



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:22 PM UTC

PDB ID : 1T9H / pdb_00001t9h
Title : The crystal structure of YloQ, a circularly permuted GTPase.
Authors : Levnikov, V.M.; Blagova, E.V.; Brannigan, J.A.; Cladiere, L.; Antson, A.A.;
Isupov, M.N.; Seror, S.J.; Wilkinson, A.J.
Deposited on : 2004-05-17
Resolution : 1.60 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

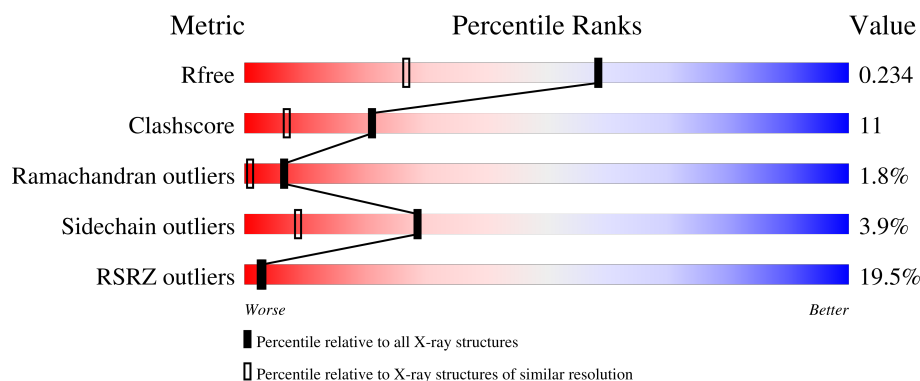
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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	307	18% 80% 10% 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	IUM	A	401	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	IUM	A	402	-	-	X	-
2	IUM	A	403	-	-	X	-
2	IUM	A	404	-	-	X	-
2	IUM	A	405	-	-	X	-
2	IUM	A	406	-	-	X	-
2	IUM	A	407	-	-	X	-
2	IUM	A	408	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 2666 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.

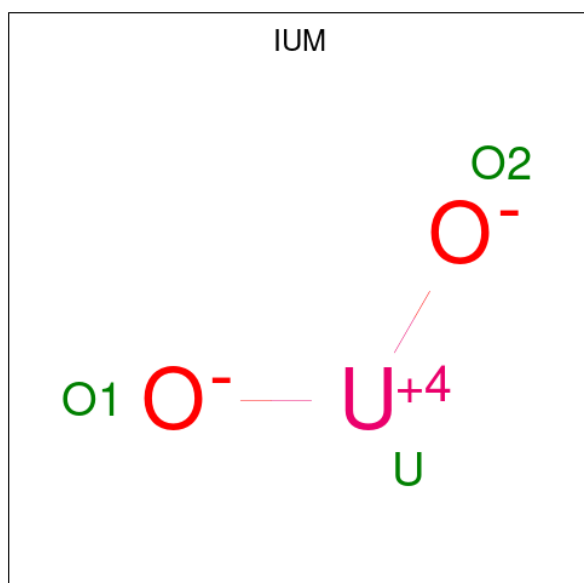
- Molecule 1 is a protein called Probable GTPase engC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	287	2322	1464	395	451	12	0	5	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	MET	-	initiating methionine	UNP O34530
A	-7	ALA	-	expression tag	UNP O34530
A	-6	ARG	-	expression tag	UNP O34530
A	-5	HIS	-	expression tag	UNP O34530
A	-4	HIS	-	expression tag	UNP O34530
A	-3	HIS	-	expression tag	UNP O34530
A	-2	HIS	-	expression tag	UNP O34530
A	-1	HIS	-	expression tag	UNP O34530
A	0	HIS	-	expression tag	UNP O34530

- Molecule 2 is URANYL (VI) ION (CCD ID: IUM) (formula: O₂U).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O U 3 2 1	0	0
2	A	1	Total O U 3 2 1	0	0
2	A	1	Total O U 3 2 1	0	0
2	A	1	Total O U 3 2 1	0	0
2	A	1	Total O U 3 2 1	0	0
2	A	1	Total O U 3 2 1	0	0
2	A	1	Total O U 3 2 1	0	0
2	A	1	Total O U 3 2 1	0	0

