



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:21 PM GMT

PDB ID : 4OKV
Title : Crystal structure of anopheline anti-platelet protein with Fab antibody
Authors : Park, S.Y.; Sugiyama, K.
Deposited on : 2014-01-23
Resolution : 1.80 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

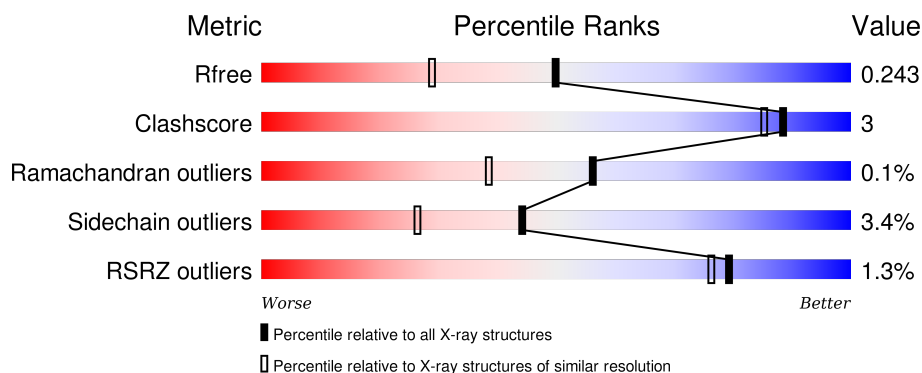
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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. 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	215	<div> <div>4%</div> <div> <div></div> <div>92%</div> <div>7%</div> </div> </div>
1	C	215	<div> <div></div> <div>96%</div> <div>.</div> </div>
2	B	218	<div> <div></div> <div>88%</div> <div>10%</div> <div>.</div> </div>
2	D	218	<div> <div>%</div> <div>94%</div> <div>6%</div> <div>.</div> </div>
3	E	66	<div> <div>2%</div> <div>91%</div> <div>5%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	66	<div><div></div><div>3%</div><div>82%</div><div>17%</div><div></div></div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8351 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 heavy chain of 8H7 mAb.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1614	1019	265	322	8			
1	C	215	Total	C	N	O	S	0	0	0
			1620	1022	266	323	9			

- Molecule 2 is a protein called light chain of 8H7 mAb.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	218	Total	C	N	O	S	0	0	0
			1688	1059	284	339	6			
2	D	218	Total	C	N	O	S	0	0	0
			1688	1059	284	339	6			

- Molecule 3 is a protein called Anti-platelet aggregation protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	65	Total	C	N	O	S	0	0	0
			517	312	87	114	4			
3	F	66	Total	C	N	O	S	0	0	0
			526	318	89	115	4			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	145	Total	O	0	0
			145	145		
4	B	142	Total	O	0	0
			142	142		
4	C	179	Total	O	0	0
			179	179		
4	D	163	Total	O	0	0
			163	163		

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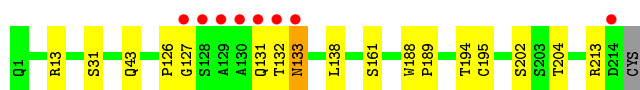
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	36	Total	O	0	0
			36	36		
4	F	33	Total	O	0	0
			33	33		

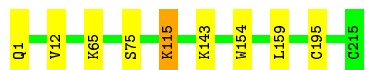
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: heavy chain of 8H7 mAb



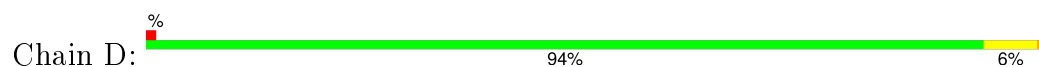
- Molecule 1: heavy chain of 8H7 mAb



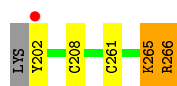
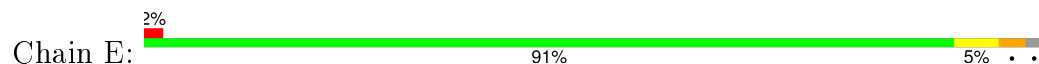
- Molecule 2: light chain of 8H7 mAb



