wwPDB X-ray Structure Validation Summary Report (i)

Jun 13, 2024 – 09:59 AM EDT

PDB ID	:	1NOY
Title	:	DNA POLYMERASE (E.C.2.7.7.7)/DNA COMPLEX
Authors	:	Wang, J.; Yu, P.; Lin, T.C.; Konigsberg, W.H.; Steitz, T.A.
Deposited on	:	1996-02-16
Resolution	:	2.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (1)) were used in the production of this report:

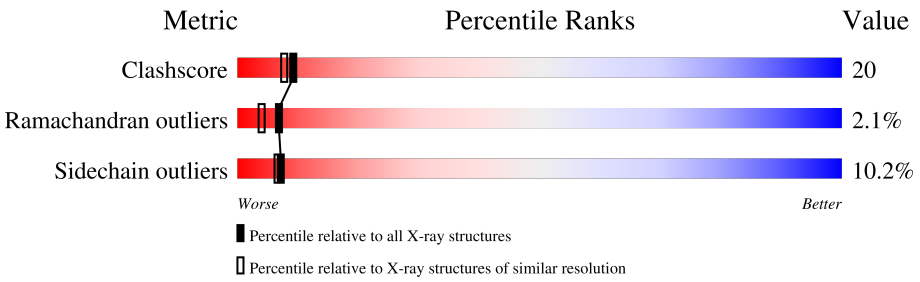
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

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))
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (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 <=5%

Mol	Chain	Length	Quality of chain
1	S	3	67% 33%
2	A	388	60% 32% • •
2	B	388	48% 35% • • 11%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 6130 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 DNA chain called DNA (5'-D(*TP*TP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	S	3	57	30	6	19	2	0	0	0

- Molecule 2 is a protein called PROTEIN (DNA POLYMERASE (E.C.2.7.7.7)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	A	372	3052	1956	497	578	21	0	0	0
2	B	346	2840	1822	466	533	19	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ASP	LYS	CONFLICT	UNP P04415
A	250	LEU	ILE	CONFLICT	UNP P04415
B	2	ASP	LYS	CONFLICT	UNP P04415
B	250	LEU	ILE	CONFLICT	UNP P04415

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mn		
4	B	1	1	1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	S	4	4	4	0	0
			Total	O		
5	A	112	112	112	0	0
			Total	O		
5	B	63	63	63	0	0

3 Residue-property plots (i)

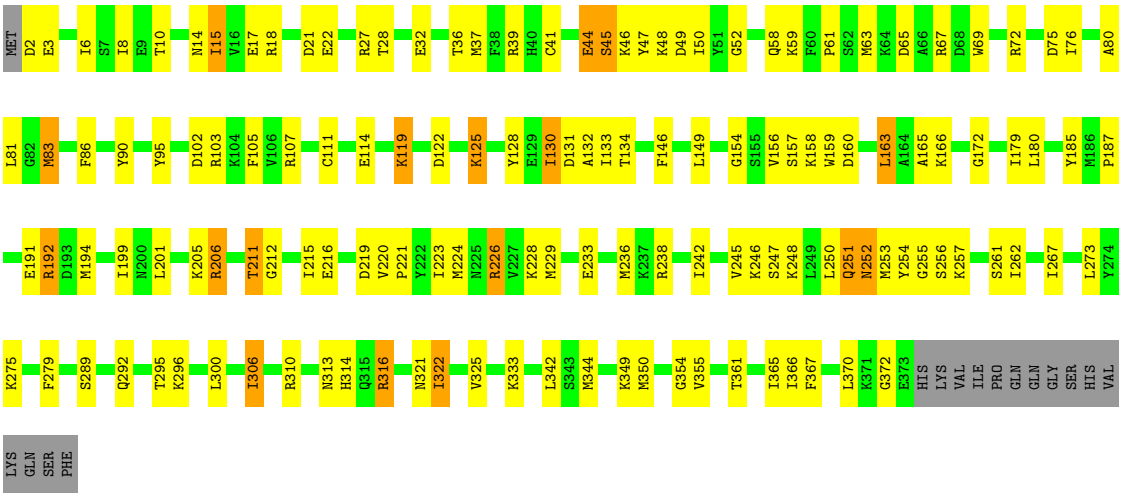
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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 (5'-D(*TP*TP*T)-3')

Chain S:

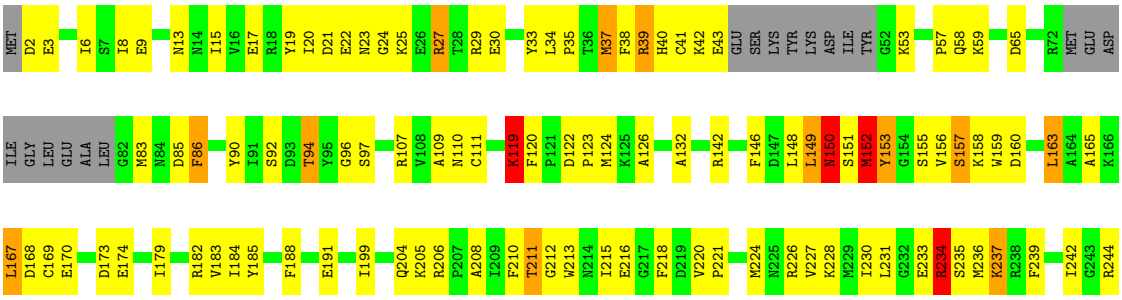
• Molecule 2: PROTEIN (DNA POLYMERASE (E.C.2.7.7.7))

