



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 01:35 am BST

PDB ID : 5H7B
Title : Crystal structure of a repeat protein with five Protein A repeat modules
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Deposited on : 2016-11-17
Resolution : 3.10 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	:	2.11
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

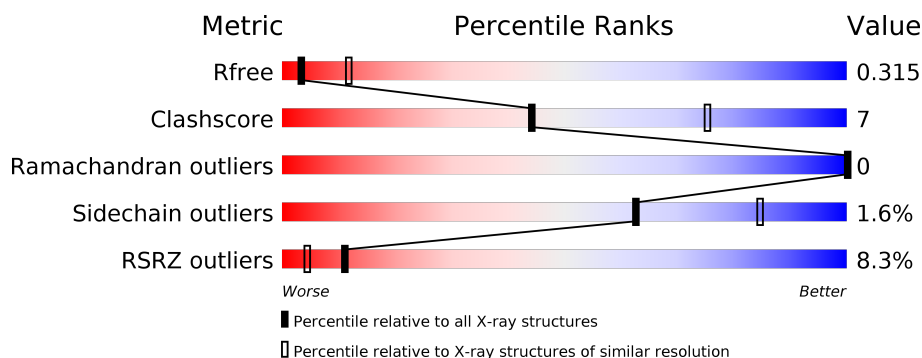
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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 $\leq 5\%$. 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	237	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 15%, orange 84%, yellow 12%, grey 1%);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> 15% 84% 12% •• </div> </div>
1	B	237	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 15%, orange 62%, yellow 18%, grey 19%);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> 15% 62% 18% 19% </div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3387 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 Immunoglobulin G-binding protein A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	230	Total	C	N	O	0	0	0
			1845	1149	321	375			
1	B	191	Total	C	N	O	0	0	0
			1542	964	266	312			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	GLY	-	expression tag	UNP P38507
A	31	SER	-	expression tag	UNP P38507
A	32	HIS	-	expression tag	UNP P38507
A	33	MET	-	expression tag	UNP P38507
A	59	ALA	GLY	engineered mutation	UNP P38507
A	86	ALA	ASN	engineered mutation	UNP P38507
A	93	SER	HIS	engineered mutation	UNP P38507
A	104	ALA	GLY	engineered mutation	UNP P38507
A	131	ALA	ASN	engineered mutation	UNP P38507
A	149	ALA	GLY	engineered mutation	UNP P38507
A	176	ALA	ASN	engineered mutation	UNP P38507
A	194	ALA	GLY	engineered mutation	UNP P38507
A	221	ALA	ASN	engineered mutation	UNP P38507
A	239	ALA	GLY	engineered mutation	UNP P38507
B	30	GLY	-	expression tag	UNP P38507
B	31	SER	-	expression tag	UNP P38507
B	32	HIS	-	expression tag	UNP P38507
B	33	MET	-	expression tag	UNP P38507
B	59	ALA	GLY	engineered mutation	UNP P38507
B	86	ALA	ASN	engineered mutation	UNP P38507
B	93	SER	HIS	engineered mutation	UNP P38507
B	104	ALA	GLY	engineered mutation	UNP P38507
B	131	ALA	ASN	engineered mutation	UNP P38507
B	149	ALA	GLY	engineered mutation	UNP P38507
B	176	ALA	ASN	engineered mutation	UNP P38507

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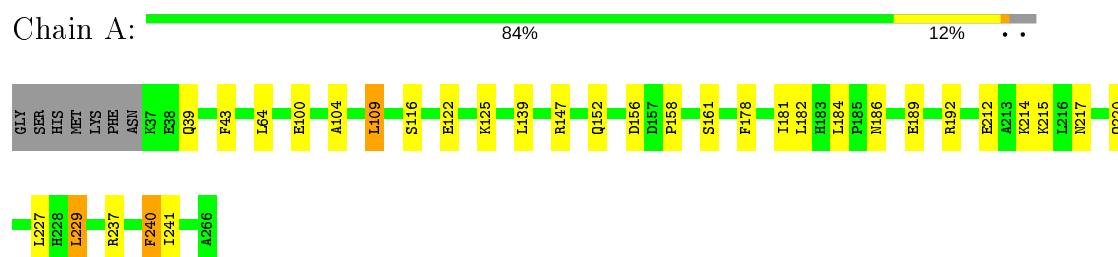
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Chain	Residue	Modelled	Actual	Comment	Reference
B	194	ALA	GLY	engineered mutation	UNP P38507
B	221	ALA	ASN	engineered mutation	UNP P38507
B	239	ALA	GLY	engineered mutation	UNP P38507

