Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2020 – 02:59 PM EST

PDB ID	:	7JKB
Title	:	2xVH Fab
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Deposited on	:	2020-07-28
Resolution	:	2.55 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 following versions of software and data (see references ⓘ) were used in the production of this report:

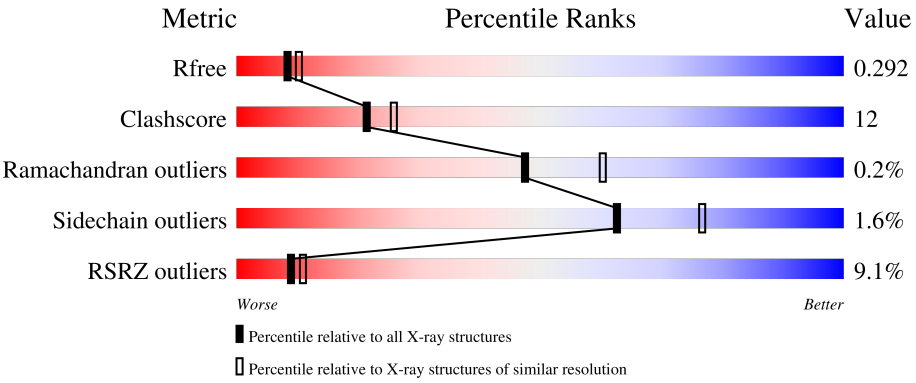
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.14.6
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.14.6

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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 >=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 <=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	226	6% 76% 23%
2	B	239	12% 70% 24% 5%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3356 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 Anti-lysozyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	223	1658	1041	277	334	6	0	0	0

- Molecule 2 is a protein called Anti-Her2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	226	1679	1051	287	332	9	0	0	0

- Molecule 3 is water.

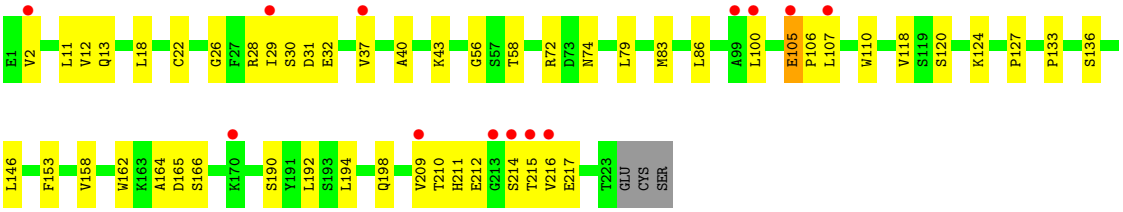
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	9	9	9	0	0
			Total	O		
3	B	10	10	10	0	0

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.

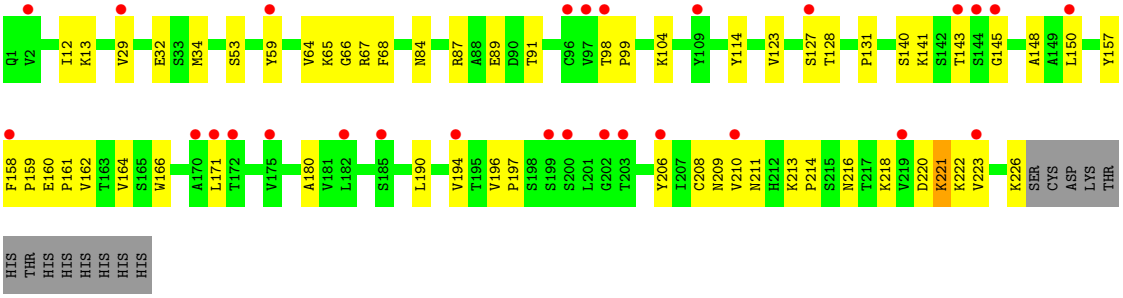
• Molecule 1: Anti-lysozyme

Chain A:



• Molecule 2: Anti-Her2

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	70.48Å 195.00Å 91.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.75 – 2.55 48.75 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.75-2.55) 99.8 (48.75-2.55)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.99 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.14_3211	Depositor
R, R_{free}	0.254 , 0.292 0.254 , 0.292	Depositor DCC
R_{free} test set	2000 reflections (9.54%)	wwPDB-VP
Wilson B-factor (Å^2)	72.8	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 54.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	3356	wwPDB-VP
Average B, all atoms (Å^2)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	A	0.26	0/1698	0.49	0/2318
2	B	0.28	0/1717	0.52	0/2335
All	All	0.27	0/3415	0.50	0/4653

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	A	1658	0	1591	35	2
2	B	1679	0	1656	48	2
3	A	9	0	0	1	0
3	B	10	0	0	0	0
All	All	3356	0	3247	79	2

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 12.

