



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 29, 2024 – 04:11 PM EDT

PDB ID : 3U7F  
Title : Crystal structure of mPNKP catalytic fragment (D170A) bound to single-stranded DNA (TCCTCp)  
Authors : Coquelle, N.; Havali, Z.; Bernstein, N.; Green, R.; Glover, J.N.M.  
Deposited on : 2011-10-13  
Resolution : 1.80 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

<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>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

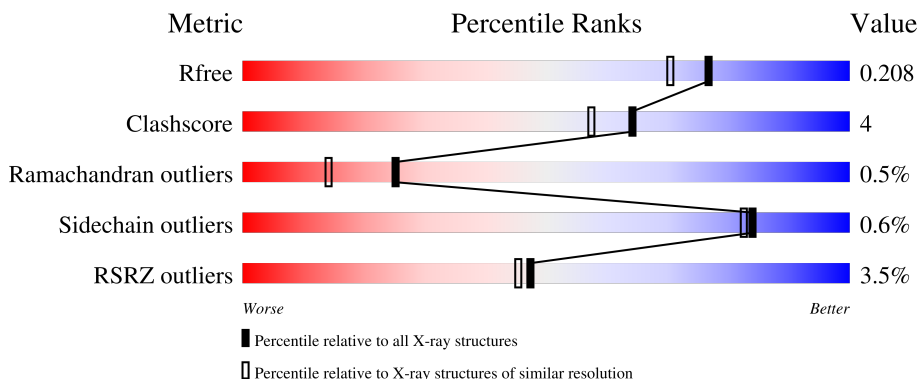
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

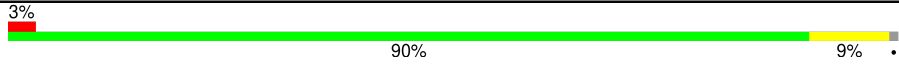

The reported resolution of this entry is 1.80 Å.

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))
$R_{free}$	164625	7108 (1.80-1.80)
Clashscore	180529	8162 (1.80-1.80)
Ramachandran outliers	177936	8077 (1.80-1.80)
Sidechain outliers	177891	8076 (1.80-1.80)
RSRZ outliers	164620	7108 (1.80-1.80)

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  $\geq 3$ , 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	B	381	
2	X	5	

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6427 atoms, of which 2998 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 Bifunctional polynucleotide phosphatase/kinase.

Mol	Chain	Residues	Atoms								ZeroOcc	AltConf	Trace
1	B	379	Total	C	H	N	O	S	Se		0	4	0
			5890	1893	2943	506	534	8	6				

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	170	ALA	ASP	engineered mutation	UNP Q9JLV6

- Molecule 2 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
2	X	2	Total	C	H	N	O	P		0	0	0
			62	19	23	5	13	2				

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

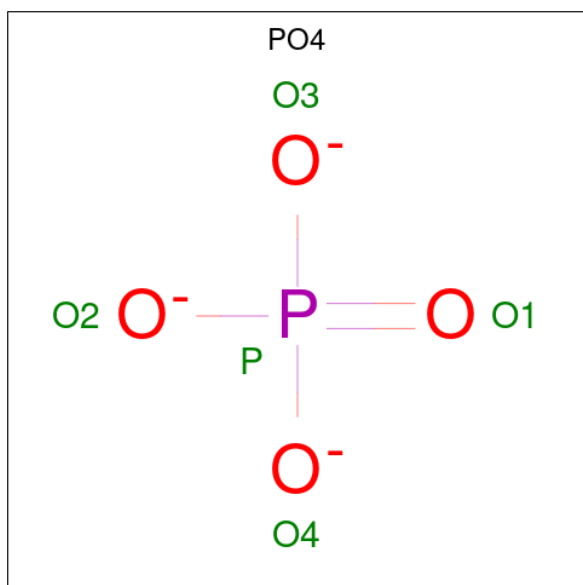
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula:  $\text{O}_4\text{P}$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	P	0	0
			5	4	1		
5	X	1	Total	O	P	0	0
			4	3	1		

