



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 03:30 am GMT

PDB ID : 3K59  
Title : Crystal structure of E.coli Pol II-normal DNA-dCTP ternary complex  
Authors : Yang, W.; Wang, F.  
Deposited on : 2009-10-06  
Resolution : 1.92 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.

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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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

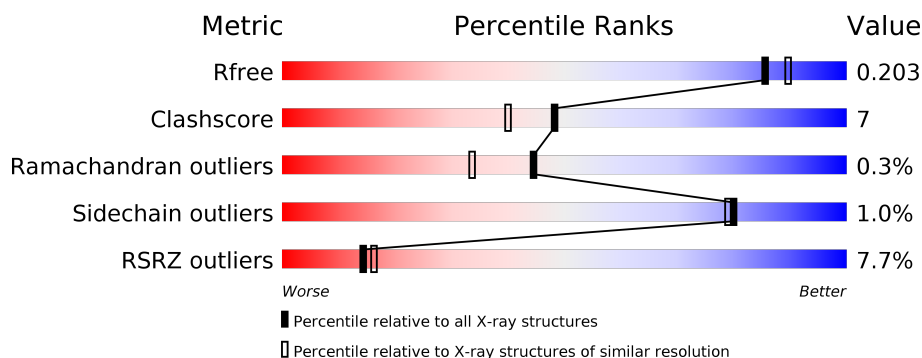
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.92 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))
$R_{free}$	100719	6276 (1.94-1.90)
Clashscore	112137	7025 (1.94-1.90)
Ramachandran outliers	110173	6947 (1.94-1.90)
Sidechain outliers	110143	6948 (1.94-1.90)
RSRZ outliers	101464	6332 (1.94-1.90)

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.

Mol	Chain	Length	Quality of chain
1	A	786	<div> <div>7%</div> <div> <div></div> <div>85%</div> <div>14%</div> <div></div> </div> </div>
2	T	17	<div> <div>29%</div> <div> <div></div> <div>82%</div> <div>18%</div> </div> </div>
3	P	13	<div> <div>15%</div> <div> <div></div> <div>69%</div> <div>31%</div> </div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7507 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 II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	781	Total	C	N	O	S	0	0	0
			6342	4039	1130	1148	25			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P21189
A	-1	PRO	-	EXPRESSION TAG	UNP P21189
A	0	HIS	-	EXPRESSION TAG	UNP P21189
A	335	ASN	ASP	ENGINEERED	UNP P21189

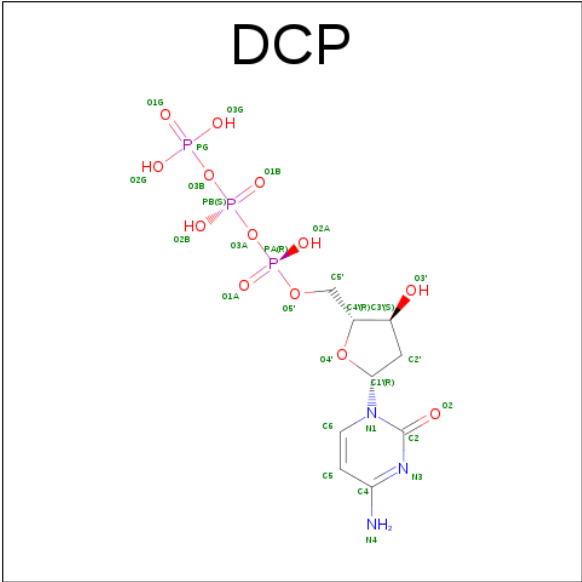
- Molecule 2 is a DNA chain called DNA (5'-D(\*TP\*AP\*GP\*GP\*TP\*AP\*CP\*GP\*CP\*TP\*AP\*GP\*GP\*CP\*AP\*CP\*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	17	Total	C	N	O	P	0	0	0
			348	166	68	98	16			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	13	Total	C	N	O	P	0	0	0
			262	126	48	76	12			

- Molecule 4 is 2'-DEOXYCYTIDINE-5'-TRIPHOSPHATE (three-letter code: DCP) (formula: C<sub>9</sub>H<sub>16</sub>N<sub>3</sub>O<sub>13</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			28	9	3	13	3		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Mg	0	0
			2	2		