All (79) 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:B:211:ASN:CB	2:B:218:LYS:HD2	2.03	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:150:LEU:HD13	2:B:223:VAL:HG21	1.56	0.85
2:B:171:LEU:HD21	2:B:194:VAL:HG21	1.57	0.84
1:A:100:LEU:HD13	1:A:107:LEU:HD21	1.62	0.80
2:B:211:ASN:HB2	2:B:218:LYS:HD2	1.63	0.80
2:B:211:ASN:CG	2:B:218:LYS:HE2	2.03	0.79
2:B:221:LYS:HA	2:B:221:LYS:HE2	1.62	0.79
1:A:43:LYS:NZ	3:A:301:HOH:O	2.14	0.77
2:B:91:THR:HG22	2:B:123:VAL:H	1.50	0.76
1:A:124:LYS:HE2	1:A:212:GLU:HG3	1.66	0.75
2:B:211:ASN:HD21	2:B:213:LYS:HG2	1.52	0.73
1:A:162:TRP:HE1	1:A:190:SER:HG	1.35	0.72
2:B:211:ASN:ND2	2:B:218:LYS:HE2	2.07	0.70
1:A:100:LEU:HD12	2:B:99:PRO:HB3	1.75	0.69
1:A:209:VAL:HG13	1:A:216:VAL:HG13	1.77	0.67
2:B:180:ALA:HB2	2:B:190:LEU:HD23	1.80	0.64
2:B:29:VAL:HG23	2:B:114:TYR:CZ	2.33	0.64
1:A:158:VAL:HG12	1:A:211:HIS:HB2	1.78	0.64
1:A:107:LEU:HB2	2:B:59:TYR:CZ	2.34	0.62
2:B:211:ASN:CA	2:B:218:LYS:HD2	2.33	0.58
1:A:210:THR:HG22	1:A:215:THR:HA	1.86	0.58
2:B:211:ASN:HA	2:B:218:LYS:HD2	1.87	0.56
1:A:164:ALA:O	1:A:166:SER:N	2.38	0.56
2:B:208:CYS:O	2:B:220:ASP:HA	2.05	0.56
2:B:221:LYS:CE	2:B:221:LYS:HA	2.31	0.55
2:B:127:SER:OG	2:B:128:THR:N	2.39	0.55
1:A:37:VAL:HG11	1:A:110:TRP:HZ3	1.71	0.55
1:A:100:LEU:O	1:A:107:LEU:HG	2.08	0.54
1:A:13:GLN:HG2	1:A:120:SER:HA	1.90	0.52
2:B:211:ASN:CG	2:B:218:LYS:CE	2.77	0.52
1:A:210:THR:HA	1:A:214:SER:O	2.10	0.52
1:A:194:LEU:HD21	1:A:198:GLN:HB2	1.92	0.51
1:A:2:VAL:HG12	1:A:26:GLY:HA3	1.92	0.51
2:B:221:LYS:HD3	2:B:222:LYS:N	2.26	0.51
1:A:127:PRO:HB3	1:A:153:PHE:HB3	1.94	0.49
2:B:148:ALA:HB3	2:B:196:VAL:O	2.12	0.49
1:A:100:LEU:CD1	2:B:99:PRO:HB3	2.44	0.47
2:B:211:ASN:CG	2:B:218:LYS:CD	2.83	0.46
1:A:28:ARG:NH2	1:A:74:ASN:O	2.48	0.46
2:B:104:LYS:HA	2:B:104:LYS:HD2	1.74	0.46
2:B:12:ILE:HG13	2:B:13:LYS:H	1.81	0.45
1:A:127:PRO:HG2	1:A:216:VAL:HG11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:162:VAL:HG23	2:B:211:ASN:O	2.17	0.44
2:B:87:ARG:HB2	2:B:89:GLU:HG2	1.99	0.44
2:B:164:VAL:HA	2:B:209:ASN:O	2.18	0.44
2:B:166:TRP:CH2	2:B:208:CYS:HB3	2.53	0.44
2:B:160:GLU:HG3	2:B:161:PRO:HA	2.00	0.44
2:B:211:ASN:CG	2:B:218:LYS:HD2	2.37	0.44
2:B:143:THR:HG22	2:B:145:GLY:H	1.83	0.43
2:B:210:VAL:O	2:B:218:LYS:HA	2.17	0.43
2:B:221:LYS:HD3	2:B:222:LYS:H	1.83	0.43
2:B:140:SER:HB2	2:B:141:LYS:HD3	2.00	0.43
2:B:141:LYS:HE2	2:B:141:LYS:HB2	1.81	0.43
1:A:107:LEU:HD13	2:B:59:TYR:CD1	2.54	0.43
1:A:29:ILE:HD12	1:A:30:SER:N	2.34	0.43
1:A:133:PRO:HA	1:A:146:LEU:HD23	2.02	0.42
1:A:18:LEU:HB2	1:A:83:MET:HE2	2.01	0.42
2:B:158:PHE:HA	2:B:159:PRO:HA	1.75	0.42
1:A:40:ALA:HB3	1:A:43:LYS:HB2	2.02	0.42
1:A:105:GLU:HB3	1:A:106:PRO:HD3	2.02	0.42
2:B:65:LYS:NZ	2:B:66:GLY:H	2.17	0.42
1:A:217:GLU:N	1:A:217:GLU:OE1	2.53	0.42
2:B:131:PRO:HB3	2:B:157:TYR:HB3	2.01	0.42
1:A:12:VAL:O	1:A:118:VAL:HA	2.19	0.41
2:B:64:VAL:HG13	2:B:68:PHE:CG	2.55	0.41
1:A:83:MET:HB3	1:A:86:LEU:HD21	2.02	0.41
2:B:196:VAL:HG21	2:B:206:TYR:CZ	2.55	0.41
1:A:29:ILE:HD13	1:A:72:ARG:HB3	2.02	0.41
1:A:146:LEU:HD12	1:A:192:LEU:HD23	2.02	0.41
2:B:32:GLU:HA	2:B:53:SER:HB2	2.02	0.41
2:B:34:MET:SD	2:B:98:THR:HG22	2.61	0.41
2:B:211:ASN:CB	2:B:218:LYS:CD	2.89	0.41
1:A:32:GLU:O	1:A:72:ARG:NH2	2.50	0.41
1:A:56:GLY:O	1:A:58:THR:HG23	2.20	0.40
1:A:22:CYS:HB3	1:A:79:LEU:HB3	2.02	0.40
2:B:67:ARG:HB3	2:B:84:ASN:O	2.21	0.40
2:B:213:LYS:N	2:B:214:PRO:HD2	2.36	0.40
1:A:11:LEU:HD22	1:A:12:VAL:H	1.86	0.40
2:B:166:TRP:CZ3	2:B:208:CYS:HB3	2.56	0.40

