



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:34 PM UTC

PDB ID : 4DB8 / pdb_00004db8
Title : Designed Armadillo-repeat Protein
Authors : Madhurantakam, C.; Varadamsetty, G.; Grutter, M.G.; Pluckthun, A.; Mittl, P.R.E.
Deposited on : 2012-01-13
Resolution : 2.50 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

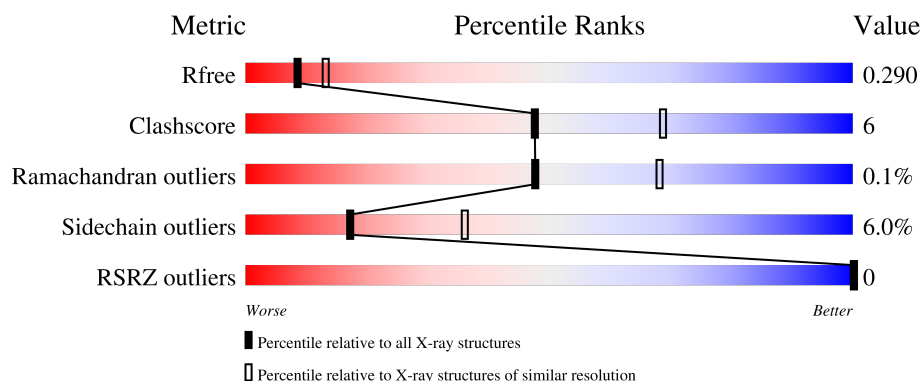
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

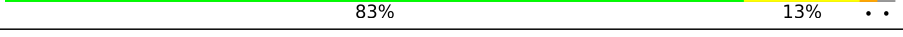
The reported resolution of this entry is 2.50 Å.

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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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	252	 83% 15% ..
1	B	252	 80% 14% • 5%
1	C	252	 86% 6% • 6%
1	D	252	 83% 13% ..

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7385 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 Armadillo-repeat Protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	247	Total	C	N	O	S	0	1	0
			1856	1151	323	380	2			
1	B	239	Total	C	N	O	S	0	1	0
			1783	1108	304	369	2			
1	C	238	Total	C	N	O	S	0	0	0
			1763	1096	299	366	2			
1	D	246	Total	C	N	O	S	0	0	0
			1845	1145	321	377	2			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	2	Total	Mg	0	0
			2	2		
2	D	1	Total	Mg	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	28	Total	O	0	0
			28	28		
3	B	25	Total	O	0	0
			25	25		
3	C	35	Total	O	0	0
			35	35		
3	D	46	Total	O	0	0
			46	46		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.00Å 113.64Å 85.59Å 90.00° 106.83° 90.00°	Depositor
Resolution (Å)	40.00 – 2.50 40.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	95.2 (40.00-2.50) 95.3 (40.00-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.232 , 0.299 0.234 , 0.290	Depositor DCC
R_{free} test set	1752 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	36.8	Xtriage
Anisotropy	0.487	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 22.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.078 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7385	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% 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

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.45	0/1883	0.91	3/2561 (0.1%)
1	B	0.44	0/1801	0.86	0/2450
1	C	0.44	0/1781	0.85	0/2426
1	D	0.43	0/1869	0.87	0/2542
All	All	0.44	0/7334	0.87	3/9979 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	26	GLN	N-CA-C	-11.66	98.86	113.55
1	A	229	LEU	N-CA-C	-5.43	107.91	114.75
1	A	25	MET	CB-CA-C	-5.31	100.21	110.35

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	1856	0	1832	32	0
1	B	1783	0	1778	20	0
1	C	1763	0	1745	16	0
1	D	1845	0	1821	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	D	1	0	0	0	0
3	A	28	0	0	7	0
3	B	25	0	0	1	0
3	C	35	0	0	2	0
3	D	46	0	0	1	0
All	All	7385	0	7176	81	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 6.