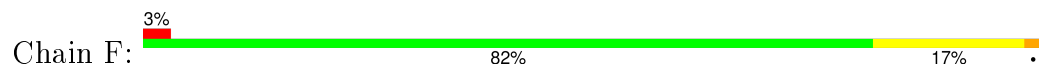
- Molecule 2: light chain of 8H7 mAb



- Molecule 3: Anti-platelet aggregation protein



- Molecule 3: Anti-platelet aggregation protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	93.81Å 99.43Å 166.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.62 – 1.80 47.62 – 1.80	Depositor EDS
% Data completeness (in resolution range)	94.3 (47.62-1.80) 94.3 (47.62-1.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 1.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.218 , 0.250 0.211 , 0.243	Depositor DCC
R_{free} test set	6804 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	17.1	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 135293 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8351	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.38	0/1657	0.57	0/2269
1	C	0.39	0/1663	0.56	0/2277
2	B	0.38	0/1726	0.59	2/2345 (0.1%)
2	D	0.39	0/1726	0.53	0/2345
3	E	0.41	0/521	0.45	0/694
3	F	0.41	0/530	0.47	0/705
All	All	0.39	0/7823	0.55	2/10635 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	51	ARG	NE-CZ-NH2	-5.76	117.42	120.30
2	B	38	LEU	CA-CB-CG	-5.31	103.08	115.30

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	1614	0	1574	7	0
1	C	1620	0	1579	6	0
2	B	1688	0	1644	12	0
2	D	1688	0	1644	7	0
3	E	517	0	492	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	526	0	505	7	0
4	A	145	0	0	0	0
4	B	142	0	0	1	0
4	C	179	0	0	1	0
4	D	163	0	0	0	0
4	E	36	0	0	0	0
4	F	33	0	0	0	0
All	All	8351	0	7438	38	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 (38) 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:127:GLY:HA2	1:A:213:ARG:HD2	1.65	0.78
3:E:202:TYR:HD1	3:F:244:THR:HG21	1.55	0.71
2:B:154:LYS:NZ	2:B:200:GLU:OE2	2.23	0.68
1:A:133:ASN:OD1	1:A:133:ASN:N	2.31	0.62
2:B:51:ARG:NH2	3:F:233:ASP:OD1	2.29	0.60
2:D:44:ARG:NH1	2:D:86:GLU:O	2.35	0.59
2:B:113:ARG:HD3	2:B:114:ALA:O	2.04	0.58
1:C:1:GLN:N	1:C:1:GLN:OE1	2.32	0.55
2:B:31:ASP:OD1	3:F:234:LYS:NZ	2.44	0.51
2:B:217:ASN:OD1	2:B:218:ALA:N	2.44	0.51
1:A:132:THR:HG22	1:A:133:ASN:H	1.76	0.50
1:C:154:TRP:CZ3	1:C:195:CYS:HB3	2.47	0.49
1:A:132:THR:HG22	1:A:133:ASN:N	2.27	0.49
2:D:86:GLU:H	2:D:86:GLU:CD	2.17	0.49
2:B:94:TRP:HE1	2:B:101:TYR:HB3	1.78	0.48
2:B:127:SER:O	2:B:131:THR:HG23	2.15	0.47
2:B:52:LEU:HA	2:B:63:VAL:HG21	1.96	0.46
3:E:266:ARG:NH1	3:F:252:ILE:HD13	2.31	0.46
2:D:42:LEU:HB2	2:D:52:LEU:HD11	1.98	0.46
1:A:188:TRP:CG	1:A:189:PRO:HA	2.52	0.45
2:B:186:LEU:HD13	2:B:191:TYR:HB2	1.99	0.45
2:B:44:ARG:NH2	4:B:367:HOH:O	2.46	0.45
3:E:261:CYS:SG	3:E:265:LYS:NZ	2.91	0.44
1:C:143:LYS:HE2	4:C:426:HOH:O	2.18	0.44
2:B:23:CYS:HB2	2:B:40:TRP:CH2	2.53	0.44
2:B:38:LEU:HG	2:B:39:ASN:N	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:94:TRP:HE1	2:D:101:TYR:HB3	1.83	0.43
1:C:154:TRP:CH2	1:C:195:CYS:HB3	2.53	0.43
3:F:263:VAL:HA	3:F:266:ARG:HD3	2.00	0.43
1:C:65:LYS:HE3	1:C:65:LYS:HB2	1.77	0.42
1:A:31:SER:OG	3:F:242:GLU:OE1	2.29	0.42
1:C:115:LYS:CE	1:C:115:LYS:H	2.33	0.42
1:A:202:SER:O	1:A:204:THR:HG23	2.19	0.42
3:F:264:GLU:HB3	3:F:265:LYS:NZ	2.35	0.42
2:D:152:LYS:HB3	2:D:152:LYS:HE3	1.85	0.42
2:D:155:ILE:HD11	2:D:184:LEU:HD21	2.01	0.42
2:D:88:LEU:HD21	2:D:173:SER:HA	2.02	0.41
3:E:208:CYS:SG	3:E:265:LYS:NZ	2.94	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	A	212/215 (99%)	208 (98%)	3 (1%)	1 (0%)	34	17
1	C	213/215 (99%)	212 (100%)	1 (0%)	0	100	100
2	B	216/218 (99%)	212 (98%)	4 (2%)	0	100	100
2	D	216/218 (99%)	214 (99%)	2 (1%)	0	100	100
3	E	63/66 (96%)	63 (100%)	0	0	100	100
3	F	64/66 (97%)	64 (100%)	0	0	100	100
All	All	984/998 (99%)	973 (99%)	10 (1%)	1 (0%)	56	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	126	PRO

