



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:09 PM UTC

PDB ID : 8QOH / pdb_00008qoh
Title : Crystal structure of the kinetoplastid kinetochore protein KKT14 C-terminal domain from *Apiculatamorpha spiralis*
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Deposited on : 2023-09-28
Resolution : 2.20 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

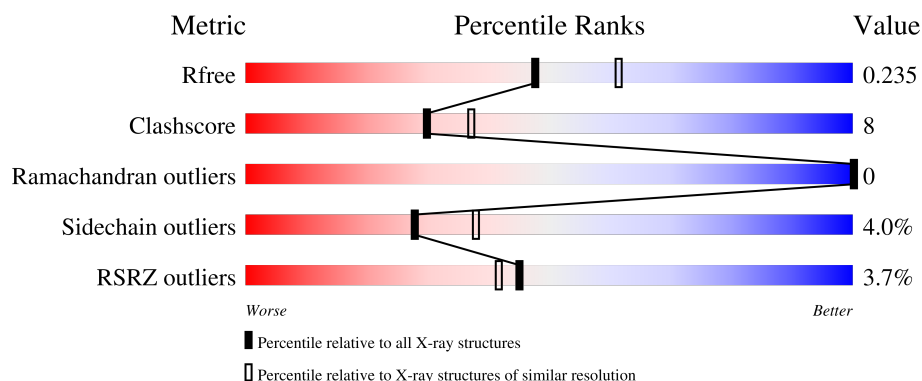
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	278	4% 82% 14% ..
1	B	278	3% 80% 16% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4575 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 kinetochore protein KKT14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	274	Total	C	N	O	S	0	0	0
			2161	1370	365	404	22			
1	A	273	Total	C	N	O	S	0	0	0
			2151	1364	362	403	22			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	117	Total	O	0	0
			117	117		
2	A	146	Total	O	0	0
			146	146		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.76Å 87.58Å 71.88Å 90.00° 93.61° 90.00°	Depositor
Resolution (Å)	43.79 – 2.20 43.79 – 2.20	Depositor EDS
% Data completeness (in resolution range)	91.9 (43.79-2.20) 92.2 (43.79-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.18.2-3874	Depositor
R, R_{free}	0.219 , 0.236 0.219 , 0.235	Depositor DCC
R_{free} test set	1356 reflections (4.33%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.979	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 24.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4575	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.02% 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

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	0.47	0/2198	0.72	0/2984
1	B	0.66	2/2209 (0.1%)	1.12	10/2999 (0.3%)
All	All	0.57	2/4407 (0.0%)	0.94	10/5983 (0.2%)

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	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	568	LEU	C-N	9.07	1.46	1.33
1	B	460	PRO	C-O	-5.30	1.19	1.24

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	568	LEU	O-C-N	17.32	134.98	121.47
1	B	567	SER	CA-C-N	13.38	146.04	121.69
1	B	567	SER	C-N-CA	13.38	146.04	121.69
1	B	568	LEU	CA-C-N	-13.30	104.04	119.47
1	B	568	LEU	C-N-CA	-13.30	104.04	119.47
1	B	567	SER	O-C-N	-11.73	106.19	122.37
1	B	429	GLY	N-CA-C	-7.90	104.45	111.95
1	B	429	GLY	CA-C-O	-6.58	117.94	122.22
1	B	428	GLU	CB-CA-C	5.55	116.63	109.80
1	B	577	LYS	N-CA-C	-5.44	106.02	113.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	568	LEU	Mainchain

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	2151	0	2096	37	1
1	B	2161	0	2106	35	1
2	A	146	0	0	12	2
2	B	117	0	0	4	1
All	All	4575	0	4202	72	3

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 8.

