



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:31 PM UTC

PDB ID : 4LIK / pdb_00004lik
Title : Crystal structure of the catalytic subunit of human primase
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Deposited on : 2013-07-02
Resolution : 1.70 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

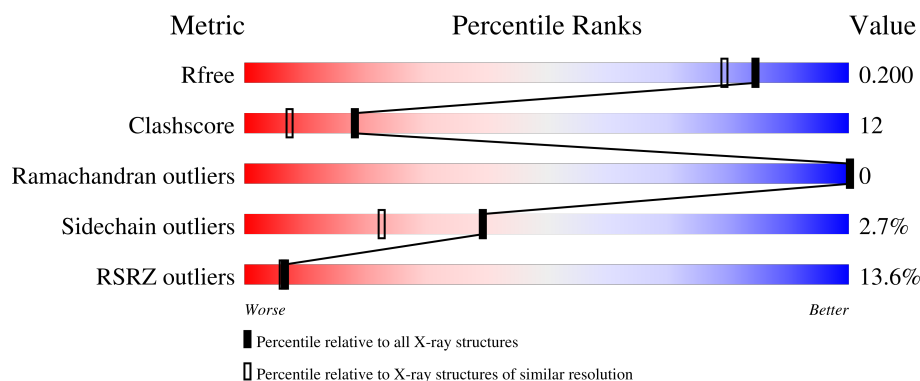
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	392	

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
3	CIT	A	502	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3514 atoms, of which 15 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 DNA primase small subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	369	3137	2022	541	559	15	0	13	0

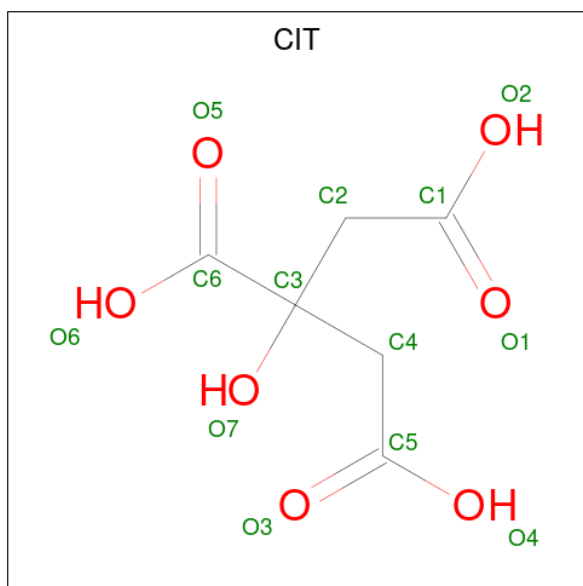
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P49642
A	-2	PRO	-	expression tag	UNP P49642
A	-1	GLY	-	expression tag	UNP P49642
A	0	SER	-	expression tag	UNP P49642
A	?	-	ILE	deletion	UNP P49642
A	?	-	SER	deletion	UNP P49642
A	?	-	THR	deletion	UNP P49642
A	?	-	ASN	deletion	UNP P49642
A	?	-	GLU	deletion	UNP P49642
A	?	-	GLU	deletion	UNP P49642
A	?	-	GLU	deletion	UNP P49642
A	?	-	LYS	deletion	UNP P49642
A	?	-	GLU	deletion	UNP P49642
A	?	-	GLU	deletion	UNP P49642
A	?	-	ASN	deletion	UNP P49642
A	?	-	GLU	deletion	UNP P49642
A	?	-	ALA	deletion	UNP P49642
A	?	-	GLU	deletion	UNP P49642
A	?	-	SER	deletion	UNP P49642
A	?	-	ASP	deletion	UNP P49642
A	?	-	VAL	deletion	UNP P49642
A	?	-	LYS	deletion	UNP P49642
A	?	-	HIS	deletion	UNP P49642
A	?	-	ARG	deletion	UNP P49642

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			18	6	5	7		
3	A	1	Total	C	H	O	0	0
			18	6	5	7		
3	A	1	Total	C	H	O	0	0
			18	6	5	7		

