



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:17 PM GMT

PDB ID : 1JN3  
Title : FIDELITY PROPERTIES AND STRUCTURE OF M282L MUTATOR MUTANT OF DNA POLYMERASE: SUBTLE STRUCTURAL CHANGES INFLUENCE THE MECHANISM OF NUCLEOTIDE DISCRIMINATION  
Authors : Conn, D.A.; Sweasy, J.B.; Jaeger, J.  
Deposited on : 2001-07-22  
Resolution : 2.35 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](mailto: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  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

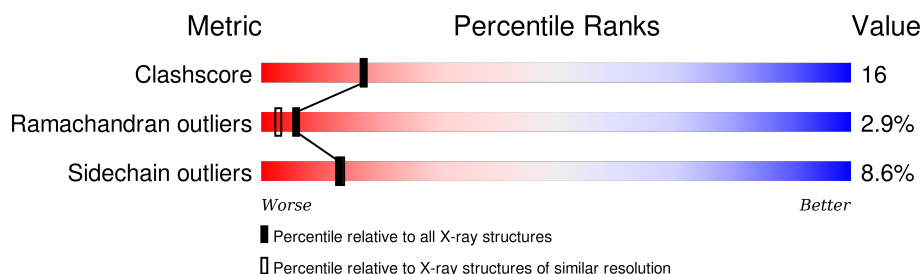
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 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	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)

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  $\geq 3$ , 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	251	 69% 22% 5% •

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2154 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 POLYMERASE BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	0	0
			1969	1239	345	378	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	282	LEU	MET	ENGINEERED	UNP P06766

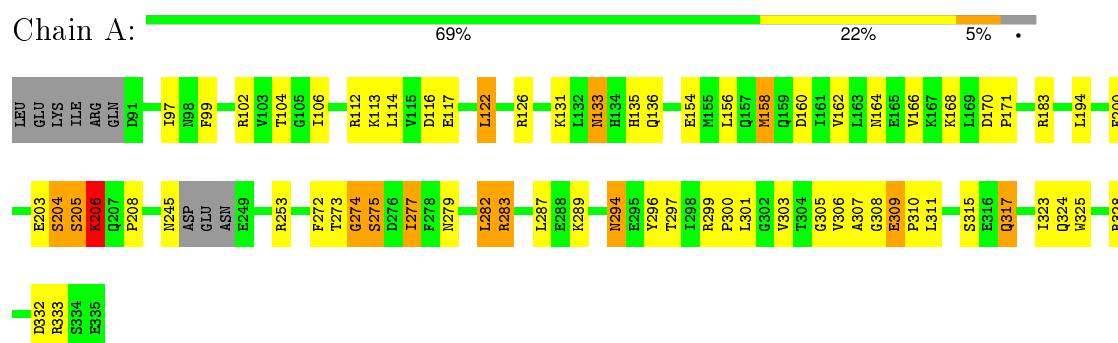
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	185	Total	O	0	0
			185	185		



Note EDS was not executed.

- Molecule 1: DNA POLYMERASE BETA



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	119.54Å 63.62Å 37.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.48 – 2.35	Depositor
% Data completeness (in resolution range)	93.4 (22.48-2.35)	Depositor
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.215 , 0.256	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2154	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

## 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	0.48	0/2007	0.72	0/2703

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	1969	0	1949	63	0
2	A	185	0	0	5	0
All	All	2154	0	1949	63	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 16.

All (63) 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:294:ASN:ND2	1:A:296:TYR:H	1.73	0.86
1:A:299:ARG:HH11	1:A:299:ARG:HG3	1.42	0.84
1:A:279:ASN:O	1:A:283:ARG:HG2	1.78	0.84
1:A:317:GLN:NE2	1:A:317:GLN:H	1.84	0.76