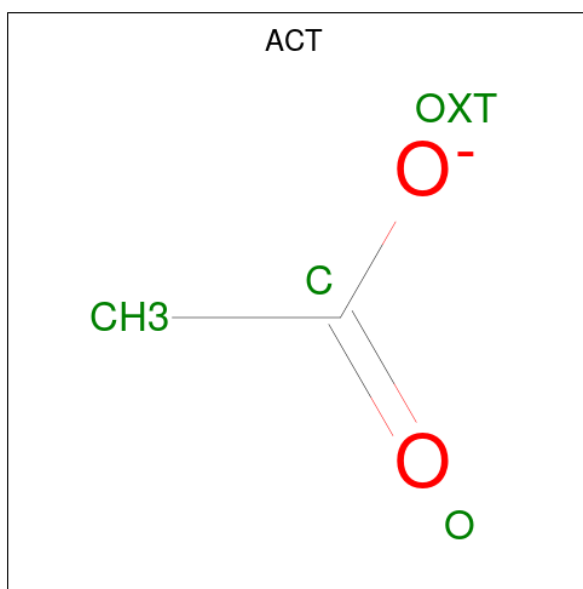
- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Zn 1 1	0	0

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total Ca 4 4	0	0

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		

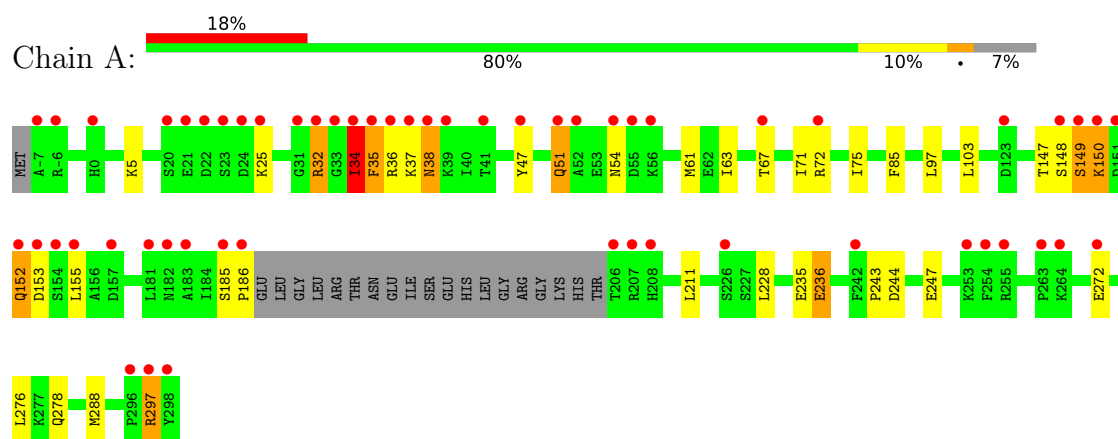
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	303	Total	O	0	0
			303	303		

3 Residue-property plots

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: Probable GTPase engC



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	47.97Å 47.97Å 279.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.50 – 1.60 20.50 – 1.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.50-1.60) 99.3 (20.50-1.60)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.1.9999	Depositor
R, R_{free}	0.144 , 0.178 0.208 , 0.234	Depositor DCC
R_{free} test set	2243 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	14.3	Xtriage
Anisotropy	0.269	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 53.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2666	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.88% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, CA, ACT, IUM

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	0.57	1/2394 (0.0%)	0.79	0/3231

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	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	34	ILE	C-N	-5.69	1.25	1.33

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	34	ILE	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	A	2322	0	2274	35	1
2	A	24	0	0	16	0
3	A	1	0	0	0	0
4	A	4	0	0	0	0
5	A	12	0	8	0	0
6	A	303	0	0	2	2
All	All	2666	0	2282	51	2

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 11.