All (81) 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:247:LYS:HB2	3:A:413:HOH:O	1.71	0.88
1:C:55:LEU:HD21	1:C:80:ILE:HD13	1.61	0.82
1:C:201:TRP:HZ3	3:C:317:HOH:O	1.63	0.81
1:C:232:HIS:HD2	1:C:234:ASN:H	1.37	0.72
1:B:192:ASN:HD22	1:B:192:ASN:C	2.02	0.68
1:A:55:LEU:HD21	1:A:80:ILE:HD13	1.75	0.67
1:A:225:LYS:O	1:A:229:LEU:HB2	1.95	0.67
1:A:247:LYS:N	3:A:413:HOH:O	2.27	0.67
1:A:9:HIS:HD2	1:B:198:GLU:OE1	1.76	0.66
1:A:223:LEU:HB2	3:A:409:HOH:O	1.96	0.65
1:B:150:ASN:HB3	1:B:153:ILE:HD12	1.83	0.60
1:A:16:MET:HA	1:A:19:GLN:HE21	1.68	0.59
1:A:43:ASN:HD21	1:D:79:ASN:HD21	1.50	0.57
1:D:101:VAL:O	1:D:104:LEU:HB2	2.04	0.57
1:A:193:GLU:HA	1:A:196:LEU:HD12	1.87	0.57
1:B:192:ASN:HD22	1:B:194:GLN:H	1.52	0.57
1:A:247:LYS:CA	3:A:413:HOH:O	2.54	0.56
1:C:55:LEU:CD2	1:C:80:ILE:HD13	2.35	0.56
1:A:247:LYS:CB	3:A:413:HOH:O	2.43	0.54
1:A:24:ASP:O	1:A:25:MET:HB3	2.08	0.54
1:D:4:HIS:N	3:D:445:HOH:O	2.41	0.54
1:D:66:ASN:HD22	1:D:69:ILE:H	1.56	0.54
1:B:120:SER:HB3	1:B:160:ALA:HB2	1.91	0.53
1:A:28:GLN:O	1:A:32:THR:HG23	2.08	0.53
1:C:232:HIS:CD2	1:C:234:ASN:H	2.22	0.53
1:D:55:LEU:HD21	1:D:80:ILE:HD13	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:202:ALA:O	1:D:206:ILE:HD12	2.09	0.52
1:C:102:GLN:NE2	3:C:333:HOH:O	2.36	0.52
1:B:16:MET:HA	1:B:19:GLN:HE21	1.74	0.51
1:B:139:LEU:HB2	1:B:140:PRO:HD3	1.92	0.51
1:A:192:ASN:C	1:A:192:ASN:HD22	2.20	0.49
1:D:12:GLU:O	1:D:16:MET:HG3	2.12	0.49
1:B:230:GLN:HE22	1:B:245:LEU:HD22	1.78	0.49
1:B:19:GLN:O	1:B:22:SER:HB3	2.12	0.49
1:D:13:LEU:HA	1:D:16:MET:HE2	1.94	0.49
1:C:66:ASN:ND2	1:C:68:GLN:H	2.12	0.48
1:B:209:GLY:O	1:B:214:LYS:HE3	2.12	0.48
1:B:232:HIS:HD2	1:B:234:ASN:H	1.61	0.48
1:A:247:LYS:C	1:A:249:GLN:H	2.22	0.47
1:C:66:ASN:HD22	1:C:68:GLN:N	2.11	0.47
1:A:214:LYS:HE2	1:A:247:LYS:O	2.14	0.47
1:B:60:GLN:NE2	3:B:424:HOH:O	2.45	0.47
1:A:25:MET:HG3	1:D:69:ILE:CD1	2.45	0.47
1:B:158:LEU:HD12	1:B:198:GLU:HG3	1.96	0.47
1:A:226:LEU:HB2	1:A:245:LEU:HD13	1.96	0.46
1:A:68:GLN:HE21	1:D:33:ARG:NH2	2.14	0.46
1:A:25:MET:HA	1:A:28:GLN:HB2	1.97	0.46
1:D:113:GLN:HG3	1:D:156:GLU:HG3	1.98	0.46
1:A:25:MET:O	1:A:25:MET:CG	2.63	0.45
1:A:18:GLN:HG2	1:C:93:ASP:HA	1.97	0.45
1:D:13:LEU:HB2	1:D:14:PRO:HD3	1.98	0.45
1:A:126:GLY:H	1:A:129:GLN:HE21	1.64	0.45
1:A:186:GLN:NE2	3:A:423:HOH:O	2.49	0.45
1:B:232:HIS:CD2	1:B:234:ASN:H	2.34	0.45
1:D:246:GLU:C	1:D:248:LEU:H	2.25	0.45
1:B:235:GLU:HA	1:B:238:GLN:HB2	1.98	0.45
1:A:24:ASP:O	1:A:25:MET:CB	2.66	0.44
1:A:192:ASN:HD22	1:A:193:GLU:N	2.15	0.44
1:A:18:GLN:CG	1:C:93:ASP:HA	2.47	0.44
1:B:192:ASN:ND2	1:B:194:GLN:H	2.15	0.44
1:D:215:GLN:HA	1:D:218:LYS:HD2	2.00	0.43
1:D:192:ASN:HB3	1:D:195:ILE:HB	1.99	0.43
1:C:66:ASN:HD22	1:C:68:GLN:H	1.66	0.43
1:D:108:ASN:HD21	1:D:110:GLN:HG2	1.83	0.43
1:A:68:GLN:HE21	1:D:33:ARG:HH22	1.65	0.43
1:D:108:ASN:HD22	1:D:108:ASN:C	2.27	0.43
1:C:66:ASN:HD22	1:C:66:ASN:C	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:LEU:HD21	1:B:187:LEU:HD13	2.01	0.42
1:B:192:ASN:C	1:B:192:ASN:ND2	2.73	0.42
1:A:126:GLY:H	1:A:129:GLN:NE2	2.18	0.42
1:A:181:LEU:HD23	1:A:184:LEU:HD12	2.01	0.42
1:C:215:GLN:O	1:C:219:GLU:HB2	2.21	0.41
1:B:106:SER:HA	1:B:107:PRO:HD3	1.92	0.41
1:C:66:ASN:HD21	1:C:68:GLN:HB2	1.85	0.41
1:A:232:HIS:HD2	1:A:234:ASN:H	1.67	0.41
1:C:186:GLN:HE21	1:C:186:GLN:HA	1.86	0.41
1:D:184:LEU:HD11	1:D:202:ALA:HB3	2.03	0.41
1:A:194:GLN:NE2	3:A:412:HOH:O	2.55	0.40
1:B:247:LYS:C	1:B:249:GLN:H	2.30	0.40
1:C:112:LEU:O	1:C:116:LEU:HB2	2.22	0.40
1:D:230:GLN:HA	1:D:238:GLN:HG3	2.04	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	246/252 (98%)	241 (98%)	4 (2%)	1 (0%)	30	49
1	B	238/252 (94%)	235 (99%)	3 (1%)	0	100	100
1	C	236/252 (94%)	233 (99%)	3 (1%)	0	100	100
1	D	244/252 (97%)	241 (99%)	3 (1%)	0	100	100
All	All	964/1008 (96%)	950 (98%)	13 (1%)	1 (0%)	48	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	248	LEU