All (72) 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:414:ARG:HG3	1:A:414:ARG:HH11	1.34	0.93
1:A:511:THR:HG21	1:A:543:THR:HB	1.62	0.80
1:B:564:PHE:HB3	1:B:571:LEU:HD11	1.63	0.79
1:A:550:ASP:OD2	1:A:573:ARG:NH2	2.17	0.77
1:B:565:VAL:CG2	1:B:568:LEU:CD1	2.65	0.75
1:A:600:CYS:N	2:A:706:HOH:O	2.21	0.74
1:B:565:VAL:HG23	1:B:568:LEU:CD1	2.22	0.69
1:B:493:GLN:OE1	2:B:701:HOH:O	2.09	0.69
1:A:564:PHE:HB3	1:A:571:LEU:HD11	1.73	0.69
1:B:565:VAL:CG2	1:B:568:LEU:HD13	2.22	0.68
1:B:577:LYS:O	1:B:578:GLU:HG3	1.94	0.66
1:A:627:GLU:O	2:A:702:HOH:O	2.14	0.65
1:B:565:VAL:HG23	1:B:568:LEU:HD13	1.78	0.65
1:B:441:ASN:OD1	1:B:443:ASN:ND2	2.31	0.64
1:B:584:ILE:HG12	2:B:739:HOH:O	1.96	0.64
1:A:505:LEU:HD11	1:A:596:ILE:HD11	1.79	0.64
1:B:618:MET:HG3	2:B:734:HOH:O	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:565:VAL:HG22	1:B:568:LEU:CD1	2.29	0.63
1:B:546:VAL:O	1:B:548:TRP:CD1	2.53	0.62
1:A:493:GLN:CG	1:A:494:ARG:H	2.12	0.60
1:A:414:ARG:HG3	1:A:414:ARG:NH1	2.11	0.59
1:A:469:TYR:O	2:A:703:HOH:O	2.17	0.59
1:A:514:SER:OG	2:A:701:HOH:O	2.06	0.59
1:A:492:LEU:HD22	1:A:596:ILE:HD13	1.85	0.58
1:B:622:THR:O	1:B:626:THR:HG23	2.06	0.55
1:B:424:VAL:HG22	1:B:435:VAL:HG22	1.89	0.55
1:A:493:GLN:HG3	1:A:494:ARG:HG2	1.88	0.54
1:A:430:ASP:OD2	2:A:704:HOH:O	2.18	0.53
1:B:374:GLU:OE1	2:B:702:HOH:O	2.19	0.53
1:B:444:SER:HB2	1:B:549:ASP:HB3	1.90	0.53
1:B:565:VAL:HG22	1:B:568:LEU:HD13	1.91	0.52
1:B:565:VAL:HG22	1:B:568:LEU:HD11	1.91	0.52
1:A:493:GLN:CG	1:A:494:ARG:HG2	2.39	0.52
1:A:576:VAL:O	1:A:577:LYS:HB2	2.10	0.50
1:A:603:MET:HG2	1:A:627:GLU:HB2	1.93	0.50
1:A:618:MET:HG3	2:A:780:HOH:O	2.11	0.50
1:A:424:VAL:HG22	1:A:435:VAL:HG22	1.93	0.50
1:A:559:GLU:OE1	2:A:705:HOH:O	2.19	0.49
1:B:507:TYR:OH	1:B:630:TRP:NE1	2.41	0.49
1:A:637:ASN:ND2	2:A:722:HOH:O	2.46	0.48
1:A:492:LEU:HD22	1:A:596:ILE:CD1	2.43	0.48
1:B:431:GLU:HB3	1:B:432:PHE:H	1.46	0.48
1:A:538:GLY:O	1:A:636:MET:HE2	2.14	0.48
1:B:427:THR:HG22	1:B:431:GLU:HB3	1.95	0.47
1:A:603:MET:HE3	1:A:607:VAL:CG2	2.44	0.47
1:B:380:VAL:HG21	1:B:448:ALA:HB1	1.96	0.47
1:A:370:GLN:O	1:A:374:GLU:HG3	2.15	0.47
1:A:388:MET:HG2	1:A:469:TYR:CZ	2.50	0.47
1:A:551:CYS:SG	2:A:812:HOH:O	2.47	0.46
1:B:577:LYS:C	1:B:578:GLU:HG3	2.41	0.46
1:B:406:ILE:HB	1:B:428:GLU:HB2	1.98	0.45
1:A:507:TYR:O	1:A:511:THR:HG23	2.17	0.45
1:A:501:THR:HG21	1:A:596:ILE:HD12	1.99	0.44
1:A:414:ARG:HH11	1:A:414:ARG:CG	2.18	0.44
1:B:420:ARG:NH1	1:B:547:CYS:HB3	2.32	0.44
1:B:609:LYS:N	1:B:609:LYS:HD3	2.33	0.44
1:A:583:TYR:HD2	1:A:584:ILE:HD12	1.83	0.44
1:A:603:MET:O	1:A:607:VAL:HG23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:511:THR:OG1	1:B:543:THR:OG1	2.07	0.44
1:B:434:VAL:HG12	1:B:479:ILE:HG12	2.00	0.43
1:B:427:THR:HG22	1:B:431:GLU:CB	2.48	0.43
1:B:565:VAL:CG2	1:B:568:LEU:HD11	2.44	0.43
1:A:507:TYR:OH	2:A:707:HOH:O	2.21	0.43
1:A:493:GLN:CG	1:A:494:ARG:N	2.80	0.43
1:A:601:SER:HB3	1:A:604:CYS:HB2	2.01	0.43
1:B:460:PRO:HD3	1:B:542:LYS:HD3	1.99	0.42
1:A:520:ILE:HG23	2:A:812:HOH:O	2.19	0.42
1:B:369:PHE:CZ	1:B:568:LEU:HD21	2.56	0.41
1:B:375:GLU:OE2	1:B:560:ARG:NH1	2.43	0.41
1:B:488:LEU:O	1:B:492:LEU:HD13	2.21	0.41
1:A:494:ARG:NH2	2:A:723:HOH:O	2.48	0.41
1:B:469:TYR:H	1:B:469:TYR:HD1	1.70	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:577:LYS:NZ	1:A:470:ASP:OD2[2_655]	1.55	0.65
2:B:800:HOH:O	2:A:727:HOH:O[1_554]	1.97	0.23
2:A:781:HOH:O	2:A:826:HOH:O[1_455]	2.08	0.12

