



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2024 – 12:04 PM EDT

PDB ID : 8VDS
Title : DNA Ligase 1 with nick DNA 3'rG:C
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Deposited on : 2023-12-17
Resolution : 2.79 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 : 4.02b-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

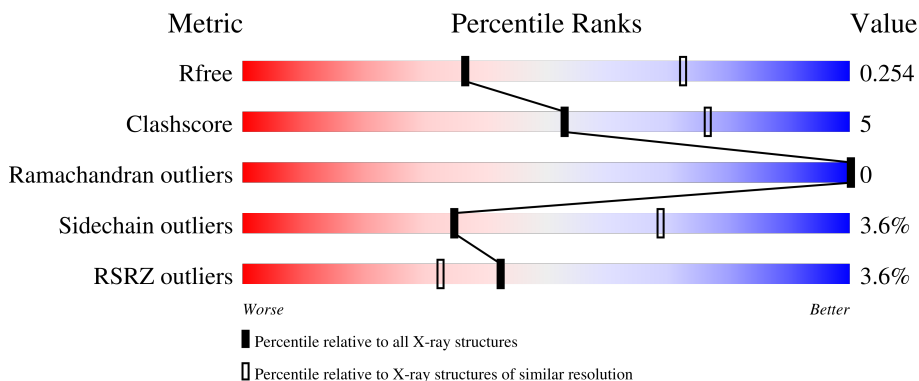
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 $\leq 5\%$. 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	658	 4% 82% 15%
2	B	18	 83% 17%
3	C	18	 100%

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 5502 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 DNA ligase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	640	4756	3033	826	882	15	0	4	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	346	ALA	GLU	conflict	UNP P18858
A	592	ALA	GLU	conflict	UNP P18858

- Molecule 2 is DNA/RNA hybrid called DNA/RNA (5'-D(*GP*CP*TP*GP*AP*TP*GP*C P*GP*T)-R(P*G)-D(P*GP*TP*CP*GP*GP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	B	18	371	176	70	108	17	0	0	0

- Molecule 3 is a DNA chain called DNA (5'-D(*GP*TP*CP*CP*GP*AP*CP*CP*AP*CP*GP*CP*AP*TP*CP*AP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	C	18	361	172	68	104	17	0	0	0

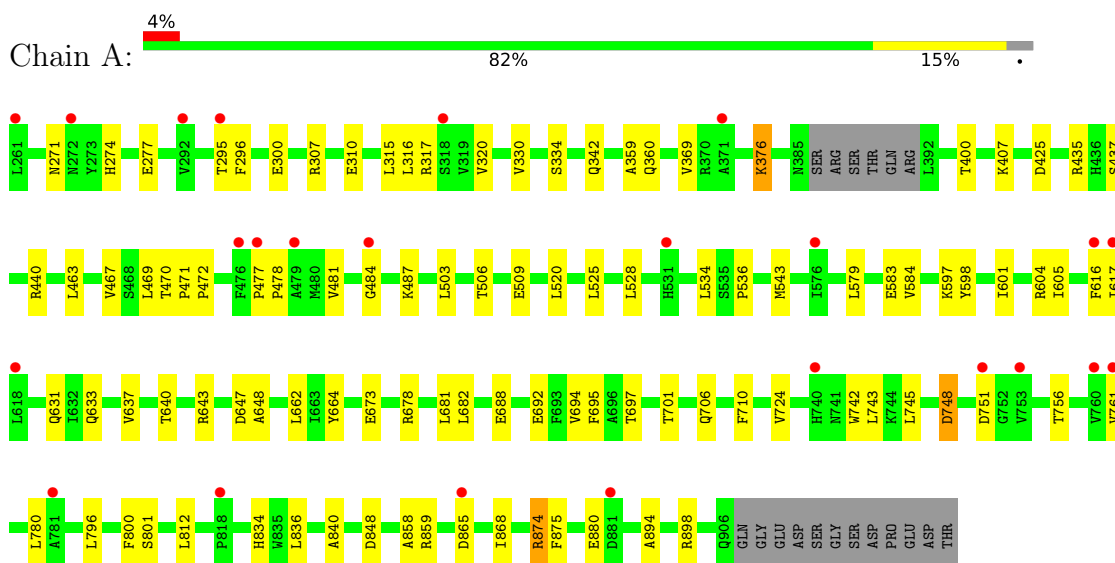
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	10	Total	O	0	0
			10	10		
4	B	1	Total	O	0	0
			1	1		
4	C	3	Total	O	0	0
			3	3		