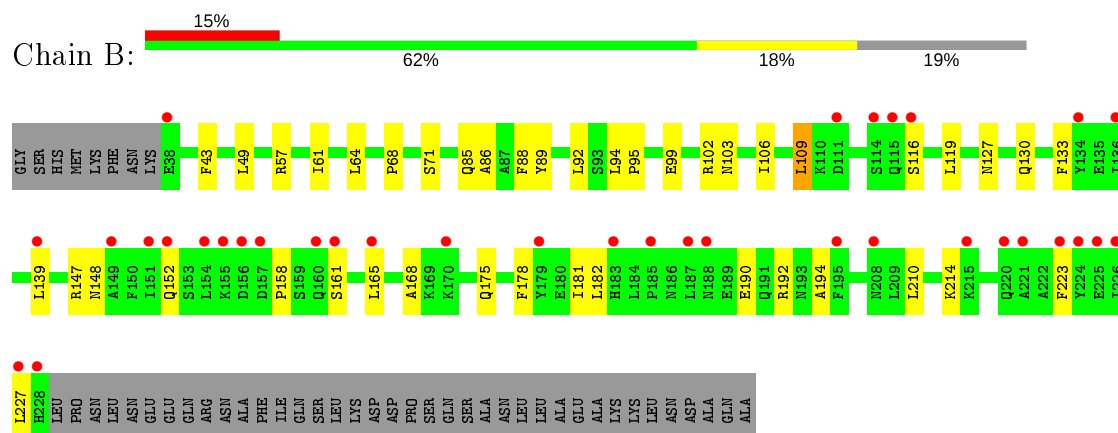
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: Immunoglobulin G-binding protein A



- Molecule 1: Immunoglobulin G-binding protein A



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	80.89Å 45.64Å 82.14Å 90.00° 105.72° 90.00°	Depositor
Resolution (Å)	39.53 – 3.10 39.53 – 3.07	Depositor EDS
% Data completeness (in resolution range)	74.5 (39.53-3.10) 74.6 (39.53-3.07)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.56 (at 3.06Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.269 , 0.315 0.270 , 0.315	Depositor DCC
R_{free} test set	420 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 81.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.022 for l,-k,h	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	3387	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.52% 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, 0.333 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	0.27	0/1873	0.43	0/2532
1	B	0.25	0/1567	0.37	0/2119
All	All	0.26	0/3440	0.40	0/4651

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	1845	0	1793	18	0
1	B	1542	0	1491	26	0
All	All	3387	0	3284	44	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 7.

