Full wwPDB X-ray Structure Validation Report (i)

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PDB ID	:	5EID
Title	:	AKR2A ankyrin repeat domain
Authors	:	Pu, H.; Li, H.
Deposited on	:	2015-10-29
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:

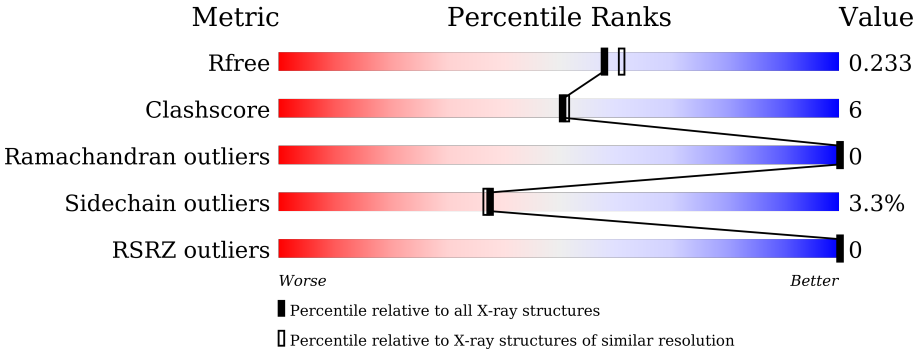
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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	A	168	62% 10% • 27%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 928 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 Ankyrin repeat domain-containing protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	123	894	556	154	181	3	0	1	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	175	MET	-	expression tag	UNP Q9SAR5
A	176	GLY	-	expression tag	UNP Q9SAR5
A	177	SER	-	expression tag	UNP Q9SAR5
A	178	SER	-	expression tag	UNP Q9SAR5
A	179	HIS	-	expression tag	UNP Q9SAR5
A	180	HIS	-	expression tag	UNP Q9SAR5
A	181	HIS	-	expression tag	UNP Q9SAR5
A	182	HIS	-	expression tag	UNP Q9SAR5
A	183	HIS	-	expression tag	UNP Q9SAR5
A	184	HIS	-	expression tag	UNP Q9SAR5
A	185	SER	-	expression tag	UNP Q9SAR5
A	186	SER	-	expression tag	UNP Q9SAR5
A	187	GLY	-	expression tag	UNP Q9SAR5
A	188	LEU	-	expression tag	UNP Q9SAR5
A	189	VAL	-	expression tag	UNP Q9SAR5
A	190	PRO	-	expression tag	UNP Q9SAR5
A	191	ARG	-	expression tag	UNP Q9SAR5
A	192	GLY	-	expression tag	UNP Q9SAR5
A	193	SER	-	expression tag	UNP Q9SAR5
A	194	HIS	-	expression tag	UNP Q9SAR5
A	195	MET	-	expression tag	UNP Q9SAR5
A	196	ALA	-	expression tag	UNP Q9SAR5
A	197	SER	-	expression tag	UNP Q9SAR5
A	198	MET	-	expression tag	UNP Q9SAR5
A	199	THR	-	expression tag	UNP Q9SAR5
A	200	GLY	-	expression tag	UNP Q9SAR5
A	201	GLY	-	expression tag	UNP Q9SAR5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	202	GLN	-	expression tag	UNP Q9SAR5
A	203	GLN	-	expression tag	UNP Q9SAR5
A	204	MET	-	expression tag	UNP Q9SAR5
A	205	GLY	-	expression tag	UNP Q9SAR5
A	206	ARG	-	expression tag	UNP Q9SAR5
A	207	GLY	-	expression tag	UNP Q9SAR5
A	208	SER	-	expression tag	UNP Q9SAR5
A	209	GLU	-	expression tag	UNP Q9SAR5
A	210	PHE	-	expression tag	UNP Q9SAR5

- Molecule 2 is water.

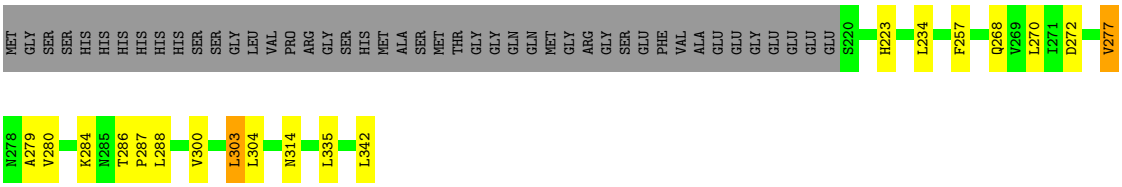
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	34	34	34	0	0

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: Ankyrin repeat domain-containing protein 2

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	124.14Å 39.19Å 28.50Å 90.00° 96.29° 90.00°	Depositor
Resolution (Å)	37.35 – 2.00 37.35 – 1.99	Depositor EDS
% Data completeness (in resolution range)	97.0 (37.35-2.00) 96.5 (37.35-1.99)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.31 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.187 , 0.235 0.189 , 0.233	Depositor DCC
R_{free} test set	437 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å^2)	22.4	Xtriage
Anisotropy	0.779	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 50.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	928	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 11.21% 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	A	0.36	0/906	0.51	0/1227

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 (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	894	0	876	10	0
2	A	34	0	0	0	0
All	All	928	0	876	10	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 (10) 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:280:VAL:HB	1:A:284:LYS:HA	1.98	0.46
1:A:279:ALA:O	1:A:287:PRO:HD3	2.17	0.44
1:A:279:ALA:O	1:A:286:THR:HA	2.18	0.44
1:A:268:GLN:NE2	1:A:272:ASP:OD1	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LEU:HD21	1:A:270:LEU:HG	2.01	0.42
1:A:277:VAL:HG22	1:A:303:LEU:HD12	2.02	0.42
1:A:284:LYS:HB3	1:A:314:ASN:HA	2.01	0.42
1:A:223:HIS:HD2	1:A:257:PHE:CE2	2.39	0.41
1:A:288:LEU:HD11	1:A:300:VAL:HG13	2.02	0.41
1:A:304:LEU:CD1	1:A:335:LEU:HD21	2.51	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	122/168 (73%)	119 (98%)	3 (2%)	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	91/131 (70%)	88 (97%)	3 (3%)	38	37

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	277	VAL
1	A	303	LEU
1	A	342	LEU

Sometimes 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 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	A	123/168 (73%)	-0.26	0	100	100	21, 29, 42, 49	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 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.