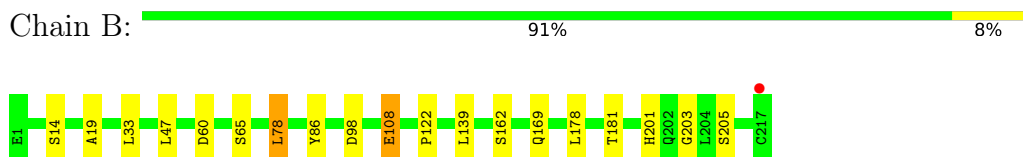
### 3 Residue-property plots [i](#)

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.

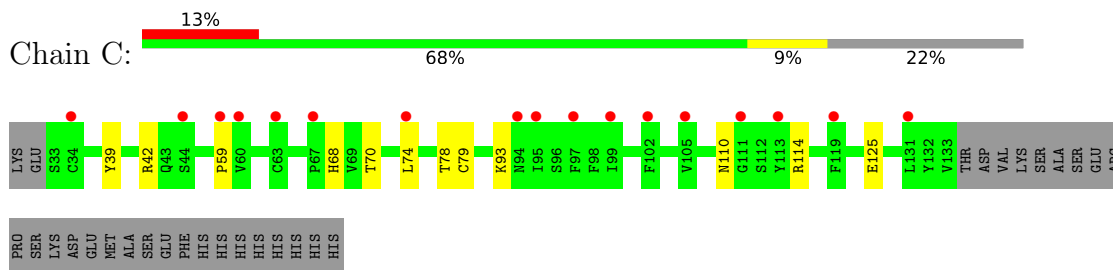
- Molecule 1: h22B3 Fab heavy chain



- Molecule 2: h22B3 Fab light chain



- Molecule 3: B- and T-lymphocyte attenuator



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	68.95Å 88.06Å 233.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.00 – 2.31 54.29 – 2.31	Depositor EDS
% Data completeness (in resolution range)	99.9 (54.00-2.31) 99.9 (54.29-2.31)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.42 (at 2.32Å)	Xtrriage
Refinement program	BUSTER v2.11.6	Depositor
R, $R_{free}$	0.215 , 0.266 0.233 , 0.286	Depositor DCC
$R_{free}$ test set	1555 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.6	Xtrriage
Anisotropy	0.902	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 60.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4258	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PCA

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	A	0.55	0/1723	0.78	0/2353
2	B	0.50	0/1681	0.75	0/2286
3	C	0.43	0/792	0.69	0/1081
All	All	0.51	0/4196	0.75	0/5720

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	A	1687	0	1635	11	0
2	B	1644	0	1585	8	0
3	C	769	0	685	4	0
4	A	93	0	0	1	0
4	B	63	0	0	0	0
4	C	2	0	0	0	0
All	All	4258	0	3905	22	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 (22) 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:A:97:ARG:O	1:A:110:LEU:HA	2.07	0.54
2:B:201[B]:HIS:CD2	2:B:203:GLY:H	2.26	0.54
1:A:71:ARG:HB2	1:A:78:VAL:HG12	1.89	0.53
1:A:154:ASP:HB3	1:A:185:LEU:HD13	1.95	0.49
1:A:34:VAL:HG11	1:A:78:VAL:HG21	1.95	0.49
1:A:129:PRO:HB3	1:A:155:TYR:HB3	1.96	0.47
1:A:96:ALA:HB1	1:A:110:LEU:HB3	1.95	0.47
1:A:22:CYS:HB3	1:A:78:VAL:HG23	1.97	0.47
1:A:146:ALA:HB2	1:A:196:SER:HB3	1.98	0.46
2:B:162:SER:HA	2:B:181:THR:O	2.16	0.46
1:A:158:GLU:HB2	4:A:326:HOH:O	2.16	0.45
1:A:135:ALA:O	2:B:122:PRO:HD2	2.17	0.44
3:C:39:TYR:HB2	3:C:59:PRO:HB2	1.98	0.44
3:C:114:ARG:HD3	3:C:125:GLU:HB3	2.00	0.43
2:B:108:GLU:HB2	2:B:169:GLN:OE1	2.19	0.43
2:B:47:LEU:HD11	2:B:86:TYR:HE2	1.83	0.42
2:B:139:LEU:HB2	2:B:178:LEU:HB3	2.03	0.41
3:C:74:LEU:HD23	3:C:79:CYS:SG	2.61	0.41
2:B:19:ALA:HB2	2:B:78:LEU:HD11	2.03	0.41
3:C:68:HIS:CE1	3:C:70:THR:HG23	2.57	0.40
1:A:130:SER:HB3	1:A:132:PHE:CZ	2.56	0.40
2:B:47:LEU:HD11	2:B:86:TYR:CE2	2.56	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	A	223/232 (96%)	220 (99%)	3 (1%)	0	100 100
2	B	216/217 (100%)	206 (95%)	9 (4%)	1 (0%)	29 35

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	100/130 (77%)	87 (87%)	12 (12%)	1 (1%)	15	17
All	All	539/579 (93%)	513 (95%)	24 (4%)	2 (0%)	34	41

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	93	LYS
2	B	78	LEU

### 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	A	190/197 (96%)	182 (96%)	8 (4%)	30	41
2	B	185/186 (100%)	178 (96%)	7 (4%)	33	46
3	C	81/122 (66%)	78 (96%)	3 (4%)	34	47
All	All	456/505 (90%)	438 (96%)	18 (4%)	31	45

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	65	SER
1	A	67	VAL
1	A	71	ARG
1	A	86	ARG
1	A	130	SER
1	A	207	ASN
1	A	219	LYS
2	B	14	SER
2	B	33	LEU
2	B	60	ASP
2	B	65	SER
2	B	98	ASP

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Mol	Chain	Res	Type
2	B	108	GLU
2	B	205	SER
3	C	42	ARG
3	C	78	THR
3	C	110	ASN

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

Mol	Chain	Res	Type
1	A	43	GLN
2	B	150	GLN
3	C	68	HIS
3	C	110	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](#)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
1	PCA	A	1	1	7,8,9	0.93	0	9,10,12	0.82	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	PCA	A	1	1	-	0/0/11/13	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	A	223/232 (96%)	0.14	2 (0%) 84 88	16, 33, 54, 91	0
2	B	217/217 (100%)	0.22	1 (0%) 91 94	21, 41, 59, 68	0
3	C	101/130 (77%)	1.18	17 (16%) 1 2	47, 67, 82, 92	0
All	All	541/579 (93%)	0.37	20 (3%) 41 48	16, 40, 76, 92	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	119	PHE	4.9
3	C	113	TYR	3.8
3	C	102	PHE	3.7
3	C	111	GLY	3.6
3	C	67	PRO	3.1
3	C	60	VAL	3.0
3	C	131	LEU	2.9
3	C	34	CYS	2.7
3	C	97	PHE	2.6
3	C	59	PRO	2.4
3	C	95	ILE	2.4
3	C	94	ASN	2.4
3	C	105	VAL	2.4
3	C	74	LEU	2.4
2	B	217	CYS	2.4
3	C	63	CYS	2.3
1	A	106	TYR	2.2
3	C	44	SER	2.2
1	A	104	TYR	2.1
3	C	99	ILE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
1	PCA	A	1	8/9	0.81	0.25	60,63,64,64	0

## 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.