



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 04:48 am BST

PDB ID : 5U4R  
Title : Crystal structure of the broadly neutralizing Influenza A antibody VRC 315 53-1A09 Fab.  
Authors : Joyce, M.G.; Andrews, S.F.; Mascola, J.R.; McDermott, A.B.; Kwong, P.D.  
Deposited on : 2016-12-05  
Resolution : 1.76 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	:	2.11
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)	:	2.11

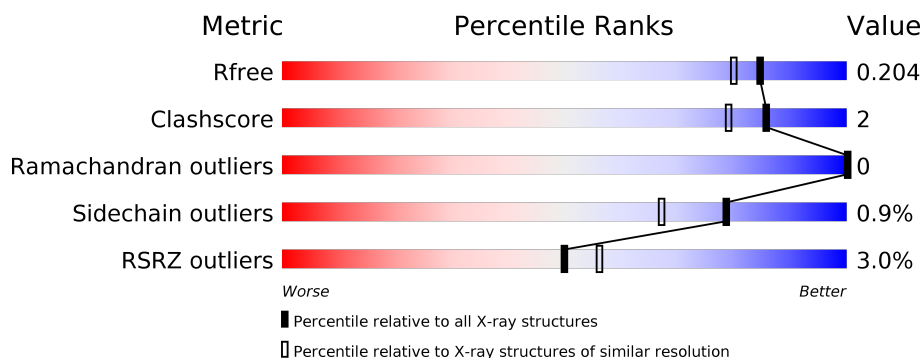
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.76 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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 on 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	235	<div> <div>5%</div> <div>96%</div> <div>••</div> </div>
1	H	235	<div> <div>3%</div> <div>94%</div> <div>5% •</div> </div>
2	B	214	<div> <div>3%</div> <div>93%</div> <div>6% •</div> </div>
2	L	214	<div> <div>%</div> <div>94%</div> <div>5%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14571 atoms, of which 6621 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 VRC 315 53-1A09 Fab Heavy chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	232	Total	C	H	N	O	S	0	0	0
			3405	1088	1680	286	345	6			
1	H	233	Total	C	H	N	O	S	0	1	0
			3444	1098	1703	289	347	7			

- Molecule 2 is a protein called VRC 315 53-1A09 Fab Light chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	212	Total	C	H	N	O	S	0	1	0
			3257	1026	1609	287	330	5			
2	L	214	Total	C	H	N	O	S	0	1	0
			3295	1036	1629	291	333	6			

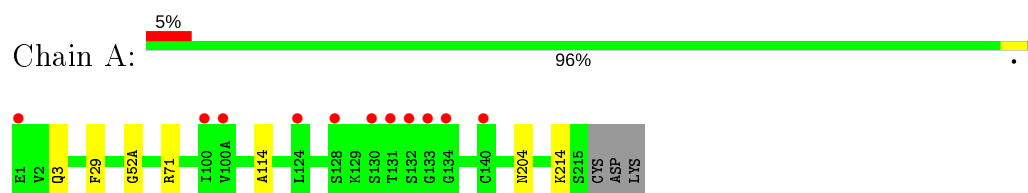
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	272	Total	O	0	0
			272	272		
3	B	242	Total	O	0	0
			242	242		
3	H	385	Total	O	0	0
			385	385		
3	L	271	Total	O	0	0
			271	271		

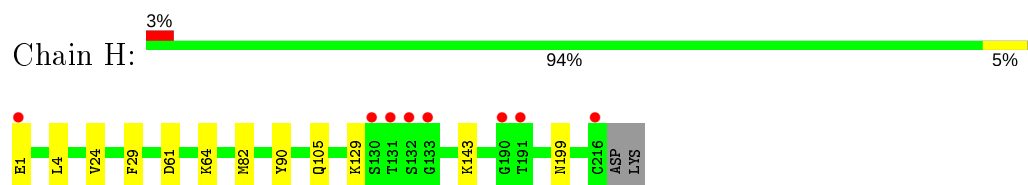
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