All (44) 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:186:ASN:O	1:A:220:GLN:NE2	2.13	0.80
1:B:130:GLN:O	1:B:133:PHE:HB3	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:GLU:OE2	1:A:192:ARG:NH1	2.27	0.66
1:A:227:LEU:HD21	1:A:241:ILE:HD11	1.79	0.65
1:A:39:GLN:O	1:A:43:PHE:HB2	1.96	0.65
1:B:148:ASN:O	1:B:152:GLN:HB2	1.97	0.64
1:A:152:GLN:NE2	1:A:156:ASP:OD1	2.33	0.61
1:B:109:LEU:HD21	1:B:116:SER:HA	1.83	0.60
1:A:229:LEU:O	1:A:237:ARG:NH2	2.36	0.59
1:A:158:PRO:O	1:A:161:SER:OG	2.21	0.57
1:A:139:LEU:O	1:A:147:ARG:NH2	2.37	0.56
1:B:139:LEU:O	1:B:147:ARG:NH2	2.32	0.56
1:A:212:GLU:HA	1:A:215:LYS:HE3	1.87	0.56
1:A:214:LYS:O	1:A:217:ASN:HB3	2.08	0.53
1:B:99:GLU:O	1:B:103:ASN:ND2	2.44	0.51
1:B:88:PHE:HE1	1:B:109:LEU:HD22	1.75	0.50
1:B:223:PHE:O	1:B:227:LEU:HB2	2.12	0.50
1:A:122:GLU:HA	1:A:125:LYS:HE3	1.95	0.48
1:A:109:LEU:HD21	1:A:116:SER:HA	1.97	0.46
1:B:109:LEU:HG	1:B:119:LEU:HD13	1.97	0.46
1:B:175:GLN:O	1:B:178:PHE:HB2	2.16	0.46
1:B:95:PRO:HA	1:B:102:ARG:HH21	1.81	0.46
1:B:68:PRO:O	1:B:71:SER:OG	2.25	0.45
1:B:210:LEU:HG	1:B:214:LYS:HE2	1.99	0.44
1:B:178:PHE:O	1:B:182:LEU:HB2	2.17	0.44
1:A:100:GLU:O	1:A:104:ALA:HB2	2.17	0.44
1:B:148:ASN:O	1:B:152:GLN:CB	2.62	0.44
1:B:88:PHE:CD1	1:B:109:LEU:HD13	2.53	0.44
1:B:94:LEU:HB3	1:B:127:ASN:OD1	2.18	0.43
1:B:158:PRO:HA	1:B:161:SER:HB3	2.00	0.43
1:B:92:LEU:HD13	1:B:106:ILE:HD11	2.00	0.43
1:B:165:LEU:O	1:B:168:ALA:HB3	2.19	0.43
1:A:229:LEU:HA	1:A:229:LEU:HD12	1.71	0.42
1:B:181:ILE:O	1:B:192:ARG:HD3	2.20	0.42
1:B:190:GLU:O	1:B:194:ALA:HB2	2.20	0.41
1:B:43:PHE:CD1	1:B:61:ILE:HD12	2.56	0.41
1:A:109:LEU:HD21	1:A:116:SER:CA	2.50	0.41
1:A:181:ILE:HA	1:A:184:LEU:HG	2.03	0.41
1:B:86:ALA:O	1:B:89:TYR:HB3	2.21	0.41
1:B:85:GLN:HA	1:B:88:PHE:CD2	2.56	0.41
1:A:178:PHE:O	1:A:182:LEU:HB2	2.21	0.41
1:B:49:LEU:O	1:B:57:ARG:NH1	2.54	0.41
1:B:139:LEU:HD23	1:B:139:LEU:HA	1.91	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:PHE:HD1	1:A:240:PHE:HA	1.76	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	228/237 (96%)	217 (95%)	11 (5%)	0	100	100
1	B	189/237 (80%)	175 (93%)	14 (7%)	0	100	100
All	All	417/474 (88%)	392 (94%)	25 (6%)	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	A	199/205 (97%)	195 (98%)	4 (2%)	55	80
1	B	166/205 (81%)	164 (99%)	2 (1%)	71	88
All	All	365/410 (89%)	359 (98%)	6 (2%)	62	84

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

Mol	Chain	Res	Type
1	A	64	LEU
1	A	109	LEU
1	A	229	LEU
1	A	240	PHE
1	B	64	LEU
1	B	109	LEU

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, 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	230/237 (97%)	-0.20	0 100 100	8, 34, 81, 103	0
1	B	191/237 (80%)	0.94	35 (18%) 1 0	25, 97, 146, 186	0
All	All	421/474 (88%)	0.32	35 (8%) 11 4	8, 51, 137, 186	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	208	ASN	5.4
1	B	223	PHE	4.6
1	B	111	ASP	4.5
1	B	149	ALA	4.1
1	B	185	PRO	4.1
1	B	226	ILE	4.0
1	B	156	ASP	3.8
1	B	136	ILE	3.8
1	B	225	GLU	3.8
1	B	179	TYR	3.7
1	B	227	LEU	3.6
1	B	116	SER	3.6
1	B	115	GLN	3.5
1	B	160	GLN	3.5
1	B	38	GLU	3.5
1	B	155	LYS	3.2
1	B	224	TYR	3.1
1	B	151	ILE	3.0
1	B	114	SER	3.0
1	B	183	HIS	2.9
1	B	221	ALA	2.8
1	B	139	LEU	2.7
1	B	134	TYR	2.7
1	B	188	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	215	LYS	2.5
1	B	161	SER	2.4
1	B	195	PHE	2.4
1	B	152	GLN	2.3
1	B	187	LEU	2.3
1	B	220	GLN	2.3
1	B	228	HIS	2.2
1	B	154	LEU	2.2
1	B	157	ASP	2.1
1	B	165	LEU	2.1
1	B	170	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.