



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 08:32 pm GMT

PDB ID : 1MF7
Title : INTEGRIN ALPHA M I DOMAIN
Authors : McCleverty, C.J.; Liddington, R.C.
Deposited on : 2002-08-09
Resolution : 1.25 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

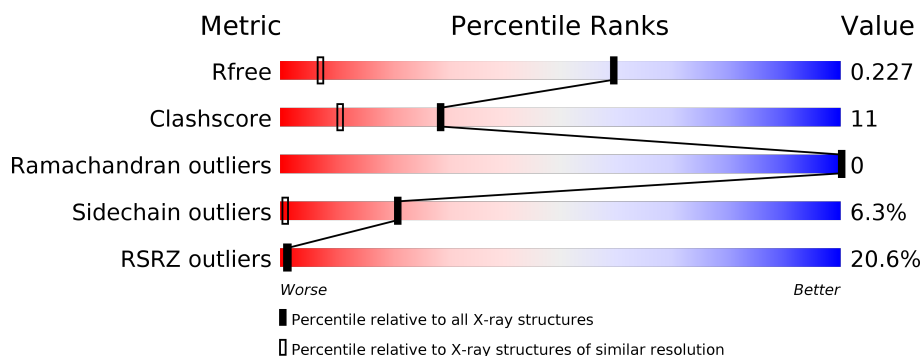
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.25 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}	100719	1690 (1.30-1.22)
Clashscore	112137	1781 (1.30-1.22)
Ramachandran outliers	110173	1716 (1.30-1.22)
Sidechain outliers	110143	1714 (1.30-1.22)
RSRZ outliers	101464	1693 (1.30-1.22)

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	194	<div> <div>21%</div> <div>79%</div> <div>19%</div> <div>.</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1806 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 INTEGRIN ALPHA M.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	194	Total	C	N	O	S	37	6	0
			1577	998	280	294	5			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	318	CYS	ALA	ENGINEERED	UNP P11215
A	320	GLY	-	CLONING ARTIFACT	UNP P11215
A	321	SER	-	CLONING ARTIFACT	UNP P11215

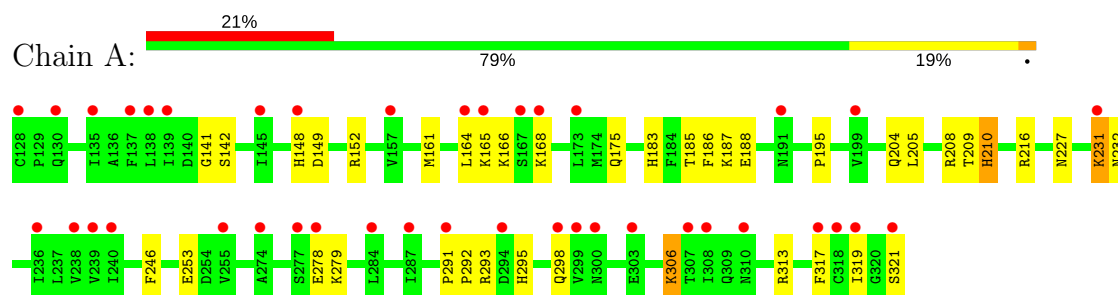
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	229	Total	O	0	0
			229	229		

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: INTEGRIN ALPHA M



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	38.64Å 51.03Å 102.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.25 24.76 – 1.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-1.25) 87.4 (24.76-1.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 1.20Å)	Xtriage
Refinement program	SHELXL	Depositor
R, R_{free}	0.166 , 0.223 0.176 , 0.227	Depositor DCC
R_{free} test set	2609 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 85.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	1806	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.82% 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 [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	0.63	0/1637	1.24	7/2203 (0.3%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	208	ARG	NE-CZ-NH2	7.98	124.29	120.30
1	A	208	ARG	CD-NE-CZ	6.64	132.89	123.60
1	A	208	ARG	NE-CZ-NH1	-5.62	117.49	120.30
1	A	148	HIS	CA-CB-CG	5.39	122.76	113.60
1	A	216	ARG	NE-CZ-NH2	-5.35	117.62	120.30
1	A	246	PHE	CB-CG-CD2	5.19	124.44	120.80
1	A	149	ASP	CB-CG-OD2	-5.12	113.69	118.30

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	1577	0	1591	34	0
2	A	229	0	0	13	0
All	All	1806	0	1591	34	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 11.