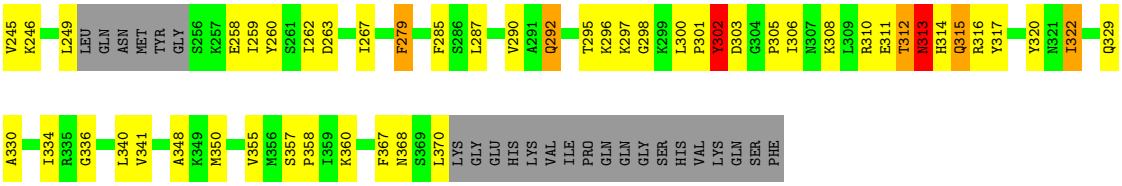
Chain A:



• Molecule 2: PROTEIN (DNA POLYMERASE (E.C.2.7.7.7))

Chain B:




V245
K246
L249
LEU
GLN
ASN
MET
TYR
GLY
S256
K257
E258
I259
Y260
S261
I262
D263
I267
F279
F285
S286
L287
V290
A291
Q292
T295
K296
K297
G298
K299
L300
P301
Y302
D303
G304
P305
I306
N307
K308
L309
R310
E311
T312
N313
H314
Q315
R316
Y317
Y320
N321
I322
Q329
A330
I334
R335
G336
L340
V341
A348
K349
M350
V355
M356
S357
P358
I359
K360
F367
N368
S369
L370
LYS
GLY
GLU
HIS
LYS
VAL
ILE
PRO
GLN
GLN
GLY
SER
HIS
VAL
LYS
GLN
SER
PHE

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	114.70Å 109.66Å 70.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.20 59.16 – 2.49	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-2.20) 79.8 (59.16-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.75 (at 2.48Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.222 , (Not available) (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	45.1	Xtriage
Anisotropy	0.440	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 87.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	6130	wwPDB-VP
Average B, all atoms (Å^2)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.91% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: MN, 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	S	3.81	8/62 (12.9%)	4.48	20/94 (21.3%)
2	A	0.54	0/3122	0.73	0/4206
2	B	0.49	0/2904	0.71	0/3911
All	All	0.64	8/6088 (0.1%)	0.86	20/8211 (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	S	0	1
2	A	1	0
2	B	1	0
All	All	2	1

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	S	3	DT	C5-C7	13.34	1.58	1.50
1	S	5	DT	C5-C7	12.16	1.57	1.50
1	S	4	DT	C5'-C4'	9.10	1.61	1.51
1	S	4	DT	C5-C7	8.99	1.55	1.50
1	S	3	DT	C5-C6	7.29	1.39	1.34

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	S	3	DT	OP2-P-O3'	14.82	137.80	105.20
1	S	4	DT	O5'-P-OP1	-14.27	92.86	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	S	4	DT	O3'-P-O5'	-13.68	78.01	104.00
1	S	4	DT	OP2-P-O3'	12.91	133.60	105.20
1	S	4	DT	O5'-P-OP2	11.74	124.79	110.70

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	211	THR	CB
2	B	211	THR	CB

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	S	5	DT	Sidechain

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	S	57	0	38	12	0
2	A	3052	0	2993	102	0
2	B	2840	0	2783	135	0
3	B	1	0	0	0	0
4	B	1	0	0	1	0
5	A	112	0	0	3	0
5	B	63	0	0	6	0
5	S	4	0	0	3	0
All	All	6130	0	5814	234	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 20.

The worst 5 of 234 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:S:4:DT:H4'	5:S:108:HOH:O	1.42	1.18

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:3:DT:O4'	5:S:116:HOH:O	1.67	1.13
1:S:5:DT:O2	2:B:119:LYS:NZ	1.84	1.10
2:B:119:LYS:HA	2:B:119:LYS:HZ3	1.15	1.08
2:A:37:MET:SD	2:A:372:GLY:HA2	2.08	0.93

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
2	A	370/388 (95%)	344 (93%)	24 (6%)	2 (0%)	29 31
2	B	338/388 (87%)	294 (87%)	31 (9%)	13 (4%)	3 1
All	All	708/776 (91%)	638 (90%)	55 (8%)	15 (2%)	7 4

5 of 15 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	45	SER
2	B	22	GLU
2	B	152	MET
2	B	313	ASN
2	B	150	ASN

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
2	A	335/350 (96%)	303 (90%)	32 (10%)	8 8
2	B	312/350 (89%)	278 (89%)	34 (11%)	6 5
All	All	647/700 (92%)	581 (90%)	66 (10%)	7 6

5 of 66 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	279	PHE
2	B	292	GLN
2	B	322	ILE
2	A	289	SER
2	A	254	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	292	GLN
2	B	313	ASN
2	B	321	ASN
2	A	313	ASN
2	A	321	ASN

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)

Of 2 ligands modelled in this entry, 2 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 (i)

6.1 Protein, DNA and RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.