



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 6CF6 / pdb_00006cf6
Title : RNF146 TBM-Tankyrase ARC2-3 complex
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Deposited on : 2018-02-13
Resolution : 1.93 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

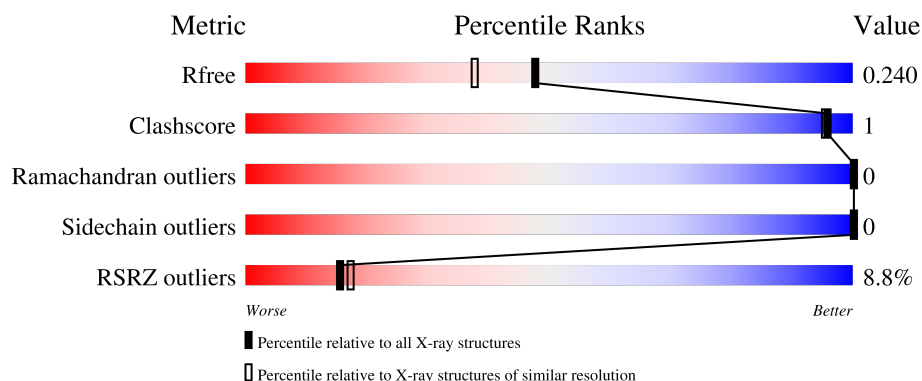
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.93 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	348	4% 89% 8%
1	B	348	10% 90% 8%
2	C	13	46% 92% 8%
2	D	13	23% 77% 23%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5338 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 Tankyrase-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	0	0
			2445	1527	450	454	14			
1	B	319	Total	C	N	O	S	0	2	0
			2453	1532	452	454	15			

- Molecule 2 is a protein called RNF146.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	0	0	0
			78	42	15	21			
2	D	10	Total	C	N	O	0	0	0
			67	36	13	18			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	157	Total	O	0	0
			157	157		
3	B	128	Total	O	0	0
			128	128		
3	C	2	Total	O	0	0
			2	2		
3	D	8	Total	O	0	0
			8	8		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	134.02Å 103.28Å 75.16Å 90.00° 106.93° 90.00°	Depositor
Resolution (Å)	50.00 – 1.93 50.00 – 1.93	Depositor EDS
% Data completeness (in resolution range)	97.0 (50.00-1.93) 97.0 (50.00-1.93)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.195 , 0.228 (Not available) , 0.240	Depositor DCC
R_{free} test set	3512 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	31.7	Xtriage
Anisotropy	0.150	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 31.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5338	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.76% 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	0.71	0/2489	0.86	0/3367
1	B	0.73	0/2501	0.83	0/3384
2	C	0.65	0/77	0.92	0/102
2	D	0.62	0/66	0.78	0/87
All	All	0.72	0/5133	0.84	0/6940

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	2445	0	2470	6	0
1	B	2453	0	2476	7	0
2	C	78	0	61	0	0
2	D	67	0	54	0	0
3	A	157	0	0	0	0
3	B	128	0	0	0	0
3	C	2	0	0	0	0
3	D	8	0	0	0	0
All	All	5338	0	5061	11	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 1.

All (11) 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:B:567:MET:CE	1:B:602:LEU:HD23	2.25	0.66
1:A:493:ALA:HB3	1:B:527[A]:HIS:HB3	1.81	0.62
1:B:567:MET:HE3	1:B:602:LEU:HD23	1.89	0.55
1:B:527[A]:HIS:CG	1:B:528:PRO:HD2	2.45	0.52
1:B:567:MET:HE2	1:B:602:LEU:HD23	1.92	0.51
1:B:441:VAL:HG12	1:B:473:LEU:HD13	1.94	0.49
1:A:493:ALA:HB3	1:B:527[B]:HIS:HB2	1.96	0.47
1:A:468:ALA:O	1:A:474:ARG:NH1	2.51	0.44
1:A:348:LEU:HD11	1:A:353:VAL:HG23	2.01	0.42
1:A:458:ASN:ND2	1:A:467:MET:HE1	2.34	0.42
1:A:562:ALA:HB2	1:A:596:HIS:CG	2.57	0.40

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	318/348 (91%)	318 (100%)	0	0	100	100
1	B	319/348 (92%)	319 (100%)	0	0	100	100
2	C	10/13 (77%)	10 (100%)	0	0	100	100
2	D	8/13 (62%)	8 (100%)	0	0	100	100
All	All	655/722 (91%)	655 (100%)	0	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	260/283 (92%)	260 (100%)	0	100	100
1	B	262/283 (93%)	262 (100%)	0	100	100
2	C	7/9 (78%)	7 (100%)	0	100	100
2	D	6/9 (67%)	6 (100%)	0	100	100
All	All	535/584 (92%)	535 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	383	GLN
1	A	439	ASN
1	A	573	HIS
1	A	585	GLN
1	B	488	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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	320/348 (91%)	0.32	14 (4%)	39 43	20, 34, 52, 78	0
1	B	319/348 (91%)	0.78	35 (10%)	10 11	14, 40, 68, 100	2 (0%)
2	C	12/13 (92%)	1.97	6 (50%)	0 0	58, 74, 80, 81	0
2	D	10/13 (76%)	1.27	3 (30%)	1 1	45, 53, 64, 65	0
All	All	661/722 (91%)	0.59	58 (8%)	15 17	14, 37, 64, 100	2 (0%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	317	PRO	6.4
1	A	315	ALA	5.5
1	B	322	VAL	5.2
1	B	318	SER	4.8
1	B	498	VAL	4.8
1	B	496	ALA	4.7
1	B	323	LEU	4.4
1	B	327	TYR	4.4
2	C	191	LEU	4.3
1	B	321	ALA	4.1
1	B	319	ALA	3.9
1	B	497	LYS	3.7
1	B	324	THR	3.5
1	B	501	THR	3.5
2	D	192	ALA	3.5
1	B	504	LEU	3.4
1	A	318	SER	3.3
1	B	503	ALA	3.2
1	A	317	PRO	3.2
1	B	481	PHE	3.2
2	C	200	ALA	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	474	ARG	2.9
1	B	316	ASP	2.9
1	B	634	GLU	2.9
1	A	327	TYR	2.8
1	B	500	LYS	2.8
1	B	489	ALA	2.8
1	B	328	LYS	2.8
1	A	323	LEU	2.8
1	B	502	LEU	2.7
2	C	192	ALA	2.7
1	B	602	LEU	2.7
1	A	321	ALA	2.7
2	C	202	SER	2.6
1	A	634	GLU	2.6
1	B	443	VAL	2.6
1	A	515	HIS	2.5
1	B	325	GLY	2.5
1	A	319	ALA	2.5
1	B	490	ALA	2.4
1	B	499	LYS	2.4
1	A	533	VAL	2.4
1	A	438	LYS	2.4
2	C	195	SER	2.3
1	B	551	PHE	2.3
1	B	495	LEU	2.3
1	A	404	SER	2.3
1	B	358	SER	2.2
2	D	201	ASP	2.2
1	B	320	LYS	2.2
2	D	195	SER	2.2
1	B	441	VAL	2.1
1	A	526	LEU	2.1
2	C	196	SER	2.1
1	B	549	LYS	2.1
1	B	357	ALA	2.1
1	B	539	ARG	2.0
1	A	326	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.