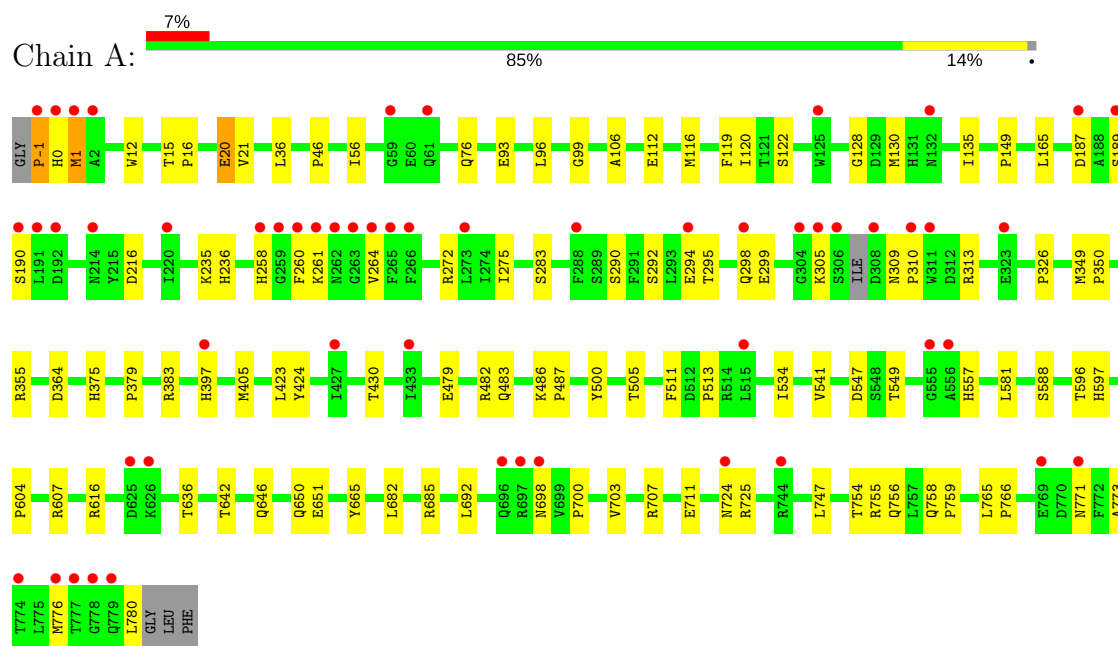
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	470	Total	O	0	0
			470	470		
6	T	35	Total	O	0	0
			35	35		
6	P	20	Total	O	0	0
			20	20		

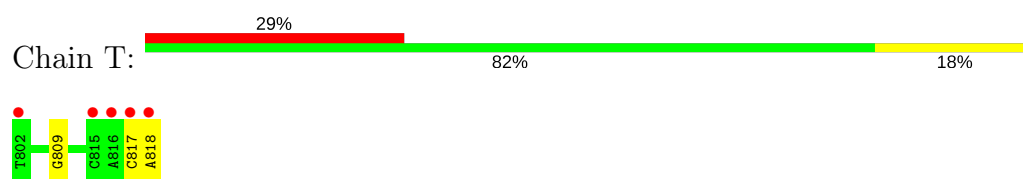
### 3 Residue-property plots [i](#)

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.

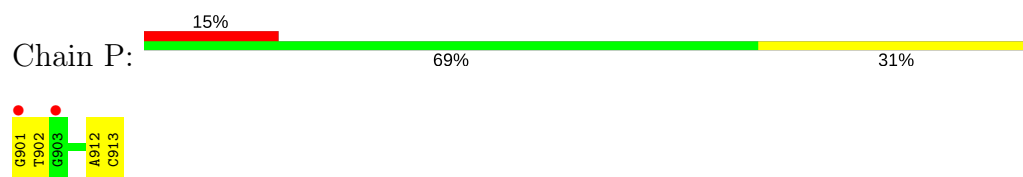
#### • Molecule 1: DNA polymerase II



#### • Molecule 2: DNA (5'-D(\*TP\*AP\*GP\*GP\*TP\*AP\*CP\*GP\*CP\*TP\*AP\*GP\*GP\*CP\*AP\*CP\*A)-3')



#### • Molecule 3: DNA (5'-D(\*GP\*TP\*GP\*CP\*CP\*TP\*AP\*GP\*CP\*GP\*TP\*AP\*(DOC))-3')



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	80.73Å 100.66Å 126.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.92 29.35 – 1.92	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-1.92) 93.4 (29.35-1.92)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.62 (at 1.92Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.198 , 0.202 0.197 , 0.203	Depositor DCC
$R_{free}$ test set	1809 reflections (2.45%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.9	Xtriage
Anisotropy	0.402	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 49.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7507	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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: DOC, MG, DCP

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	0.38	0/6511	0.69	1/8830 (0.0%)
2	T	0.59	0/391	1.00	0/602
3	P	0.45	0/273	0.81	0/420
All	All	0.40	0/7175	0.72	1/9852 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	PRO	N-CA-C	-5.55	97.66	112.10

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	6342	0	6215	86	0
2	T	348	0	192	6	0
3	P	262	0	148	3	0
4	A	28	0	9	2	0
5	A	2	0	0	0	0
6	A	470	0	0	3	0
6	P	20	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	T	35	0	0	0	0
All	All	7507	0	6564	94	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 7.