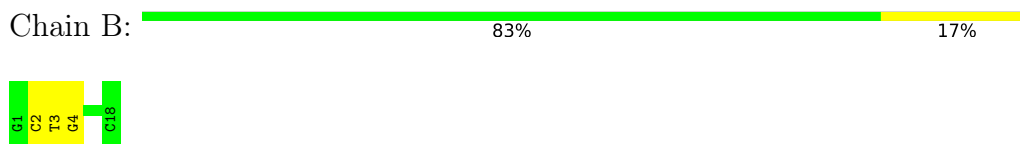
3 Residue-property plots [i](#)

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 ligase 1



- Molecule 2: DNA/RNA (5'-D(*GP*CP*TP*GP*AP*TP*GP*CP*GP*T)-R(P*G)-D(P*GP*T P*CP*GP*GP*AP*C)-3')



- Molecule 3: DNA (5'-D(*GP*TP*CP*CP*GP*AP*CP*CP*AP*CP*GP*CP*AP*TP*CP*AP *GP*C)-3')



There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.72Å 115.89Å 126.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.90 – 2.79 19.90 – 2.79	Depositor EDS
% Data completeness (in resolution range)	98.2 (19.90-2.79) 98.2 (19.90-2.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.79Å)	Xtrriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.196 , 0.252 0.198 , 0.254	Depositor DCC
R_{free} test set	1994 reflections (8.35%)	wwPDB-VP
Wilson B-factor (Å ²)	74.1	Xtrriage
Anisotropy	0.629	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 53.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5502	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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	0.24	0/4865	0.48	0/6628
2	B	0.47	0/416	0.88	0/642
3	C	0.52	0/404	0.83	0/620
All	All	0.29	0/5685	0.56	0/7890

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	4756	0	4631	55	0
2	B	371	0	204	2	0
3	C	361	0	202	0	0
4	A	10	0	0	0	0
4	B	1	0	0	0	0
4	C	3	0	0	0	0
All	All	5502	0	5037	57	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 5.