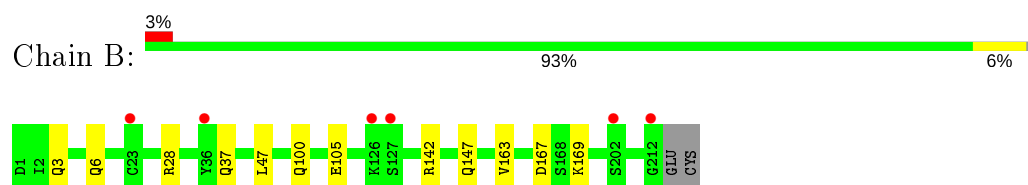
- Molecule 1: VRC 315 53-1A09 Fab Heavy chain



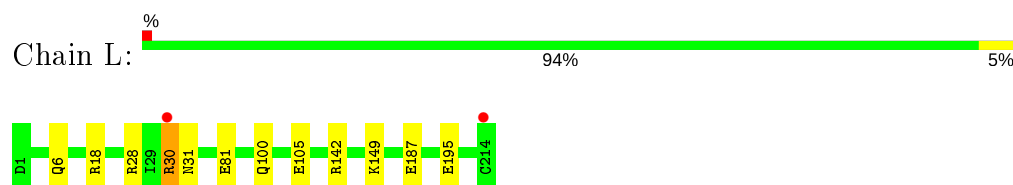
- Molecule 1: VRC 315 53-1A09 Fab Heavy chain



- Molecule 2: VRC 315 53-1A09 Fab Light chain



- Molecule 2: VRC 315 53-1A09 Fab Light chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.55Å 84.13Å 79.27Å 90.00° 111.80° 90.00°	Depositor
Resolution (Å)	33.72 – 1.76 33.72 – 1.76	Depositor EDS
% Data completeness (in resolution range)	98.7 (33.72-1.76) 98.7 (33.72-1.76)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.62 (at 1.76Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, $R_{free}$	0.177 , 0.203 0.179 , 0.204	Depositor DCC
$R_{free}$ test set	4612 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.9	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 50.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	14571	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.40% 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](#)

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	0.27	0/1766	0.48	0/2407
1	H	0.29	0/1785	0.51	0/2431
2	B	0.31	0/1685	0.51	0/2286
2	L	0.36	0/1703	0.51	0/2309
All	All	0.31	0/6939	0.50	0/9433

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	1725	1680	1680	4	0
1	H	1741	1703	1703	8	0
2	B	1648	1609	1612	8	0
2	L	1666	1629	1629	10	0
3	A	272	0	0	3	3
3	B	242	0	0	3	0
3	H	385	0	0	5	1
3	L	271	0	0	7	2
All	All	7950	6621	6624	30	3

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 2.

All (30) 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:L:81:GLU:OE2	3:L:301:HOH:O	1.86	0.91
1:H:143:LYS:NZ	3:H:301:HOH:O	2.00	0.89
2:L:30:ARG:NE	3:L:302:HOH:O	1.92	0.85
1:H:105[A]:GLN:NE2	3:H:302:HOH:O	2.08	0.84
2:L:149:LYS:NZ	2:L:195:GLU:OE1	2.14	0.80
2:L:81:GLU:OE1	3:L:303:HOH:O	2.00	0.79
1:A:114:ALA:O	3:A:301:HOH:O	2.06	0.73
2:B:147:GLN:OE1	3:B:301:HOH:O	2.07	0.72
1:A:214:LYS:O	3:A:302:HOH:O	2.06	0.71
2:B:28:ARG:NH2	3:B:302:HOH:O	2.24	0.65
2:B:6:GLN:O	2:B:100:GLN:NE2	2.34	0.58
1:A:3:GLN:HG3	3:A:495:HOH:O	2.04	0.57
1:H:199:ASN:OD1	3:H:303:HOH:O	2.17	0.56
2:L:6:GLN:O	2:L:100:GLN:NE2	2.39	0.56
2:L:187:GLU:OE1	3:L:304:HOH:O	2.19	0.50
1:H:129:LYS:NZ	3:H:313:HOH:O	2.45	0.48
2:L:30:ARG:CD	3:L:302:HOH:O	2.54	0.46
2:L:18:ARG:NH1	3:L:311:HOH:O	2.48	0.46
2:B:167:ASP:OD2	2:B:169:LYS:HG2	2.16	0.44
2:B:142:ARG:NH2	2:B:163:VAL:HG11	2.32	0.44
1:H:61:ASP:HA	1:H:64:LYS:HG3	2.00	0.43
2:L:28:ARG:NE	3:L:308:HOH:O	2.37	0.43
1:A:52(A):GLY:HA2	1:A:71:ARG:NH1	2.33	0.42
2:B:3:GLN:NE2	3:B:309:HOH:O	2.42	0.42
1:H:1:GLU:HA	3:H:539:HOH:O	2.21	0.41
2:B:167:ASP:OD2	2:B:169:LYS:CG	2.68	0.41
2:B:37:GLN:HB2	2:B:47:LEU:HD11	2.03	0.41
1:H:4:LEU:HG	1:H:24:VAL:HG12	2.03	0.41
1:H:82:MET:HE1	1:H:90:TYR:CZ	2.56	0.41
2:L:30:ARG:HB3	2:L:31:ASN:H	1.78	0.40