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:THR:HG22	1:A:274:GLY:H	1.50	0.75
1:A:294:ASN:HD22	1:A:296:TYR:H	1.34	0.73
1:A:194:LEU:HD21	1:A:272:PHE:CD1	2.24	0.73
1:A:317:GLN:HE21	1:A:317:GLN:H	1.37	0.72
1:A:282:LEU:HD12	1:A:282:LEU:O	1.90	0.72
1:A:294:ASN:C	1:A:294:ASN:HD22	1.93	0.71
1:A:113:LYS:O	1:A:117:GLU:HG2	1.91	0.69
1:A:205:SER:O	1:A:206:LYS:HB3	1.93	0.68
1:A:133:ASN:ND2	1:A:136:GLN:H	1.91	0.67
1:A:289:LYS:HA	1:A:289:LYS:HE2	1.76	0.67
1:A:299:ARG:NH1	1:A:299:ARG:HG3	2.12	0.62
1:A:122:LEU:C	1:A:122:LEU:HD12	2.19	0.62
1:A:315:SER:HB2	1:A:317:GLN:NE2	2.15	0.62
1:A:194:LEU:HD21	1:A:272:PHE:HD1	1.63	0.62
1:A:274:GLY:O	1:A:275:SER:HB3	1.99	0.62
1:A:277:ILE:H	1:A:277:ILE:HD13	1.63	0.61
1:A:282:LEU:HD12	1:A:282:LEU:C	2.24	0.58
1:A:274:GLY:HA2	1:A:333:ARG:HB3	1.87	0.56
1:A:303:VAL:C	1:A:305:GLY:H	2.07	0.56
1:A:104:THR:HG22	2:A:337:HOH:O	2.04	0.56
1:A:133:ASN:H	1:A:136:GLN:NE2	2.04	0.55
1:A:300:PRO:HD3	1:A:311:LEU:HD21	1.87	0.55
1:A:205:SER:O	1:A:206:LYS:HD3	2.06	0.55
1:A:154:GLU:O	1:A:158:MET:HG2	2.07	0.54
1:A:204:SER:HB3	1:A:206:LYS:HZ1	1.73	0.54
1:A:299:ARG:NE	2:A:338:HOH:O	2.22	0.53
1:A:183:ARG:HD3	1:A:274:GLY:O	2.09	0.53
1:A:273:THR:HG22	1:A:274:GLY:N	2.21	0.51
1:A:301:LEU:O	1:A:307:ALA:HB3	2.10	0.50
1:A:300:PRO:HD3	1:A:311:LEU:CD2	2.42	0.49
1:A:162:VAL:O	1:A:166:VAL:HG23	2.12	0.49
1:A:97:ILE:HD13	1:A:112:ARG:HG2	1.93	0.49
1:A:133:ASN:C	1:A:133:ASN:HD22	2.16	0.49
1:A:168:LYS:HB2	2:A:419:HOH:O	2.13	0.48
1:A:205:SER:O	1:A:206:LYS:CB	2.61	0.47
1:A:315:SER:HB2	1:A:317:GLN:HE21	1.78	0.47
1:A:122:LEU:O	1:A:126:ARG:HG3	2.14	0.46
1:A:282:LEU:CD1	1:A:323:ILE:HD13	2.45	0.46
1:A:283:ARG:NH1	2:A:383:HOH:O	2.47	0.46
1:A:294:ASN:ND2	1:A:297:THR:H	2.13	0.46
1:A:183:ARG:CD	1:A:274:GLY:O	2.64	0.45

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:PHE:O	1:A:102:ARG:HB2	2.17	0.45
1:A:106:ILE:HG12	1:A:136:GLN:HG2	1.97	0.45
1:A:294:ASN:ND2	1:A:294:ASN:C	2.67	0.45
1:A:200:PHE:HE1	1:A:205:SER:HA	1.82	0.45
1:A:289:LYS:CA	1:A:289:LYS:HE2	2.47	0.44
1:A:277:ILE:N	1:A:277:ILE:HD13	2.31	0.44
1:A:206:LYS:O	1:A:208:PRO:HD3	2.18	0.44
1:A:308:GLY:O	1:A:309:GLU:CB	2.66	0.43
1:A:299:ARG:NH2	2:A:338:HOH:O	2.51	0.43
1:A:328:ARG:HG2	1:A:332:ASP:HB2	2.00	0.43
1:A:279:ASN:HB3	1:A:283:ARG:NH1	2.34	0.42
1:A:282:LEU:HD13	1:A:323:ILE:CD1	2.49	0.42
1:A:164:ASN:HD21	1:A:168:LYS:HE3	1.85	0.42
1:A:160:ASP:O	1:A:164:ASN:HB3	2.21	0.41
1:A:309:GLU:HA	1:A:310:PRO:HD2	1.89	0.41
1:A:133:ASN:HD21	1:A:135:HIS:HB3	1.87	0.40
1:A:301:LEU:HD12	1:A:301:LEU:HA	1.98	0.40
1:A:170:ASP:HA	1:A:171:PRO:HD2	1.96	0.40

There are no symmetry-related clashes.

## 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	238/251 (95%)	221 (93%)	10 (4%)	7 (3%)	<b>6</b> <b>3</b>

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	204	SER
1	A	206	LYS
1	A	309	GLU

Continued on next page...

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	275	SER
1	A	274	GLY
1	A	306	VAL
1	A	205	SER

### 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	220/229 (96%)	201 (91%)	19 (9%)	13	13

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	LEU
1	A	116	ASP
1	A	122	LEU
1	A	131	LYS
1	A	133	ASN
1	A	156	LEU
1	A	158	MET
1	A	203	GLU
1	A	206	LYS
1	A	245	ASN
1	A	253	ARG
1	A	277	ILE
1	A	282	LEU
1	A	283	ARG
1	A	287	LEU
1	A	294	ASN
1	A	317	GLN
1	A	324	GLN
1	A	325	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	ASN
1	A	136	GLN
1	A	164	ASN
1	A	217	GLN
1	A	279	ASN
1	A	281	ASN
1	A	294	ASN
1	A	317	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](#)

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.