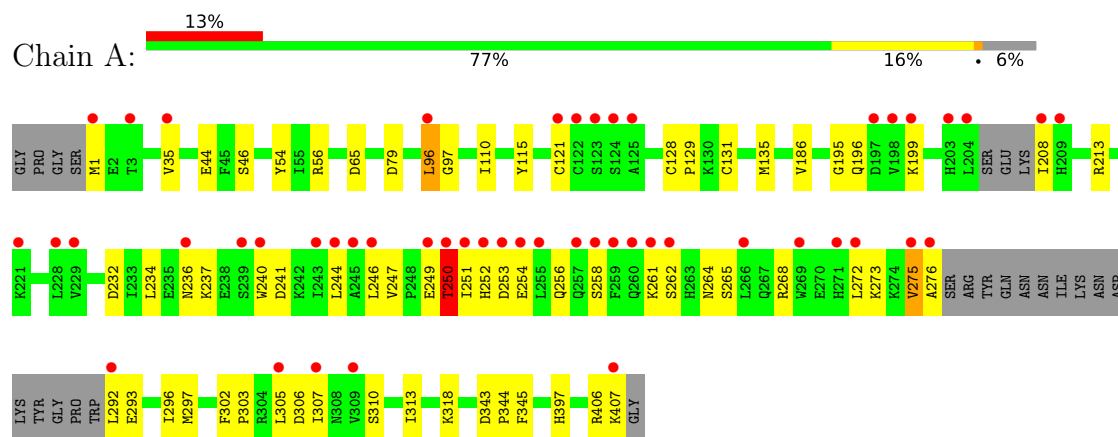
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	322	Total	O	0	0
			322	322		

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: DNA primase small subunit



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.81Å 71.92Å 84.80Å 90.00° 122.04° 90.00°	Depositor
Resolution (Å)	30.40 – 1.70 30.40 – 1.70	Depositor EDS
% Data completeness (in resolution range)	98.5 (30.40-1.70) 98.7 (30.40-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 1.70Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.175 , 0.198 0.179 , 0.200	Depositor DCC
R_{free} test set	2865 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	23.0	Xtriage
Anisotropy	0.505	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 57.6	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	3514	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% 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: CIT, ZN

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	0.48	0/3240	0.82	4/4377 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	273	LYS	N-CA-C	7.50	119.09	111.07
1	A	250	THR	N-CA-C	7.26	121.45	112.59
1	A	273	LYS	CB-CA-C	-5.95	101.53	110.88
1	A	253	ASP	N-CA-C	5.39	122.27	110.80

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	3137	0	3102	76	0
2	A	1	0	0	0	0
3	A	39	15	15	1	0
4	A	322	0	0	6	1
All	All	3499	15	3117	76	1

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 12.