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	271/278 (98%)	260 (96%)	11 (4%)	0	100	100
1	B	272/278 (98%)	262 (96%)	10 (4%)	0	100	100
All	All	543/556 (98%)	522 (96%)	21 (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	236/242 (98%)	227 (96%)	9 (4%)	29	40
1	B	237/242 (98%)	227 (96%)	10 (4%)	26	36
All	All	473/484 (98%)	454 (96%)	19 (4%)	28	38

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	394	THR
1	B	410	ARG
1	B	427	THR
1	B	430	ASP
1	B	431	GLU
1	B	443	ASN
1	B	476	THR
1	B	563	LYS
1	B	609	LYS
1	B	627	GLU
1	A	371	SER
1	A	380	VAL
1	A	414	ARG
1	A	443	ASN
1	A	493	GLN
1	A	494	ARG
1	A	511	THR
1	A	578	GLU
1	A	584	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	605	GLN

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Mol	Chain	Res	Type
1	A	605	GLN
1	A	638	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 oligosaccharides 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	273/278 (98%)	0.43	12 (4%) 39 36	26, 36, 52, 68	0
1	B	274/278 (98%)	0.57	8 (2%) 53 50	25, 38, 62, 109	0
All	All	547/556 (98%)	0.50	20 (3%) 45 42	25, 37, 57, 109	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	575	ALA	5.4
1	B	429	GLY	5.1
1	B	430	ASP	4.4
1	A	368	HIS	3.9
1	B	383	VAL	3.7
1	A	640	CYS	3.5
1	A	576	VAL	3.4
1	A	575	ALA	3.4
1	B	443	ASN	3.4
1	B	546	VAL	3.3
1	A	369	PHE	3.2
1	A	562	LEU	2.9
1	A	383	VAL	2.7
1	A	639	PHE	2.7
1	A	370	GLN	2.6
1	B	495	SER	2.4
1	A	495	SER	2.4
1	A	469	TYR	2.2
1	B	599	ASN	2.1
1	A	493	GLN	2.1

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.