All (94) 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:1:MET:HE2	1:A:135:ILE:HD12	1.37	1.06
1:A:1:MET:SD	1:A:1:MET:C	2.55	0.85
1:A:379:PRO:O	1:A:383:ARG:HG2	1.79	0.81
2:T:817:DC:H2''	2:T:818:DA:C8	2.18	0.78
1:A:298:GLN:OE1	1:A:305:LYS:HG2	1.84	0.77
1:A:754:THR:HG22	1:A:755:ARG:HD3	1.67	0.77
1:A:1:MET:CE	1:A:135:ILE:HD12	2.15	0.77
1:A:581:LEU:HD12	1:A:588:SER:HB2	1.68	0.76
1:A:258:HIS:HB2	1:A:264:VAL:CG2	2.16	0.75
1:A:1:MET:O	1:A:1:MET:SD	2.44	0.75
1:A:258:HIS:CE1	1:A:261:LYS:HG3	2.23	0.72
1:A:1:MET:HE2	1:A:135:ILE:CD1	2.22	0.68
3:P:912:DA:H2''	3:P:913:DOC:H5'	1.77	0.66
1:A:120:ILE:O	1:A:375:HIS:HE1	1.81	0.64
1:A:1:MET:HE1	1:A:135:ILE:HG13	1.81	0.62
1:A:292:SER:OG	1:A:294:GLU:HG2	2.01	0.61
1:A:190:SER:HB2	6:A:1203:HOH:O	2.01	0.59
1:A:1:MET:CE	1:A:135:ILE:CD1	2.79	0.59
1:A:479:GLU:OE2	1:A:483:GLN:NE2	2.35	0.59
1:A:692:LEU:HD21	1:A:707:ARG:HG2	1.85	0.59
1:A:258:HIS:HB2	1:A:264:VAL:HG23	1.83	0.58
1:A:298:GLN:OE1	1:A:305:LYS:CG	2.50	0.58
1:A:692:LEU:HD21	1:A:707:ARG:HE	1.68	0.58
1:A:479:GLU:OE2	1:A:482:ARG:NH1	2.36	0.58
1:A:534:ILE:CD1	1:A:549:THR:HG21	2.34	0.57
1:A:112:GLU:O	1:A:116:MET:HB2	2.05	0.57
1:A:294:GLU:CG	1:A:295:THR:N	2.67	0.56
1:A:776:MET:O	1:A:780:LEU:N	2.29	0.55
1:A:557:HIS:HA	6:A:844:HOH:O	2.05	0.55
1:A:283:SER:HB2	1:A:355:ARG:HH12	1.71	0.55
4:A:914:DCP:H6	3:P:913:DOC:H2'	1.88	0.54
1:A:698:ASN:C	1:A:698:ASN:OD1	2.45	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:GLU:HG3	1:A:295:THR:N	2.22	0.53
1:A:758:GLN:HB3	1:A:759:PRO:HD3	1.91	0.52
1:A:295:THR:O	1:A:299:GLU:HG3	2.10	0.52
1:A:754:THR:CG2	1:A:755:ARG:HH11	2.22	0.52
1:A:724:ASN:O	1:A:725:ARG:HB2	2.11	0.51
3:P:901:DG:H2'	3:P:902:DT:H72	1.94	0.51
2:T:817:DC:H2''	2:T:818:DA:H8	1.72	0.50
1:A:216:ASP:HA	1:A:272:ARG:NH2	2.26	0.50
1:A:310:PRO:HA	1:A:313:ARG:HB3	1.93	0.50
1:A:486:LYS:HB2	1:A:487:PRO:HD3	1.94	0.49
1:A:423:LEU:HD22	1:A:547:ASP:HB3	1.94	0.49
1:A:642:THR:O	1:A:646:GLN:HG3	2.13	0.48
1:A:258:HIS:NE2	1:A:261:LYS:HE2	2.28	0.48
1:A:235:LYS:HE2	1:A:235:LYS:HB3	1.68	0.48
1:A:216:ASP:HA	1:A:272:ARG:HH21	1.78	0.48
1:A:21:VAL:HB	1:A:36:LEU:HD12	1.96	0.48
1:A:692:LEU:CD2	1:A:707:ARG:HE	2.27	0.48
1:A:430:THR:HG21	1:A:581:LEU:HD13	1.96	0.47
1:A:424:TYR:CD2	4:A:914:DCP:H2'2	2.50	0.47
1:A:581:LEU:HD12	1:A:588:SER:CB	2.42	0.47