All (51) 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:152:GLN:O	1:A:155:LEU:HD12	1.75	0.87
1:A:147:THR:HB	1:A:152:GLN:HB2	1.73	0.70
1:A:150:LYS:HG2	1:A:150:LYS:O	1.91	0.70
1:A:34:ILE:O	1:A:36:ARG:N	2.22	0.70
1:A:147:THR:CB	1:A:152:GLN:HB2	2.22	0.70
1:A:152:GLN:O	1:A:155:LEU:CD1	2.45	0.64
2:A:406:IUM:O2	2:A:406:IUM:U	1.78	0.63
2:A:407:IUM:U	2:A:407:IUM:O2	1.80	0.63
2:A:405:IUM:O1	2:A:405:IUM:U	1.80	0.62
2:A:408:IUM:O1	2:A:408:IUM:U	1.80	0.61
2:A:404:IUM:O1	2:A:404:IUM:U	1.81	0.61
2:A:401:IUM:U	2:A:401:IUM:O1	1.81	0.61
2:A:404:IUM:U	2:A:404:IUM:O2	1.80	0.61
2:A:406:IUM:U	2:A:406:IUM:O1	1.82	0.61
2:A:402:IUM:O1	2:A:402:IUM:U	1.82	0.61
2:A:401:IUM:U	2:A:401:IUM:O2	1.82	0.61
2:A:402:IUM:U	2:A:402:IUM:O2	1.81	0.61
2:A:403:IUM:O2	2:A:403:IUM:U	1.81	0.61
2:A:403:IUM:U	2:A:403:IUM:O1	1.82	0.60
2:A:408:IUM:U	2:A:408:IUM:O2	1.82	0.60
2:A:405:IUM:U	2:A:405:IUM:O2	1.81	0.60
1:A:236:GLU:H	1:A:236:GLU:CD	2.10	0.60
1:A:235:GLU:HB2	1:A:288[A]:MET:HE1	1.84	0.59
2:A:407:IUM:U	2:A:407:IUM:O1	1.82	0.59
1:A:34:ILE:HG21	1:A:61:MET:HA	1.87	0.56
1:A:244:ASP:HB2	1:A:276:LEU:CD2	2.35	0.56
1:A:297:ARG:HD3	1:A:297:ARG:H	1.73	0.54
1:A:36:ARG:HA	1:A:37:LYS:C	2.35	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:GLN:OE1	1:A:61:MET:CE	2.58	0.51
1:A:35:PHE:CE2	1:A:63:ILE:HD12	2.46	0.51
1:A:36:ARG:HA	1:A:38:ASN:N	2.25	0.51
1:A:5:LYS:HE2	1:A:47:TYR:CZ	2.46	0.50
1:A:235:GLU:CB	1:A:288[A]:MET:HE1	2.42	0.49
1:A:148:SER:O	1:A:152:GLN:HB3	2.13	0.49
1:A:67:THR:O	6:A:449:HOH:O	2.20	0.47
1:A:34:ILE:HG21	1:A:61:MET:SD	2.55	0.47
1:A:75:ILE:HD11	1:A:211:LEU:HD21	1.95	0.46
1:A:147:THR:OG1	1:A:152:GLN:HB2	2.15	0.46
1:A:103:LEU:CD1	1:A:228:LEU:HD22	2.45	0.46
1:A:278:GLN:NE2	6:A:606:HOH:O	2.49	0.46
1:A:71:ILE:HG22	1:A:72:ARG:HG3	1.98	0.45
1:A:32:ARG:CG	1:A:32:ARG:HH11	2.29	0.45
1:A:35:PHE:C	1:A:36:ARG:HG3	2.41	0.45
1:A:34:ILE:HG21	1:A:61:MET:CE	2.46	0.45
1:A:185:SER:O	1:A:186:PRO:C	2.60	0.44
1:A:243:PRO:O	1:A:247:GLU:HG2	2.18	0.43
1:A:297:ARG:H	1:A:297:ARG:CD	2.33	0.42
1:A:85:PHE:CE1	1:A:97:LEU:HD13	2.55	0.42
1:A:35:PHE:O	1:A:35:PHE:CG	2.74	0.41

All (2) 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)
6:A:433:HOH:O	6:A:552:HOH:O[1_545]	1.93	0.27
1:A:272:GLU:OE2	6:A:433:HOH:O[1_565]	2.12	0.08

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	288/307 (94%)	275 (96%)	8 (3%)	5 (2%)	7 1

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	34	ILE
1	A	35	PHE
1	A	38	ASN
1	A	149	SER
1	A	150	LYS

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	261/273 (96%)	251 (96%)	10 (4%)	29 9

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	LYS
1	A	32	ARG
1	A	34	ILE
1	A	51	GLN
1	A	54	ASN
1	A	149	SER
1	A	152	GLN
1	A	153	ASP
1	A	236	GLU
1	A	297	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN

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Mol	Chain	Res	Type
1	A	152	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 [i](#)