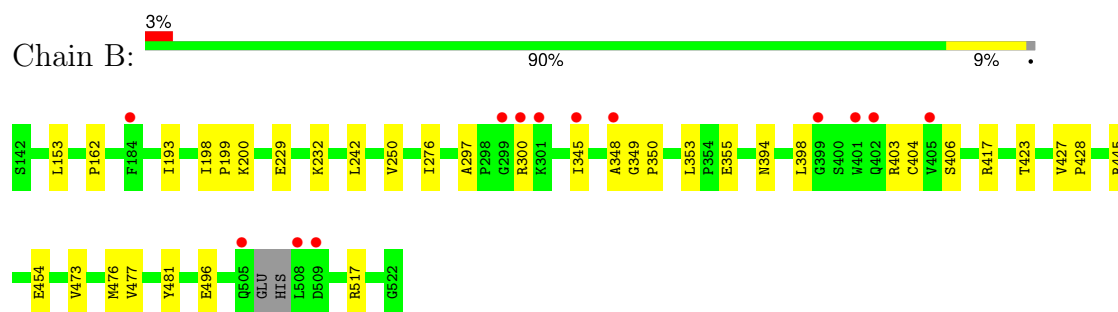
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	403	Total	O	0	0
			403	403		
6	X	6	Total	O	0	0
			6	6		

### 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: Bifunctional polynucleotide phosphatase/kinase



- Molecule 2: DNA



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	42.60Å 62.33Å 67.97Å 90.00° 91.81° 90.00°	Depositor
Resolution (Å)	45.93 – 1.80 45.93 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.93-1.80) 99.8 (45.93-1.80)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.03 (at 1.79Å)	Xtriage
Refinement program	PHENIX 1.7.2_869	Depositor
R, $R_{free}$	0.169 , 0.214 0.163 , 0.208	Depositor DCC
$R_{free}$ test set	1650 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	7.7	Xtriage
Anisotropy	0.090	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.44 , 48.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.034 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6427	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.40% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL, PO4, MG, CSO

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	B	0.42	0/3022	0.56	0/4093
2	X	1.86	1/42 (2.4%)	1.99	1/62 (1.6%)
All	All	0.47	1/3064 (0.0%)	0.61	1/4155 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	X	401	DT	P-O5'	-7.87	1.51	1.59

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	X	401	DT	N3-C2-O2	-5.49	119.00	122.30

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	B	2947	2943	2937	23	0
2	X	39	23	23	0	0
3	B	1	0	0	0	0
4	B	24	32	32	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	5	0	0	0	0
5	X	4	0	0	0	0
6	B	403	0	0	6	2
6	X	6	0	0	0	0
All	All	3429	2998	2992	23	2

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 4.

All (23) 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:300:ARG:NH1	6:B:795:HOH:O	2.32	0.62
1:B:394:ASN:O	1:B:398:LEU:HD13	2.01	0.60
1:B:355:GLU:OE2	6:B:707:HOH:O	2.15	0.60
1:B:398:LEU:O	1:B:403:ARG:HD3	2.06	0.56
1:B:297:ALA:O	1:B:300:ARG:HG2	2.09	0.52
1:B:454:GLU:CD	6:B:804:HOH:O	2.48	0.51
1:B:350:PRO:HG3	1:B:353:LEU:HD21	1.95	0.48
1:B:398:LEU:HD11	6:B:677:HOH:O	2.14	0.48
1:B:476[A]:MSE:HG2	6:B:49:HOH:O	2.14	0.47
1:B:398:LEU:HD23	1:B:404:CYS:HA	1.97	0.46
1:B:345:ILE:O	1:B:345:ILE:HG23	2.16	0.45
1:B:162:PRO:HB2	1:B:276:ILE:HD12	1.97	0.45
1:B:345:ILE:O	1:B:517:ARG:NH1	2.44	0.45
1:B:350:PRO:CG	1:B:353:LEU:HD21	2.47	0.44
1:B:153:LEU:HB2	1:B:250:VAL:HB	2.00	0.44
1:B:200:LYS:HE2	6:B:135:HOH:O	2.18	0.43
1:B:445:ARG:HD2	1:B:496:GLU:HG2	2.02	0.42
1:B:427:VAL:HB	1:B:428:PRO:HD3	2.03	0.41
1:B:193:ILE:HD12	1:B:242:LEU:HD21	2.02	0.41
1:B:473:VAL:CG1	1:B:477:VAL:CG2	3.00	0.40
1:B:198:ILE:HB	1:B:199:PRO:HD3	2.02	0.40
1:B:229:GLU:HA	1:B:232:LYS:HE2	2.03	0.40
1:B:423:THR:HB	1:B:481:TYR:CE1	2.56	0.40