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	198/202 (98%)	192 (97%)	6 (3%)	36	64
1	B	190/202 (94%)	176 (93%)	14 (7%)	13	27
1	C	186/202 (92%)	172 (92%)	14 (8%)	12	26
1	D	196/202 (97%)	184 (94%)	12 (6%)	17	35
All	All	770/808 (95%)	724 (94%)	46 (6%)	17	36

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	LEU
1	A	60	GLN
1	A	70	LEU
1	A	110	GLN
1	A	112	LEU
1	A	246	GLU
1	B	29	LEU
1	B	34	LYS
1	B	41	ASP
1	B	74	LEU
1	B	108	ASN
1	B	112	LEU
1	B	155	GLN
1	B	158	LEU
1	B	192	ASN
1	B	196	LEU
1	B	223	LEU
1	B	224	GLU
1	B	230	GLN
1	B	233	GLU
1	C	29	LEU
1	C	44	GLU
1	C	66	ASN
1	C	67	GLU

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Mol	Chain	Res	Type
1	C	70	LEU
1	C	111	ILE
1	C	116	LEU
1	C	186	GLN
1	C	196	LEU
1	C	201	TRP
1	C	208	SER
1	C	215	GLN
1	C	219	GLU
1	C	233	GLU
1	D	29	LEU
1	D	74	LEU
1	D	108	ASN
1	D	159	TRP
1	D	171	GLN
1	D	184	LEU
1	D	185	VAL
1	D	196	LEU
1	D	213	GLN
1	D	215	GLN
1	D	219	GLU
1	D	238	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	19	GLN
1	A	21	ASN
1	A	43	ASN
1	A	60	GLN
1	A	68	GLN
1	A	79	ASN
1	A	87	GLN
1	A	89	GLN
1	A	129	GLN
1	A	152	GLN
1	A	163	ASN
1	A	169	ASN
1	A	186	GLN
1	A	192	ASN
1	A	194	GLN

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Mol	Chain	Res	Type
1	A	232	HIS
1	B	19	GLN
1	B	21	ASN
1	B	60	GLN
1	B	108	ASN
1	B	163	ASN
1	B	192	ASN
1	B	215	GLN
1	B	230	GLN
1	B	232	HIS
1	C	21	ASN
1	C	37	GLN
1	C	66	ASN
1	C	87	GLN
1	C	131	GLN
1	C	152	GLN
1	C	163	ASN
1	C	186	GLN
1	C	197	GLN
1	C	230	GLN
1	C	232	HIS
1	D	66	ASN
1	D	68	GLN
1	D	79	ASN
1	D	87	GLN
1	D	108	ASN
1	D	113	GLN
1	D	131	GLN
1	D	152	GLN
1	D	169	ASN
1	D	232	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	247/252 (98%)	-1.44	0 100 100	24, 39, 75, 78	1 (0%)
1	B	239/252 (94%)	-1.45	0 100 100	18, 36, 59, 61	1 (0%)
1	C	238/252 (94%)	-1.48	0 100 100	27, 35, 51, 54	0
1	D	246/252 (97%)	-1.30	0 100 100	24, 43, 95, 98	0
All	All	970/1008 (96%)	-1.42	0 100 100	18, 37, 73, 98	2 (0%)

There are no RSRZ outliers to report.

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MG	B	301	1/1	0.98	0.07	30,30,30,30	0
2	MG	B	302	1/1	0.98	0.03	30,30,30,30	0
2	MG	D	301	1/1	0.98	0.03	36,36,36,36	0
2	MG	A	301	1/1	0.99	0.03	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.