All (2) 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:A:217:GLU:OE2	2:B:218:LYS:N[8_455]	1.98	0.22
1:A:31:ASP:OD1	2:B:104:LYS:NZ[4_555]	2.05	0.15

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	221/226 (98%)	212 (96%)	8 (4%)	1 (0%)	29 40
2	B	224/239 (94%)	207 (92%)	17 (8%)	0	100 100
All	All	445/465 (96%)	419 (94%)	25 (6%)	1 (0%)	47 60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	165	ASP

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	180/188 (96%)	178 (99%)	2 (1%)	73 83
2	B	190/203 (94%)	186 (98%)	4 (2%)	53 68
All	All	370/391 (95%)	364 (98%)	6 (2%)	62 77

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	GLU
1	A	136	SER
2	B	197	PRO
2	B	216	ASN
2	B	221	LYS
2	B	226	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	211	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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^{th} 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	A	223/226 (98%)	0.78	13 (5%) 23 27	53, 77, 104, 125	0
2	B	226/239 (94%)	0.93	28 (12%) 4 5	54, 81, 112, 144	0
All	All	449/465 (96%)	0.86	41 (9%) 9 11	53, 79, 110, 144	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	213	GLY	9.2
2	B	145	GLY	5.9
1	A	214	SER	4.8
2	B	199	SER	4.4
2	B	96	CYS	4.4
2	B	171	LEU	4.0
2	B	203	THR	4.0
2	B	144	SER	3.9
2	B	206	TYR	3.6
1	A	107	LEU	3.5
1	A	100	LEU	3.4
2	B	175	VAL	3.3
2	B	143	THR	3.2
2	B	59	TYR	3.0
2	B	219	VAL	2.9
2	B	127	SER	2.9
2	B	109	TYR	2.8
1	A	2	VAL	2.7
2	B	200	SER	2.7
2	B	29	VAL	2.7
1	A	170	LYS	2.7
2	B	185	SER	2.6
1	A	99	ALA	2.6
2	B	210	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
2	B	194	VAL	2.5
1	A	216	VAL	2.4
1	A	105	GLU	2.4
1	A	209	VAL	2.3
2	B	170	ALA	2.3
2	B	172	THR	2.3
1	A	37	VAL	2.2
2	B	223	VAL	2.1
2	B	150	LEU	2.1
2	B	2	VAL	2.1
2	B	97	VAL	2.1
2	B	98	THR	2.1
2	B	202	GLY	2.1
2	B	158	PHE	2.1
2	B	182	LEU	2.1
1	A	215	THR	2.0
1	A	29	ILE	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.