



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 05:29 pm GMT

PDB ID : 1D3Y
Title : STRUCTURE OF THE DNA TOPOISOMERASE VI A SUBUNIT
Authors : Nichols, M.D.; DeAngelis, K.A.; Keck, J.L.; Berger, J.M.
Deposited on : 1999-10-01
Resolution : 2.00 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

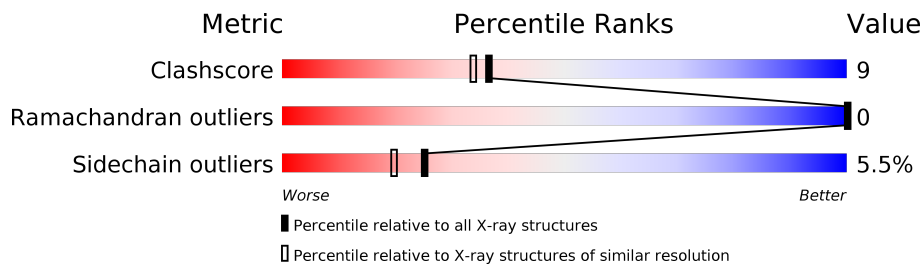
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	301	
1	B	301	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4979 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 TOPOISOMERASE VI A SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2322	1483	389	444	6			
1	B	290	Total	C	N	O	S	0	0	0
			2326	1485	391	444	6			

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Na	0	0
			1	1		
2	A	1	Total	Na	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	2	Total	Mg	0	0
			2	2		

- Molecule 4 is water.

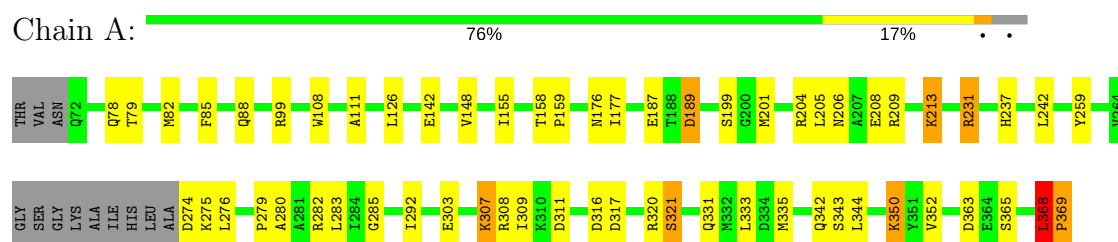
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	146	Total	O	0	0
			146	146		
4	B	180	Total	O	0	0
			180	180		

3 Residue-property plots [i](#)

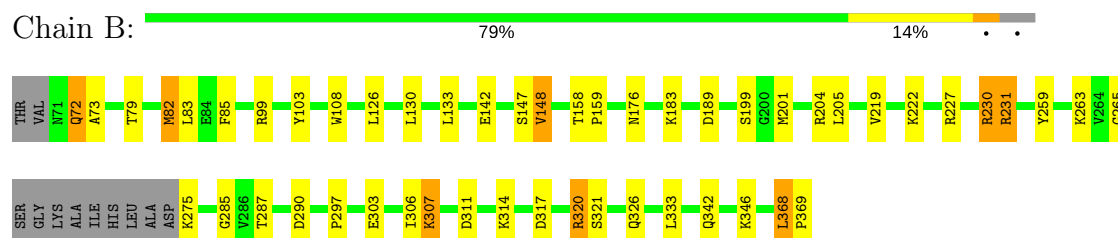
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

Note EDS was not executed.

• Molecule 1: DNA TOPOISOMERASE VI A SUBUNIT



• Molecule 1: DNA TOPOISOMERASE VI A SUBUNIT



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.51Å 59.04Å 87.26Å 90.00° 94.05° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.00)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.197 , 0.242	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4979	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, MG

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	0.37	1/2369 (0.0%)	0.88	6/3207 (0.2%)
1	B	0.36	0/2373	0.84	7/3212 (0.2%)
All	All	0.36	1/4742 (0.0%)	0.86	13/6419 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	369	PRO	N-CD	5.22	1.55	1.47

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	231	ARG	CD-NE-CZ	14.13	143.38	123.60
1	A	285	GLY	O-C-N	-12.86	102.12	122.70
1	B	285	GLY	O-C-N	-11.56	104.20	122.70
1	A	285	GLY	CA-C-N	9.65	138.42	117.20
1	B	285	GLY	CA-C-N	7.04	132.70	117.20
1	A	231	ARG	NE-CZ-NH2	6.82	123.71	120.30
1	A	231	ARG	NE-CZ-NH1	-6.46	117.07	120.30
1	B	368	LEU	C-N-CD	6.33	141.70	128.40
1	B	231	ARG	NE-CZ-NH2	-6.29	117.16	120.30
1	B	230	ARG	CD-NE-CZ	6.04	132.05	123.60
1	A	368	LEU	C-N-CD	5.90	140.79	128.40
1	B	148	VAL	CB-CA-C	-5.66	100.66	111.40
1	B	368	LEU	CB-CA-C	5.18	120.04	110.20

There are no chirality outliers.

