Full wwPDB X-ray Structure Validation Report (i)

Jun 19, 2024 – 12:28 PM EDT

PDB ID	:	4OJF
Title	:	Humanised 3D6 Fab complexed to amyloid beta 1-8
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Deposited on	:	2014-01-21
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 (i) 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 (1)) were used in the production of this report:

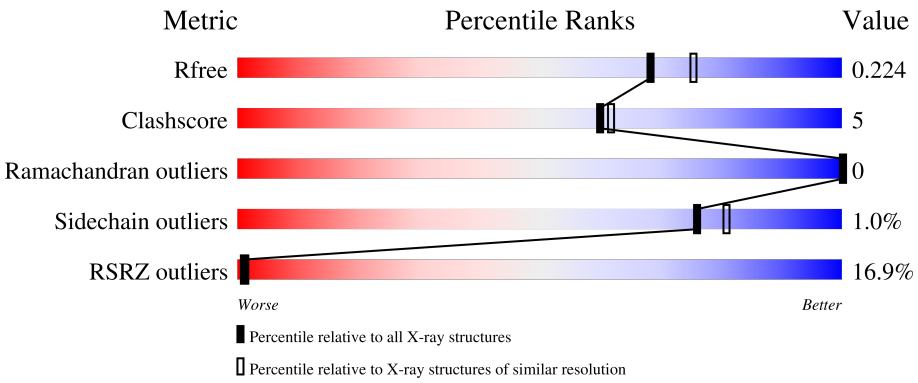
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 <=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	H	227	17% 89% 7% •
2	L	219	16% 88% 12%
3	A	8	12% 75% 25%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6984 atoms, of which 3304 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 Humanised 3D6 Fab Heavy Chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	H	219	3234	1027	1599	281	321	6	0	0	0

- Molecule 2 is a protein called Humanised 3D6 Fab Light Chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	L	219	3347	1062	1658	283	338	6	0	0	0

- Molecule 3 is a protein called Amyloid beta A4 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	H	N	O			
3	A	6	101	33	47	11	10	0	0	0

- Molecule 4 is water.

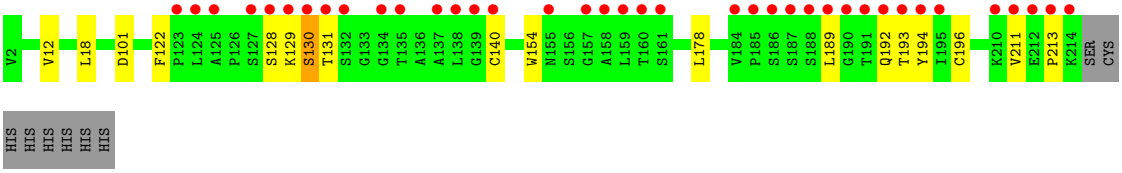
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	H	158	158	158	0	0
4	L	141	141	141	0	0
4	A	3	3	3	0	0

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.

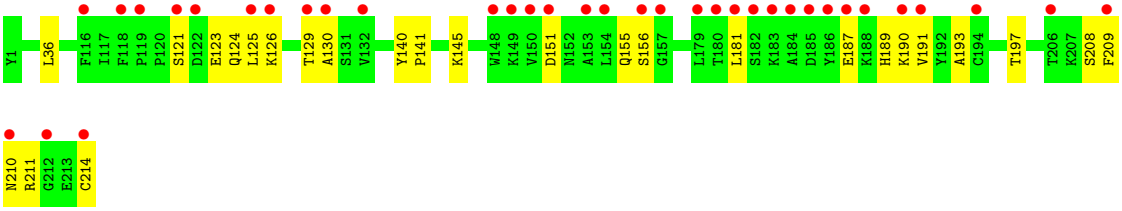
- Molecule 1: Humanised 3D6 Fab Heavy Chain

Chain H:



- Molecule 2: Humanised 3D6 Fab Light Chain

Chain L:



- Molecule 3: Amyloid beta A4 protein

Chain A:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 2 21	Depositor
Cell constants a, b, c, α, β, γ	60.16Å 83.43Å 88.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.70 – 2.00 42.70 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (42.70-2.00) 98.4 (42.70-2.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.28 (at 2.00Å)	Xtriage
Refinement program	REFMAC, PHENIX 1.8_1069	Depositor
R, R_{free}	0.172 , 0.215 0.184 , 0.224	Depositor DCC
R_{free} test set	1546 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	17.0	Xtriage
Anisotropy	0.343	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 51.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6984	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.10% 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 (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	H	0.42	0/1674	0.63	1/2276 (0.0%)
2	L	0.43	0/1727	0.60	0/2345
3	A	0.51	0/55	0.53	0/72
All	All	0.43	0/3456	0.61	1/4693 (0.0%)

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	H	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	H	101	ASP	CB-CG-OD1	5.99	123.69	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	130	SER	Peptide

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	H	1635	1599	1596	11	0
2	L	1689	1658	1655	21	0
3	A	54	47	46	0	0
4	A	3	0	0	0	0
4	H	158	0	0	1	0
4	L	141	0	0	1	0
All	All	3680	3304	3297	31	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 5.