Of 16 ligands modelled in this entry, 5 are monoatomic - leaving 11 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	IUM	A	401	-	0,2,2	-	-	-		
2	IUM	A	403	-	0,2,2	-	-	-		
2	IUM	A	406	-	0,2,2	-	-	-		
2	IUM	A	407	-	0,2,2	-	-	-		
5	ACT	A	418	-	3,3,3	0.76	0	3,3,3	1.35	0
2	IUM	A	405	-	0,2,2	-	-	-		
2	IUM	A	402	-	0,2,2	-	-	-		
5	ACT	A	417	-	3,3,3	0.89	0	3,3,3	0.82	0
2	IUM	A	404	-	0,2,2	-	-	-		
5	ACT	A	416	-	3,3,3	1.04	0	3,3,3	1.56	1 (33%)
2	IUM	A	408	-	0,2,2	-	-	-		

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	416	ACT	OXT-C-O	-2.09	114.29	122.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	IUM	2	0
2	A	403	IUM	2	0
2	A	406	IUM	2	0
2	A	407	IUM	2	0
2	A	405	IUM	2	0
2	A	402	IUM	2	0
2	A	404	IUM	2	0
2	A	408	IUM	2	0

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	287/307 (93%)	1.30	56 (19%) 3 3	11, 23, 46, 78	5 (1%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	35	PHE	8.8
1	A	34	ILE	8.4
1	A	153	ASP	8.1
1	A	152	GLN	7.6
1	A	33	GLY	7.4
1	A	154	SER	6.7
1	A	-6	ARG	6.3
1	A	150	LYS	6.3
1	A	151	ASP	5.4
1	A	254	PHE	5.1
1	A	36	ARG	5.0
1	A	155	LEU	4.8
1	A	186	PRO	4.5
1	A	32	ARG	4.3
1	A	0	HIS	4.2
1	A	-7	ALA	4.2
1	A	182	ASN	4.1
1	A	37	LYS	3.6
1	A	206	THR	3.6
1	A	181	LEU	3.6
1	A	253	LYS	3.2
1	A	39	LYS	3.2
1	A	23	SER	3.2
1	A	38	ASN	3.2
1	A	24	ASP	3.1
1	A	298	TYR	3.1
1	A	255	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	123	ASP	2.9
1	A	149	SER	2.9
1	A	72	ARG	2.9
1	A	56	LYS	2.9
1	A	185	SER	2.8
1	A	264	LYS	2.7
1	A	54	ASN	2.7
1	A	20	SER	2.6
1	A	47	TYR	2.6
1	A	157	ASP	2.6
1	A	51	GLN	2.5
1	A	297	ARG	2.5
1	A	263	PRO	2.5
1	A	31	GLY	2.5
1	A	22	ASP	2.5
1	A	208	HIS	2.5
1	A	55	ASP	2.4
1	A	41	THR	2.4
1	A	148	SER	2.3
1	A	207	ARG	2.3
1	A	226	SER	2.3
1	A	52	ALA	2.1
1	A	21	GLU	2.1
1	A	67	THR	2.1
1	A	296	PRO	2.0
1	A	272	GLU	2.0
1	A	242	PHE	2.0
1	A	183	ALA	2.0
1	A	25	LYS	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 ⓘ

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
5	ACT	A	416	4/4	0.79	0.18	15,16,16,21	0
5	ACT	A	417	4/4	0.85	0.14	11,16,18,19	0
5	ACT	A	418	4/4	0.85	0.23	20,22,23,23	0
4	CA	A	414	1/1	0.94	0.09	44,44,44,44	0
4	CA	A	415	1/1	0.96	0.12	39,39,39,39	0
4	CA	A	413	1/1	0.96	0.14	27,27,27,27	0
2	IUM	A	405	3/3	0.99	0.09	25,25,26,27	0
2	IUM	A	406	3/3	0.99	0.11	25,25,25,29	0
2	IUM	A	407	3/3	0.99	0.06	26,26,27,28	0
3	ZN	A	411	1/1	0.99	0.03	21,21,21,21	0
2	IUM	A	404	3/3	0.99	0.09	25,25,25,26	0
2	IUM	A	401	3/3	1.00	0.07	20,20,24,24	0
2	IUM	A	408	3/3	1.00	0.11	18,18,24,25	0
2	IUM	A	402	3/3	1.00	0.11	21,21,21,23	0
4	CA	A	412	1/1	1.00	0.05	21,21,21,21	0
2	IUM	A	403	3/3	1.00	0.10	22,22,24,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.