All (2) 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)
6:B:699:HOH:O	6:B:707:HOH:O[2_645]	1.97	0.23
6:B:582:HOH:O	6:B:650:HOH:O[1_655]	2.05	0.15

## 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	B	378/381 (99%)	366 (97%)	10 (3%)	2 (0%)	25	14

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	348	ALA
1	B	349	GLY

### 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	B	314/312 (101%)	312 (99%)	2 (1%)	84	82

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	406	SER
1	B	417	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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	$\# Z  > 2$
1	CSO	B	450	1	3,6,7	0.77	0	1,6,8	1.12	0

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
1	CSO	B	450	1	-	0/1/5/7	-

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.

## 5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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	$\# Z  > 2$
4	GOL	B	529	-	5,5,5	0.36	0	5,5,5	0.19	0
4	GOL	B	530	-	5,5,5	0.33	0	5,5,5	0.43	0
4	GOL	B	532	-	5,5,5	0.35	0	5,5,5	0.34	0
5	PO4	X	523	2,3	0,3,4	-	-	0,3,6	-	-
4	GOL	B	531	-	5,5,5	0.38	0	5,5,5	0.32	0
5	PO4	B	533	-	4,4,4	0.96	0	6,6,6	0.46	0

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
4	GOL	B	531	-	-	3/4/4/4	-
4	GOL	B	529	-	-	0/4/4/4	-
4	GOL	B	530	-	-	2/4/4/4	-
4	GOL	B	532	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	530	GOL	C1-C2-C3-O3
4	B	532	GOL	C1-C2-C3-O3
4	B	532	GOL	O2-C2-C3-O3
4	B	531	GOL	O1-C1-C2-C3
4	B	532	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
4	B	530	GOL	O2-C2-C3-O3
4	B	531	GOL	O1-C1-C2-O2
4	B	532	GOL	O1-C1-C2-O2
4	B	531	GOL	C1-C2-C3-O3

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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	B	373/381 (97%)	-0.33	13 (3%)	47	45	0, 9, 33, 51	3 (0%)
2	X	2/5 (40%)	0.43	0	100	100	9, 9, 9, 19	2 (100%)
All	All	375/386 (97%)	-0.33	13 (3%)	47	45	0, 9, 33, 51	5 (1%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	399	GLY	5.0
1	B	401	TRP	4.2
1	B	505	GLN	3.0
1	B	508	LEU	2.9
1	B	402	GLN	2.5
1	B	300	ARG	2.4
1	B	299	GLY	2.3
1	B	184	PHE	2.2
1	B	345	ILE	2.2
1	B	301	LYS	2.1
1	B	405	VAL	2.0
1	B	509[A]	ASP	2.0
1	B	348	ALA	2.0

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

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, 95<sup>th</sup> 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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
1	CSO	B	450	7/8	0.98	0.06	3,8,19,22	0

### 6.3 Carbohydrates [i](#)

There are no monosaccharides 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, 95<sup>th</sup> 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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	GOL	B	532	6/6	0.77	0.18	33,40,46,49	0
4	GOL	B	531	6/6	0.88	0.15	17,32,40,47	0
3	MG	B	1	1/1	0.91	0.19	38,38,38,38	0
4	GOL	B	530	6/6	0.94	0.08	7,15,28,34	0
5	PO4	B	533	5/5	0.96	0.09	13,14,16,18	0
4	GOL	B	529	6/6	0.97	0.06	0,3,20,24	0
5	PO4	X	523	4/5	0.99	0.03	0,1,1,8	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.