All (31) 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:H:129:LYS:N	1:H:130:SER:OG	2.04	0.91
2:L:187:GLU:OE2	2:L:211:ARG:NH2	2.12	0.81
2:L:190:LYS:HA	2:L:211:ARG:HD3	1.64	0.77
2:L:189:HIS:O	2:L:211:ARG:NH1	2.22	0.73
2:L:189:HIS:O	2:L:211:ARG:HD2	1.89	0.73
2:L:189:HIS:O	2:L:211:ARG:CD	2.42	0.68
2:L:190:LYS:HA	2:L:211:ARG:CD	2.29	0.62
2:L:121:SER:O	2:L:125:LEU:HB2	2.04	0.58
2:L:151:ASP:HA	2:L:191:VAL:HB	1.86	0.58
1:H:192:GLN:OE1	1:H:193:THR:N	2.31	0.56
2:L:190:LYS:O	2:L:210:ASN:HA	2.05	0.56
1:H:192:GLN:HG3	1:H:194:TYR:CE1	2.42	0.55
2:L:187:GLU:HA	2:L:211:ARG:NH2	2.22	0.53
2:L:126:LYS:NZ	4:L:440:HOH:O	2.42	0.52
2:L:208:SER:O	2:L:209:PHE:CD1	2.65	0.50
2:L:36:LEU:C	2:L:36:LEU:HD12	2.34	0.47
1:H:211:VAL:O	4:H:440:HOH:O	2.20	0.47
2:L:140:TYR:CG	2:L:141:PRO:HA	2.50	0.46
1:H:128:SER:C	1:H:130:SER:OG	2.54	0.46
1:H:128:SER:HA	1:H:129:LYS:HA	1.65	0.45
2:L:145:LYS:HB2	2:L:197:THR:HB	1.98	0.45
2:L:155:GLN:O	2:L:156:SER:OG	2.30	0.44
2:L:193:ALA:HA	2:L:208:SER:OG	2.18	0.43
2:L:193:ALA:CB	2:L:208:SER:OG	2.66	0.43
2:L:123:GLU:O	2:L:126:LYS:HG2	2.18	0.43
1:H:122:PHE:CE2	2:L:124:GLN:HG3	2.54	0.42
2:L:130:ALA:N	2:L:181:LEU:O	2.42	0.42
1:H:12:VAL:CG2	1:H:18:LEU:HD13	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:178:LEU:C	1:H:178:LEU:HD12	2.41	0.41
1:H:189:LEU:HD11	1:H:213:PRO:HG2	2.01	0.41
1:H:154:TRP:CH2	1:H:196:CYS:HB3	2.55	0.41

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	H	217/227 (96%)	206 (95%)	11 (5%)	0	100	100
2	L	217/219 (99%)	211 (97%)	6 (3%)	0	100	100
3	A	4/8 (50%)	4 (100%)	0	0	100	100
All	All	438/454 (96%)	421 (96%)	17 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	H	182/190 (96%)	180 (99%)	2 (1%)	73	78
2	L	195/195 (100%)	193 (99%)	2 (1%)	76	81
3	A	5/7 (71%)	5 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	382/392 (97%)	378 (99%)	4 (1%)	76 81

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	131	THR
1	H	140	CYS
2	L	129	THR
2	L	214	CYS

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

Mol	Chain	Res	Type
1	H	148	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 (i)

6.1 Protein, DNA and RNA chains (i)

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	H	219/227 (96%)	0.78	38 (17%)	1	1	6, 18, 71, 85	0
2	L	219/219 (100%)	0.75	36 (16%)	1	1	8, 22, 66, 91	0
3	A	6/8 (75%)	0.12	1 (16%)	1	1	8, 9, 19, 40	0
All	All	444/454 (97%)	0.76	75 (16%)	1	1	6, 20, 69, 91	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	214	CYS	13.1
2	L	154	LEU	7.7
2	L	210	ASN	7.6
2	L	125	LEU	7.4
1	H	128	SER	6.6
1	H	195	ILE	6.0
1	H	130	SER	5.9
1	H	189	LEU	5.8
1	H	193	THR	5.8
1	H	211	VAL	5.2
2	L	153	ALA	5.0
1	H	212	GLU	4.7
1	H	159	LEU	4.6
2	L	209	PHE	4.6
2	L	126	LYS	4.5
1	H	194	TYR	4.4
1	H	190	GLY	4.4
2	L	157	GLY	4.4
1	H	187	SER	4.3
1	H	192	GLN	4.3
1	H	138	LEU	4.2
2	L	156	SER	4.2
2	L	181	LEU	3.9

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Mol	Chain	Res	Type	RSRZ
2	L	184	ALA	3.9
1	H	129	LYS	3.8
2	L	129	THR	3.8
1	H	127	SER	3.7
2	L	119	PRO	3.5
1	H	131	THR	3.3
1	H	186	SER	3.3
2	L	122	ASP	3.3
1	H	132	SER	3.2
2	L	191	VAL	3.2
2	L	121	SER	3.2
2	L	183	LYS	3.1
1	H	140	CYS	3.1
1	H	191	THR	2.9
2	L	187	GLU	2.8
1	H	124	LEU	2.8
3	A	6	HIS	2.8
1	H	139	GLY	2.8
1	H	158	ALA	2.7
2	L	132	VAL	2.7
1	H	125	ALA	2.7
1	H	184	VAL	2.7
1	H	123	PRO	2.6
1	H	210	LYS	2.6
1	H	213	PRO	2.6
2	L	190	LYS	2.6
2	L	179	LEU	2.5
2	L	150	VAL	2.5
1	H	214	LYS	2.4
2	L	151	ASP	2.4
1	H	157	GLY	2.4
2	L	186	TYR	2.4
2	L	130	ALA	2.4
1	H	135	THR	2.4
2	L	149	LYS	2.3
1	H	155	ASN	2.3
2	L	194	CYS	2.3
2	L	185	ASP	2.3
2	L	116	PHE	2.3
1	H	161	SER	2.3
1	H	188	SER	2.3
2	L	182	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	H	185	PRO	2.3
1	H	134	GLY	2.2
2	L	118	PHE	2.2
2	L	212	GLY	2.2
1	H	137	ALA	2.1
1	H	160	THR	2.0
2	L	180	THR	2.0
2	L	148	TRP	2.0
2	L	188	LYS	2.0
2	L	206	THR	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.