All (34) 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:141:GLY:HA3	1:A:175:GLN:HE21	1.49	0.77
1:A:165:LYS:HG3	2:A:1168:HOH:O	1.89	0.72
1:A:161:MET:HE1	1:A:195:PRO:HB3	1.70	0.71
1:A:185[B]:THR:HG23	1:A:188:GLU:H	1.57	0.70
1:A:185[B]:THR:HG21	2:A:1051:HOH:O	1.93	0.68
1:A:185[B]:THR:HG22	1:A:188:GLU:HG3	1.79	0.64
1:A:185[B]:THR:OG1	1:A:227:ASN:HB3	1.99	0.63
1:A:185[B]:THR:CG2	1:A:188:GLU:H	2.13	0.60
1:A:291:PRO:O	1:A:295:HIS:HD2	1.87	0.58
1:A:175:GLN:NE2	1:A:205:LEU:H	2.03	0.56
1:A:175:GLN:HE22	1:A:204:GLN:HA	1.70	0.56
1:A:306:LYS:HG2	2:A:1205:HOH:O	2.06	0.54
1:A:210:HIS:HD2	2:A:1047:HOH:O	1.90	0.54
1:A:168:LYS:HA	2:A:1204:HOH:O	2.07	0.54
1:A:175:GLN:HE22	1:A:205:LEU:H	1.56	0.53
1:A:210:HIS:HE1	2:A:1036:HOH:O	1.92	0.53
1:A:183:HIS:HD2	2:A:1091:HOH:O	1.93	0.50
1:A:185[B]:THR:HG22	1:A:188:GLU:CG	2.44	0.48
1:A:185[B]:THR:OG1	1:A:227:ASN:O	2.26	0.48
1:A:298:GLN:HG2	2:A:1020:HOH:O	2.13	0.48
1:A:292:PRO:HB3	2:A:1101:HOH:O	2.13	0.48
1:A:209[A]:THR:HG21	2:A:1007:HOH:O	2.14	0.47
1:A:253:GLU:HG2	2:A:1176:HOH:O	2.15	0.47
1:A:164:LEU:HD23	1:A:164:LEU:HA	1.74	0.45
1:A:232:ASN:ND2	2:A:1201:HOH:O	2.49	0.45
1:A:231:LYS:O	1:A:232:ASN:HB2	2.17	0.44
1:A:142[A]:SER:HB3	1:A:209[A]:THR:HG22	2.00	0.43
1:A:295:HIS:HE1	2:A:1170:HOH:O	2.00	0.43
1:A:185[B]:THR:OG1	1:A:186:PHE:N	2.52	0.43
1:A:161:MET:HE1	1:A:195:PRO:CB	2.45	0.43
1:A:161:MET:CE	1:A:195:PRO:HB3	2.43	0.42
1:A:161:MET:HB3	1:A:161:MET:HE2	1.78	0.42
1:A:185[B]:THR:HG23	1:A:187:LYS:N	2.35	0.42
1:A:152:ARG:NH1	1:A:152:ARG:HG3	2.35	0.41

There are no symmetry-related clashes.

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	198/194 (102%)	190 (96%)	8 (4%)	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	180/174 (103%)	169 (94%)	11 (6%)	22	1

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	166	LYS
1	A	210	HIS
1	A	231	LYS
1	A	278	GLU
1	A	279	LYS
1	A	293	ARG
1	A	306	LYS
1	A	313	ARG
1	A	317	PHE
1	A	319	ILE
1	A	321	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	175	GLN
1	A	183	HIS
1	A	191	ASN
1	A	210	HIS
1	A	227	ASN
1	A	232	ASN
1	A	295	HIS
1	A	309	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 ⓘ

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	194/194 (100%)	1.17	40 (20%) 1 1	22, 29, 44, 56	13 (6%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	167	SER	6.1
1	A	164	LEU	4.3
1	A	238	VAL	4.1
1	A	191	ASN	4.0
1	A	321	SER	4.0
1	A	303	GLU	3.8
1	A	148	HIS	3.7
1	A	145	ILE	3.7
1	A	291	PRO	3.5
1	A	240	ILE	3.4
1	A	128	CYS	3.4
1	A	231	LYS	3.3
1	A	317	PHE	3.2
1	A	236	ILE	3.2
1	A	137	PHE	3.2
1	A	318	CYS	3.0
1	A	165	LYS	3.0
1	A	307	THR	2.8
1	A	135	ILE	2.8
1	A	287	ILE	2.8
1	A	239	VAL	2.8
1	A	310	ASN	2.7
1	A	199	VAL	2.7
1	A	308	ILE	2.6
1	A	139	ILE	2.6
1	A	277	SER	2.4
1	A	294	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	255	VAL	2.4
1	A	319	ILE	2.4
1	A	278	GLU	2.4
1	A	157	VAL	2.4
1	A	298	GLN	2.3
1	A	284	LEU	2.3
1	A	274	ALA	2.3
1	A	130	GLN	2.2
1	A	299	VAL	2.2
1	A	300	ASN	2.1
1	A	138	LEU	2.1
1	A	173	LEU	2.1
1	A	168	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.