There are no planarity outliers.

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	2322	0	2316	47	0
1	B	2326	0	2321	45	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
4	A	146	0	0	1	0
4	B	180	0	0	0	0
All	All	4979	0	4637	84	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 9.

All (84) 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:158:THR:OG1	1:B:159:PRO:HD2	1.67	0.94
1:A:201:MET:HE3	1:A:204:ARG:HB3	1.52	0.92
1:A:201:MET:HE2	1:A:205:LEU:HG	1.52	0.90
1:B:320:ARG:HA	1:B:326:GLN:HE21	1.44	0.83
1:B:201:MET:HE2	1:B:205:LEU:HG	1.61	0.82
1:B:201:MET:HE3	1:B:204:ARG:HB3	1.64	0.78
1:A:99:ARG:HH11	1:B:342:GLN:HE22	1.35	0.74
1:B:201:MET:O	1:B:201:MET:HE2	1.86	0.74
1:B:82:MET:HA	1:B:82:MET:HE3	1.69	0.74
1:B:72:GLN:HG2	1:B:73:ALA:H	1.51	0.74
1:B:158:THR:OG1	1:B:159:PRO:CD	2.35	0.73
1:A:282:ARG:HA	1:A:369:PRO:HG3	1.69	0.73
1:A:342:GLN:HE22	1:B:99:ARG:HH11	1.36	0.72
1:B:189:ASP:O	1:B:189:ASP:OD1	2.08	0.72
1:B:307:LYS:NZ	1:B:311:ASP:OD1	2.24	0.70
1:A:352:VAL:HG23	4:A:374:HOH:O	1.90	0.69
1:B:82:MET:HE2	1:B:85:PHE:HD1	1.57	0.69
1:A:201:MET:HE2	1:A:201:MET:O	1.95	0.66
1:A:331:GLN:HG2	1:A:335:MET:HE3	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:307:LYS:HE3	1:B:311:ASP:OD1	1.97	0.65
1:A:99:ARG:HD2	1:B:342:GLN:NE2	2.12	0.65
1:A:82:MET:HE3	1:A:108:TRP:HH2	1.61	0.65
1:A:237:HIS:HD2	1:A:242:LEU:O	1.79	0.64
1:A:280:ALA:O	1:A:282:ARG:HD2	1.97	0.64
1:A:307:LYS:HE3	1:A:311:ASP:OD1	1.99	0.63
1:A:158:THR:HB	1:A:159:PRO:HD2	1.82	0.61
1:B:306:ILE:HG23	1:B:333:LEU:HD11	1.83	0.61
1:A:209:ARG:O	1:A:213:LYS:HD3	2.01	0.60
1:A:82:MET:HE2	1:A:85:PHE:HD1	1.65	0.60
1:A:82:MET:HE3	1:A:82:MET:HA	1.82	0.60
1:B:307:LYS:CE	1:B:311:ASP:OD1	2.50	0.59
1:B:320:ARG:HG2	1:B:320:ARG:O	2.02	0.59
1:B:230:ARG:NH2	1:B:265:GLY:HA2	2.18	0.59
1:B:201:MET:CE	1:B:204:ARG:HB3	2.32	0.59
1:B:259:TYR:OH	1:B:369:PRO:HG2	2.03	0.58
1:A:368:LEU:O	1:A:369:PRO:O	2.22	0.58
1:A:142:GLU:O	1:B:199:SER:OG	2.22	0.57
1:B:368:LEU:O	1:B:369:PRO:OXT	2.22	0.57
1:B:297:PRO:HG3	1:B:346:LYS:HD2	1.86	0.56
1:B:82:MET:CE	1:B:85:PHE:HD1	2.18	0.56
1:B:320:ARG:HA	1:B:326:GLN:NE2	2.18	0.56
1:B:82:MET:HE3	1:B:108:TRP:HH2	1.70	0.55
1:A:259:TYR:OH	1:A:369:PRO:HG3	2.07	0.55
1:A:189:ASP:N	1:A:189:ASP:OD1	2.38	0.54
1:A:344:LEU:HD13	1:A:352:VAL:HG22	1.89	0.54
1:A:344:LEU:CD1	1:A:352:VAL:HG22	2.40	0.52
1:A:82:MET:CE	1:A:85:PHE:HD1	2.22	0.52
1:A:342:GLN:NE2	1:B:99:ARG:HD2	2.25	0.52
1:A:199:SER:OG	1:B:142:GLU:O	2.29	0.51
1:B:82:MET:HA	1:B:82:MET:CE	2.38	0.51