All (3) 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)
3:A:497:HOH:O	3:L:551:HOH:O[1_656]	2.10	0.10
3:A:353:HOH:O	3:L:394:HOH:O[2_645]	2.14	0.06
3:A:366:HOH:O	3:H:462:HOH:O[1_655]	2.16	0.04

## 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	A	230/235 (98%)	227 (99%)	3 (1%)	0	100	100
1	H	232/235 (99%)	228 (98%)	4 (2%)	0	100	100
2	B	211/214 (99%)	205 (97%)	6 (3%)	0	100	100
2	L	213/214 (100%)	207 (97%)	6 (3%)	0	100	100
All	All	886/898 (99%)	867 (98%)	19 (2%)	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	A	192/196 (98%)	190 (99%)	2 (1%)	76	63
1	H	195/196 (100%)	194 (100%)	1 (0%)	88	83
2	B	188/189 (100%)	187 (100%)	1 (0%)	88	83
2	L	190/189 (100%)	186 (98%)	4 (2%)	53	31
All	All	765/770 (99%)	757 (99%)	8 (1%)	78	63

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	PHE
1	A	204	ASN

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Mol	Chain	Res	Type
2	B	105	GLU
1	H	29	PHE
2	L	30	ARG
2	L	105	GLU
2	L	142[A]	ARG
2	L	142[B]	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 carbohydrates 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	232/235 (98%)	0.22	11 (4%) 31 37	15, 27, 48, 88	0
1	H	233/235 (99%)	-0.06	8 (3%) 45 51	9, 16, 40, 88	0
2	B	212/214 (99%)	0.38	6 (2%) 53 58	19, 30, 46, 61	0
2	L	214/214 (100%)	0.03	2 (0%) 84 89	13, 24, 44, 89	2 (0%)
All	All	891/898 (99%)	0.14	27 (3%) 50 56	9, 25, 45, 89	2 (0%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	100	ILE	7.2
1	A	1	GLU	6.5
1	H	216	CYS	5.3
1	A	132	SER	4.5
2	L	214	CYS	4.3
1	A	131	THR	4.0
1	A	133	GLY	3.9
1	A	100(A)	VAL	3.9
1	H	133	GLY	3.8
1	H	130	SER	3.8
1	H	132	SER	3.7
1	H	191	THR	3.7
2	B	212	GLY	3.5
1	H	1	GLU	3.5
1	H	131	THR	3.3
2	L	30	ARG	2.9
2	B	202	SER	2.9
1	A	130	SER	2.8
1	A	128	SER	2.8
1	A	140	CYS	2.7
1	A	134	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	H	190	GLY	2.4
1	A	124	LEU	2.4
2	B	127	SER	2.2
2	B	126	LYS	2.2
2	B	23	CYS	2.1
2	B	36	TYR	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 carbohydrates 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.