All (57) 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:597:LYS:HG3	1:A:643:ARG:HA	1.67	0.76
1:A:509:GLU:HB3	1:A:534:LEU:HB2	1.81	0.63
1:A:467:VAL:HG11	1:A:520:LEU:HD21	1.82	0.62
1:A:584:VAL:HG21	1:A:605:ILE:HG13	1.83	0.61
1:A:859:ARG:HH12	1:A:865:ASP:HA	1.69	0.57
1:A:859:ARG:HG2	1:A:859:ARG:HH11	1.69	0.56
1:A:342:GLN:OE1	1:A:342:GLN:N	2.25	0.53
1:A:697:THR:HG22	1:A:710:PHE:HZ	1.74	0.53
1:A:274:HIS:HE1	1:A:481:VAL:HG21	1.74	0.53
1:A:604:ARG:NE	1:A:692:GLU:OE1	2.42	0.52
1:A:317:ARG:HH11	1:A:469:LEU:HB3	1.73	0.52
1:A:637:VAL:O	1:A:640:THR:OG1	2.22	0.52
1:A:724:VAL:HB	1:A:743:LEU:HB2	1.92	0.52
1:A:579:LEU:HB2	1:A:583:GLU:HB3	1.90	0.52
1:A:274:HIS:HD2	1:A:277:GLU:HB2	1.76	0.51
1:A:647:ASP:OD1	1:A:648:ALA:N	2.44	0.51
1:A:682:LEU:HG	1:A:695:PHE:HE1	1.76	0.50
1:A:437:SER:HB2	1:A:440:ARG:NH1	2.27	0.50
1:A:748:ASP:HB2	1:A:874:ARG:HH22	1.77	0.50
1:A:662:LEU:HD12	1:A:682:LEU:HA	1.94	0.49
1:A:543:MET:HG2	1:A:742:TRP:CE2	2.48	0.48
1:A:274:HIS:CD2	1:A:277:GLU:HB2	2.48	0.48
1:A:598:TYR:HB3	1:A:601:ILE:HB	1.96	0.48
1:A:316:LEU:O	1:A:320:VAL:HG13	2.14	0.48
1:A:701:THR:HG21	1:A:706:GLN:HB2	1.97	0.47
1:A:673:GLU:O	1:A:678:ARG:NH1	2.48	0.47
1:A:536:PRO:HA	1:A:617:ILE:HD11	1.95	0.46
1:A:307:ARG:NH1	1:A:310:GLU:OE1	2.49	0.46
1:A:359:ALA:HB2	1:A:369:VAL:HG21	1.97	0.45
1:A:463:LEU:HD23	1:A:503:LEU:HD21	1.97	0.45
1:A:745:LEU:HD23	1:A:745:LEU:HA	1.68	0.45
1:A:274:HIS:CE1	1:A:481:VAL:HG21	2.52	0.45
1:A:376:LYS:N	1:A:376:LYS:HD2	2.32	0.45
1:A:503:LEU:HB2	1:A:525:LEU:HD11	2.00	0.44
1:A:506:THR:HG21	1:A:528:LEU:HG	1.99	0.44
1:A:780:LEU:HD11	1:A:796:LEU:HD22	1.99	0.44
1:A:471:PRO:HA	1:A:472:PRO:HD3	1.90	0.44
1:A:812:LEU:HB3	1:A:836:LEU:HD21	2.00	0.43
2:B:3:DT:H2''	2:B:4:DG:C8	2.53	0.43
1:A:330:VAL:O	1:A:334:SER:OG	2.29	0.43
1:A:536:PRO:HB2	1:A:579:LEU:HD22	2.01	0.43
1:A:484:GLY:HA2	1:A:487:LYS:HD2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:PHE:O	1:A:300:GLU:HG2	2.19	0.42
1:A:317:ARG:HD2	1:A:470:THR:OG1	2.20	0.42
1:A:761:VAL:HG22	1:A:840:ALA:HB3	2.01	0.42
1:A:407:LYS:HA	1:A:407:LYS:HD3	1.80	0.42
1:A:688:GLU:HG2	1:A:694:VAL:HA	2.01	0.42
2:B:2:DC:H6	2:B:2:DC:H5'	1.85	0.41
1:A:682:LEU:HG	1:A:695:PHE:CE1	2.54	0.41
1:A:894:ALA:O	1:A:898:ARG:HG2	2.20	0.41
1:A:477:PRO:HA	1:A:478:PRO:HD3	1.97	0.40
1:A:295:THR:OG1	1:A:315:LEU:HD12	2.21	0.40
1:A:631:GLN:OE1	1:A:631:GLN:HA	2.21	0.40
1:A:858:ALA:HB3	1:A:868:ILE:HD12	2.02	0.40
1:A:796:LEU:HD21	1:A:875:PHE:HB2	2.03	0.40
1:A:616:PHE:HA	1:A:664:TYR:O	2.21	0.40
1:A:748:ASP:HB2	1:A:874:ARG:NH2	2.37	0.40

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	A	640/658 (97%)	609 (95%)	31 (5%)	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	477/558 (86%)	460 (96%)	17 (4%)	35 69

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	271	ASN
1	A	360	GLN
1	A	376	LYS
1	A	400	THR
1	A	425	ASP
1	A	435	ARG
1	A	633	GLN
1	A	681	LEU
1	A	748	ASP
1	A	751	ASP
1	A	756	THR
1	A	800	PHE
1	A	801	SER
1	A	834	HIS
1	A	848	ASP
1	A	874	ARG
1	A	880	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	274	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

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	640/658 (97%)	0.12	24 (3%) 40 30	54, 94, 129, 159	0
2	B	18/18 (100%)	-0.83	0 100 100	56, 64, 85, 88	0
3	C	18/18 (100%)	-0.80	0 100 100	50, 62, 78, 84	0
All	All	676/694 (97%)	0.07	24 (3%) 42 32	50, 92, 129, 159	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	371	ALA	3.9
1	A	272	ASN	3.9
1	A	740	HIS	3.7
1	A	617	ILE	3.4
1	A	865	ASP	3.2
1	A	477	PRO	2.9
1	A	295	THR	2.8
1	A	760	VAL	2.7
1	A	751	ASP	2.7
1	A	618	LEU	2.5
1	A	479	ALA	2.5
1	A	616	PHE	2.4
1	A	881	ASP	2.4
1	A	761	VAL	2.3
1	A	476	PHE	2.3
1	A	261	LEU	2.2
1	A	292	VAL	2.2
1	A	781	ALA	2.2
1	A	318	SER	2.1
1	A	576	ILE	2.1
1	A	484	GLY	2.1
1	A	531	HIS	2.0
1	A	753	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	818	PRO	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 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.