

Full wwPDB X-ray Structure Validation Report (i)

Apr 27, 2024 – 06:00 pm BST

PDB ID	:	2VFY
Title	:	AKAP18 delta central domain
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Deposited on		
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 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:

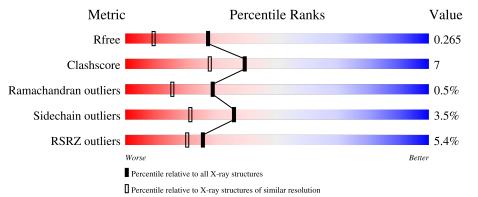
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793(1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (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 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		
			5%		
1	А	205	86%	12%	••



2VFY

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1701 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 AKAP18 DELTA.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	203	Total 1603	C 1036	N 267	O 292	S 8	0	1	0

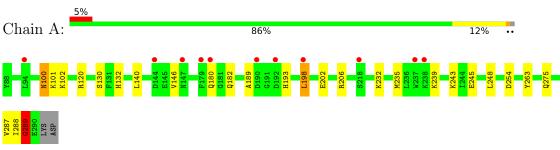
• Molecule 2 is water.

[Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
	2	А	98	Total O 98 98	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: AKAP18 DELTA



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31	Depositor
Cell constants	42.42Å 42.42 Å 96.91 Å	Denesiter
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	97.13 - 1.80	Depositor
Resolution (A)	$34.35 \ - \ 1.80$	EDS
% Data completeness	99.7 (97.13-1.80)	Depositor
(in resolution range)	99.7 (34.35 - 1.80)	EDS
R _{merge}	0.04	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.54 (at 1.81 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
D D.	0.206 , 0.238	Depositor
R, R_{free}	0.241 , 0.265	DCC
R_{free} test set	919 reflections (5.09%)	wwPDB-VP
Wilson B-factor $(Å^2)$	25.9	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.39, 44.9	EDS
L-test for twinning ²	$< L >=0.49, < L^2>=0.32$	Xtriage
	0.014 for -h,-k,l	
Estimated twinning fraction	0.067 for h,-h-k,-l	Xtriage
	0.038 for -k,-h,-l	
F_o, F_c correlation	0.94	EDS
Total number of atoms	1701	wwPDB-VP
Average B, all atoms $(Å^2)$	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.56% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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	А	0.52	0/1638	0.60	0/2207

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	А	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	288	ILE	Peptide
1	А	289	GLY	Mainchain,Peptide

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	А	1603	0	1626	24	0
2	А	98	0	0	3	0
All	All	1701	0	1626	24	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 (24) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

A + a ma 1	A + a	Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
1:A:263:TYR:CE1	1:A:287:VAL:CG1	2.78	0.66
1:A:263:TYR:CZ	1:A:287:VAL:HG11	2.32	0.64
1:A:189:ALA:O	1:A:193:HIS:HE1	1.82	0.62
1:A:263:TYR:CE1	1:A:287:VAL:HG13	2.36	0.60
1:A:263:TYR:CE1	1:A:289:GLY:HA3	2.36	0.60
1:A:100:ASN:C	1:A:100:ASN:HD22	2.06	0.59
1:A:189:ALA:O	1:A:193:HIS:CE1	2.55	0.59
1:A:263:TYR:CD1	1:A:289:GLY:HA3	2.39	0.57
1:A:202:GLU:OE1	2:A:2054:HOH:O	2.18	0.53
1:A:263:TYR:CZ	1:A:287:VAL:CG1	2.92	0.52
1:A:198:LEU:O	1:A:202:GLU:HG2	2.10	0.52
1:A:206:ARG:NH1	2:A:2055:HOH:O	2.44	0.50
1:A:130:SER:O	1:A:132:HIS:HD2	1.95	0.49
1:A:180:GLN:O	1:A:182:GLN:NE2	2.47	0.47
1:A:263:TYR:CE1	1:A:287:VAL:HG11	2.48	0.47
1:A:235:MET:O	1:A:239:LYS:HG2	2.14	0.47
1:A:275:GLN:NE2	2:A:2094:HOH:O	2.35	0.47
1:A:100:ASN:ND2	1:A:102:LYS:H	2.14	0.45
1:A:245:GLU:HB2	1:A:248:LEU:HG	1.99	0.43
1:A:100:ASN:HD21	1:A:102:LYS:HB3	1.84	0.43
1:A:100:ASN:C	1:A:100:ASN:ND2	2.71	0.42
1:A:140:LEU:HD13	1:A:146:VAL:HA	2.02	0.41
1:A:100:ASN:HD22	1:A:101:LYS:N	2.18	0.41
1:A:263:TYR:CD1	1:A:287:VAL:HG13	2.56	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	А	202/205~(98%)	198 (98%)	3~(2%)	1 (0%)	29 15

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	289	GLY

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.

Mo	l Chain	Analysed	Analysed Rotameric Outl		Percentiles
1	А	174/179~(97%)	168~(97%)	6 (3%)	37 22

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

Mol	Chain	Res	Type
1	А	100	ASN
1	А	120	ARG
1	А	198	LEU
1	А	232	LYS
1	А	243	LYS
1	А	254	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	100	ASN
1	А	112	ASN
1	А	132	HIS
1	А	139	GLN
1	А	174	HIS
1	А	193	HIS



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, 95^{th} 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 $>$	#RSI	RZ>2	$OWAB(Å^2)$	Q<0.9
1	А	203/205~(99%)	0.44	11 (5%)	25 20	24, 30, 39, 42	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	192	ASP	3.6
1	А	237	TRP	3.1
1	А	144	ASP	3.1
1	А	179	PHE	2.8
1	А	190	ASP	2.8
1	А	180	GLN	2.6
1	А	147	ASN	2.5
1	А	238	LYS	2.3
1	А	94	LEU	2.3
1	А	198	LEU	2.3
1	А	218	SER	2.3

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.