All (76) 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:237:LYS:HD2	1:A:240:TRP:HE1	1.28	0.95
1:A:250:THR:C	1:A:251:ILE:HD12	1.91	0.94
1:A:237:LYS:HD2	1:A:240:TRP:NE1	1.83	0.94
1:A:237:LYS:HA	1:A:240:TRP:CE2	2.03	0.93
1:A:318:LYS:NZ	3:A:502:CIT:O5	2.04	0.89
1:A:306[B]:ASP:OD1	1:A:306[B]:ASP:N	2.06	0.88
1:A:261:LYS:HG3	1:A:262:SER:N	1.95	0.81
1:A:264:ASN:HA	1:A:268:ARG:HH12	1.48	0.78
1:A:261:LYS:CG	1:A:262:SER:N	2.46	0.77
1:A:261:LYS:CG	1:A:262:SER:H	1.98	0.77
1:A:258:SER:O	1:A:261:LYS:HG2	1.86	0.75
1:A:264:ASN:CA	1:A:268:ARG:HH12	1.99	0.74
1:A:251:ILE:HG22	1:A:251:ILE:O	1.92	0.69
1:A:292:LEU:O	1:A:296:ILE:HG12	1.92	0.68
1:A:1:MET:N	4:A:794:HOH:O	2.27	0.68
1:A:397:HIS:ND1	4:A:753:HOH:O	2.27	0.67
1:A:244:LEU:HA	1:A:247:VAL:HG22	1.79	0.63
1:A:252:HIS:O	1:A:256:GLN:CB	2.47	0.63
1:A:244:LEU:HD23	1:A:247:VAL:HG21	1.81	0.62
1:A:261:LYS:HG2	1:A:262:SER:H	1.65	0.62
1:A:305[A]:LEU:O	1:A:306[A]:ASP:CB	2.48	0.61
1:A:313:ILE:HG22	4:A:679:HOH:O	2.00	0.61
1:A:251:ILE:HD12	1:A:251:ILE:N	2.15	0.61
1:A:196:GLN:H	1:A:196:GLN:CD	2.09	0.61
1:A:258:SER:O	1:A:262:SER:N	2.34	0.60
1:A:275:VAL:CG2	1:A:276:ALA:N	2.65	0.60
1:A:96:LEU:HD12	1:A:97:GLY:N	2.18	0.59
1:A:232:ASP:OD1	1:A:265:SER:OG	2.20	0.59
1:A:240:TRP:CD1	1:A:241:ASP:N	2.71	0.59
1:A:264:ASN:C	1:A:268:ARG:NH1	2.63	0.57
1:A:240:TRP:HD1	1:A:241:ASP:N	2.03	0.57
1:A:264:ASN:CA	1:A:268:ARG:NH1	2.68	0.56
1:A:65:ASP:OD1	4:A:802:HOH:O	2.17	0.56
1:A:46[B]:SER:OG	1:A:79:ASP:HB2	2.06	0.55
1:A:251:ILE:O	1:A:251:ILE:CG2	2.54	0.55
1:A:268:ARG:HH11	1:A:268:ARG:HG3	1.72	0.53
1:A:406:ARG:O	1:A:407:LYS:HB2	2.10	0.52
1:A:240:TRP:CD1	1:A:240:TRP:C	2.89	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:208[A]:ILE:O	1:A:208[A]:ILE:HG22	2.12	0.49
1:A:258:SER:HA	1:A:261:LYS:HG2	1.95	0.49
1:A:302:PHE:CG	1:A:303:PRO:HD2	2.48	0.48
1:A:237:LYS:HD2	1:A:240:TRP:CE2	2.46	0.48
1:A:234:LEU:HB2	1:A:265:SER:HA	1.96	0.48
1:A:275:VAL:HG22	1:A:276:ALA:N	2.28	0.47
1:A:195:GLY:O	1:A:196:GLN:C	2.56	0.46
1:A:305[A]:LEU:O	1:A:306[A]:ASP:HB2	2.15	0.46
1:A:275:VAL:HG22	1:A:276:ALA:H	1.81	0.46
1:A:252:HIS:O	1:A:256:GLN:N	2.43	0.45
1:A:268:ARG:NH1	1:A:268:ARG:HG3	2.31	0.45
1:A:244:LEU:O	1:A:247:VAL:HG22	2.17	0.44
1:A:128:CYS:HB2	1:A:129:PRO:HD2	2.00	0.44
1:A:397:HIS:CE1	4:A:753:HOH:O	2.69	0.44
1:A:115:TYR:OH	1:A:303:PRO:HA	2.18	0.44
1:A:237:LYS:HA	1:A:240:TRP:CD2	2.48	0.43
1:A:258:SER:CA	1:A:261:LYS:HG2	2.48	0.43
1:A:46[A]:SER:OG	1:A:54:TYR:CE1	2.68	0.43
1:A:186:VAL:HG21	1:A:310:SER:HB2	2.00	0.43
1:A:343:ASP:OD1	1:A:345:PHE:HD2	2.02	0.43
1:A:96:LEU:HD12	1:A:96:LEU:C	2.43	0.43
1:A:293:GLU:O	1:A:297:MET:HG3	2.19	0.43
1:A:244:LEU:HD23	1:A:247:VAL:CG2	2.49	0.43
1:A:305[A]:LEU:HD23	1:A:305[A]:LEU:HA	1.58	0.42
1:A:251:ILE:O	1:A:252:HIS:C	2.62	0.42
1:A:232:ASP:C	1:A:234:LEU:H	2.28	0.42
1:A:44:GLU:OE2	1:A:56:ARG:HB2	2.20	0.41
1:A:110[B]:ILE:HG22	1:A:135:MET:HE1	2.03	0.41
1:A:213:ARG:HD2	4:A:717:HOH:O	2.20	0.41
1:A:121:CYS:SG	1:A:131:CYS:HB3	2.61	0.41
1:A:199:LYS:O	1:A:246:LEU:HD21	2.21	0.41
1:A:264:ASN:C	1:A:268:ARG:HH11	2.28	0.41
1:A:250:THR:CA	1:A:251:ILE:HD12	2.51	0.41
1:A:258:SER:C	1:A:261:LYS:HG2	2.45	0.41
1:A:305[A]:LEU:O	1:A:306[A]:ASP:HB3	2.21	0.40
1:A:343:ASP:HA	1:A:344:PRO:HD2	1.93	0.40
1:A:244:LEU:CD2	1:A:247:VAL:HG21	2.50	0.40
1:A:250:THR:O	1:A:250:THR:OG1	2.35	0.40

All (1) 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)
4:A:856:HOH:O	4:A:877:HOH:O[4_547]	2.01	0.19

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	375/392 (96%)	354 (94%)	21 (6%)	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	347/364 (95%)	338 (97%)	9 (3%)	40	23

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	VAL
1	A	96	LEU
1	A	236	ASN
1	A	249	GLU
1	A	250	THR
1	A	254	GLU
1	A	272	LEU
1	A	275	VAL