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	184/185 (100%)	176 (96%)	8 (4%)	35	17
1	C	185/185 (100%)	181 (98%)	4 (2%)	60	45
2	B	193/193 (100%)	186 (96%)	7 (4%)	42	24
2	D	193/193 (100%)	188 (97%)	5 (3%)	54	37
3	E	61/62 (98%)	59 (97%)	2 (3%)	45	27
3	F	62/62 (100%)	58 (94%)	4 (6%)	21	7
All	All	878/880 (100%)	848 (97%)	30 (3%)	44	26

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	43	GLN
1	A	131	GLN
1	A	133	ASN
1	A	138	LEU
1	A	161	SER
1	A	194	THR
1	A	195	CYS
2	B	113	ARG
2	B	159	GLU
2	B	161	GLN
2	B	174	LYS
2	B	186	LEU
2	B	193	ARG
2	B	208	SER
1	C	12	VAL
1	C	75	SER
1	C	115	LYS
1	C	159	LEU
2	D	13	VAL
2	D	88	LEU
2	D	152	LYS

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Mol	Chain	Res	Type
2	D	180	MET
2	D	186	LEU
3	E	265	LYS
3	E	266	ARG
3	F	201	LYS
3	F	253	GLU
3	F	261	CYS
3	F	266	ARG

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

Mol	Chain	Res	Type
2	B	27	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 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 [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	A	214/215 (99%)	-0.06	8 (3%)	45	39	13, 21, 49, 98	0
1	C	215/215 (100%)	-0.44	0	100	100	11, 20, 35, 62	0
2	B	218/218 (100%)	-0.32	0	100	100	13, 24, 37, 52	0
2	D	218/218 (100%)	-0.36	2 (0%)	85	83	12, 22, 34, 58	0
3	E	65/66 (98%)	-0.09	1 (1%)	76	72	12, 27, 48, 55	0
3	F	66/66 (100%)	-0.02	2 (3%)	54	48	16, 31, 48, 58	0
All	All	996/998 (99%)	-0.27	13 (1%)	79	76	11, 23, 41, 98	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	130	ALA	8.8
1	A	131	GLN	8.6
1	A	129	ALA	8.4
1	A	127	GLY	6.7
1	A	132	THR	5.7
1	A	133	ASN	5.0
1	A	128	SER	4.9
3	F	202	TYR	4.5
3	E	202	TYR	4.1
2	D	218	ALA	3.0
1	A	214	ASP	2.9
2	D	217	ASN	2.4
3	F	201	LYS	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.