1:A:368:LEU:O	1:A:369:PRO:C	2.49	0.51
1:A:201:MET:CE	1:A:204:ARG:HB3	2.34	0.50
1:B:72:GLN:CG	1:B:73:ALA:H	2.23	0.50
1:A:317:ASP:O	1:A:321:SER:HB3	2.12	0.50
1:B:82:MET:HE3	1:B:108:TRP:CH2	2.47	0.50
1:B:259:TYR:OH	1:B:263:LYS:HD2	2.12	0.49
1:A:158:THR:HB	1:A:159:PRO:CD	2.43	0.49
1:A:231:ARG:HG3	1:A:276:LEU:HD13	1.94	0.48
1:A:283:LEU:H	1:A:369:PRO:CD	2.26	0.48
1:A:155:ILE:HD12	1:A:187:GLU:HG3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:316:ASP:O	1:A:320:ARG:HG3	2.13	0.47
1:A:368:LEU:C	1:A:369:PRO:O	2.51	0.47
1:A:208:GLU:OE2	1:A:350:LYS:NZ	2.34	0.47
1:A:282:ARG:HA	1:A:369:PRO:CG	2.39	0.47
1:A:99:ARG:HD2	1:B:342:GLN:HE22	1.80	0.46
1:A:78:GLN:HG2	1:A:111:ALA:O	2.16	0.46
1:B:368:LEU:HD23	1:B:368:LEU:HA	1.68	0.45
1:B:72:GLN:HG2	1:B:73:ALA:N	2.26	0.45
1:B:79:THR:HG23	1:B:126:LEU:HD21	1.99	0.45
1:A:343:SER:HA	1:B:103:TYR:CG	2.52	0.44
1:A:201:MET:CE	1:A:201:MET:O	2.62	0.44
1:A:148:VAL:CG2	1:A:177:ILE:HD11	2.48	0.44
1:B:201:MET:O	1:B:201:MET:CE	2.63	0.43
1:B:83:LEU:HD21	1:B:130:LEU:HD11	2.00	0.43
1:A:363:ASP:OD1	1:A:365:SER:OG	2.33	0.42
1:A:344:LEU:HB2	1:A:352:VAL:HG22	2.01	0.42
1:A:292:ILE:HD11	1:A:335:MET:HE2	2.02	0.41
1:B:82:MET:HE2	1:B:85:PHE:CD1	2.47	0.41
1:B:148:VAL:HG13	1:B:219:VAL:HG22	2.03	0.41
1:A:79:THR:HG23	1:A:126:LEU:HD21	2.04	0.40
1:B:72:GLN:CG	1:B:73:ALA:N	2.84	0.40
1:B:287:THR:O	1:B:290:ASP:HB2	2.21	0.40
1:A:309:ILE:HG21	1:A:333:LEU:HG	2.03	0.40
1:B:133:LEU:HD21	1:B:231:ARG:NE	2.36	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	285/301 (95%)	276 (97%)	9 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	286/301 (95%)	276 (96%)	10 (4%)	0	100	100
All	All	571/602 (95%)	552 (97%)	19 (3%)	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	253/261 (97%)	239 (94%)	14 (6%)	25	20
1	B	253/261 (97%)	239 (94%)	14 (6%)	25	20
All	All	506/522 (97%)	478 (94%)	28 (6%)	25	20

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	GLN
1	A	176	ASN
1	A	189	ASP
1	A	206	ASN
1	A	213	LYS
1	A	274	ASP
1	A	275	LYS
1	A	279	PRO
1	A	303	GLU
1	A	307	LYS
1	A	308	ARG
1	A	321	SER
1	A	350	LYS
1	A	368	LEU
1	B	72	GLN
1	B	82	MET
1	B	147	SER
1	B	176	ASN

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Mol	Chain	Res	Type
1	B	183	LYS
1	B	222	LYS
1	B	227	ARG
1	B	275	LYS
1	B	303	GLU
1	B	307	LYS
1	B	314	LYS
1	B	317	ASP
1	B	320	ARG
1	B	321	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	206	ASN
1	A	237	HIS
1	A	342	GLN
1	B	78	GLN
1	B	206	ASN
1	B	299	HIS
1	B	326	GLN
1	B	342	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.