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Mol	Chain	Res	Type
1	A	307	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	58	GLN
1	A	63	GLN
1	A	341	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 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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$
3	CIT	A	504	-	12,12,12	2.21	4 (33%)	17,17,17	1.41	2 (11%)
3	CIT	A	502	-	12,12,12	2.01	4 (33%)	17,17,17	1.59	4 (23%)
3	CIT	A	503	-	12,12,12	2.49	3 (25%)	17,17,17	1.68	3 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CIT	A	504	-	-	3/16/16/16	-
3	CIT	A	502	-	-	10/16/16/16	-
3	CIT	A	503	-	-	1/16/16/16	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	503	CIT	C3-C6	-6.41	1.46	1.53
3	A	504	CIT	C4-C3	-3.87	1.49	1.54
3	A	504	CIT	C3-C6	-3.71	1.49	1.53
3	A	504	CIT	C2-C3	-3.64	1.49	1.54
3	A	502	CIT	C3-C6	-3.55	1.49	1.53
3	A	502	CIT	C2-C3	-3.26	1.49	1.54
3	A	503	CIT	C2-C3	-3.07	1.50	1.54
3	A	503	CIT	C4-C3	-2.97	1.50	1.54
3	A	502	CIT	O7-C3	-2.95	1.37	1.43
3	A	502	CIT	C4-C3	-2.64	1.50	1.54
3	A	504	CIT	O7-C3	-2.50	1.38	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	504	CIT	O6-C6-C3	3.48	119.83	113.14
3	A	502	CIT	C3-C4-C5	3.14	122.52	113.92
3	A	503	CIT	C3-C2-C1	3.10	122.41	113.92
3	A	502	CIT	C2-C3-C6	-2.71	104.04	110.03
3	A	503	CIT	O7-C3-C6	-2.66	105.18	108.96
3	A	503	CIT	O2-C1-O1	-2.31	117.39	123.33
3	A	502	CIT	C3-C2-C1	2.13	119.73	113.92
3	A	502	CIT	O4-C5-C4	2.07	120.91	114.35
3	A	504	CIT	C3-C2-C1	2.01	119.42	113.92

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	CIT	C1-C2-C3-O7

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Mol	Chain	Res	Type	Atoms
3	A	502	CIT	C1-C2-C3-C4
3	A	502	CIT	C1-C2-C3-C6
3	A	502	CIT	C2-C3-C6-O5
3	A	502	CIT	C2-C3-C6-O6
3	A	502	CIT	O7-C3-C6-O5
3	A	502	CIT	O7-C3-C6-O6
3	A	504	CIT	C1-C2-C3-O7
3	A	504	CIT	C1-C2-C3-C6
3	A	502	CIT	C6-C3-C4-C5
3	A	504	CIT	C1-C2-C3-C4
3	A	502	CIT	O7-C3-C4-C5
3	A	502	CIT	C2-C3-C4-C5
3	A	503	CIT	C1-C2-C3-O7

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	CIT	1	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	369/392 (94%)	0.41	50 (13%) 7 6	10, 25, 105, 142	19 (5%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	292	LEU	10.0
1	A	208[A]	ILE	8.8
1	A	275	VAL	7.1
1	A	259	PHE	6.2
1	A	253	ASP	5.3
1	A	307	ILE	5.2
1	A	257	GLN	4.9
1	A	198	VAL	4.8
1	A	407	LYS	4.7
1	A	229	VAL	4.6
1	A	255	LEU	4.6
1	A	251	ILE	4.2
1	A	260	GLN	3.8
1	A	276	ALA	3.7
1	A	124	SER	3.5
1	A	252	HIS	3.4
1	A	246	LEU	3.3
1	A	123	SER	3.2
1	A	271	HIS	3.0
1	A	199	LYS	3.0
1	A	244	LEU	3.0
1	A	262	SER	2.9
1	A	250	THR	2.8
1	A	197	ASP	2.7
1	A	261	LYS	2.6
1	A	96	LEU	2.6
1	A	3	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	209[A]	HIS	2.5
1	A	228	LEU	2.5
1	A	272	LEU	2.5
1	A	258	SER	2.5
1	A	1	MET	2.4
1	A	305[A]	LEU	2.4
1	A	125	ALA	2.4
1	A	249	GLU	2.4
1	A	245	ALA	2.4
1	A	122	CYS	2.3
1	A	221	LYS	2.3
1	A	240	TRP	2.3
1	A	269	TRP	2.3
1	A	35	VAL	2.3
1	A	266	LEU	2.2
1	A	203	HIS	2.2
1	A	239	SER	2.2
1	A	254	GLU	2.2
1	A	243	ILE	2.1
1	A	236	ASN	2.1
1	A	121	CYS	2.0
1	A	204	LEU	2.0
1	A	309	VAL	2.0

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](#)

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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CIT	A	502	13/13	0.48	0.30	20,20,24,24	18
3	CIT	A	503	13/13	0.79	0.20	20,20,24,24	17
3	CIT	A	504	13/13	0.88	0.18	20,20,24,24	18
2	ZN	A	501	1/1	0.97	0.05	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.