1:A:604:PRO:HB3	6:A:964:HOH:O	2.15	0.47
1:A:258:HIS:CB	1:A:264:VAL:CG2	2.92	0.46
1:A:119:PHE:HE2	1:A:383:ARG:NH2	2.13	0.46
2:T:818:DA:H8	2:T:818:DA:O5'	1.98	0.46
1:A:754:THR:HG21	1:A:755:ARG:HH11	1.80	0.46
1:A:616:ARG:HD3	2:T:809:DG:H4'	1.98	0.46
2:T:817:DC:H2''	2:T:818:DA:O5'	2.16	0.46
2:T:817:DC:C2'	2:T:818:DA:C8	2.97	0.46
1:A:165:LEU:HG	1:A:236:HIS:NE2	2.31	0.45
1:A:56:ILE:HD12	1:A:96:LEU:HA	1.97	0.45
1:A:776:MET:O	1:A:780:LEU:HB2	2.16	0.45
1:A:119:PHE:CE2	1:A:383:ARG:NH2	2.84	0.45
1:A:682:LEU:HD12	1:A:747:LEU:HD13	1.99	0.44
1:A:260:PHE:CE1	1:A:261:LYS:CG	3.01	0.44
1:A:636:THR:HG21	1:A:650:GLN:CG	2.47	0.44
1:A:56:ILE:HD11	1:A:99:GLY:HA3	1.99	0.44
1:A:596:THR:HG22	1:A:597:HIS:N	2.33	0.43
1:A:534:ILE:HD13	1:A:549:THR:HG21	2.01	0.43
1:A:707:ARG:O	1:A:711:GLU:HG3	2.18	0.43
1:A:-1:PRO:HA	1:A:130:MET:HB3	2.00	0.43
1:A:423:LEU:HD22	1:A:547:ASP:CB	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:MET:HE1	1:A:541:VAL:HG12	2.01	0.43
1:A:700:PRO:HD2	1:A:703:VAL:HB	2.00	0.43
1:A:0:HIS:HA	1:A:128:GLY:O	2.18	0.42
1:A:216:ASP:OD1	1:A:272:ARG:NH2	2.53	0.42
1:A:642:THR:H	1:A:756:GLN:NE2	2.17	0.42
1:A:765:LEU:N	1:A:766:PRO:CD	2.83	0.42
1:A:283:SER:CB	1:A:355:ARG:HH12	2.32	0.42
1:A:93:GLU:HG3	1:A:106:ALA:HB3	2.02	0.42
1:A:46:PRO:HB3	1:A:76:GLN:HG2	2.01	0.41
1:A:275:ILE:HD13	1:A:364:ASP:HB3	2.02	0.41
1:A:505:THR:O	1:A:511:PHE:HB2	2.20	0.41
1:A:651:GLU:OE1	1:A:665:TYR:OH	2.29	0.41
1:A:685:ARG:HB2	1:A:685:ARG:HE	1.30	0.41
1:A:771:ASN:OD1	1:A:773:ALA:HB3	2.20	0.41
1:A:534:ILE:HD12	1:A:549:THR:HG21	2.00	0.41
1:A:12:TRP:HA	1:A:20:GLU:O	2.20	0.41
1:A:397:HIS:HB3	1:A:513:PRO:HG2	2.03	0.41
1:A:187:ASP:CG	1:A:189:SER:HG	2.25	0.41
1:A:15:THR:HB	1:A:16:PRO:CD	2.50	0.40
1:A:260:PHE:CE1	1:A:261:LYS:HG3	2.56	0.40
1:A:349:MET:HB2	1:A:350:PRO:HD3	2.02	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	777/786 (99%)	761 (98%)	14 (2%)	2 (0%)	44 33

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	309	ASN
1	A	122	SER

### 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	668/672 (99%)	661 (99%)	7 (1%)	80	79

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-1	PRO
1	A	1	MET
1	A	20	GLU
1	A	290	SER
1	A	326	PRO
1	A	500	TYR
1	A	607	ARG

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	137	ASN
1	A	332	ASN
1	A	375	HIS
1	A	449	HIS
1	A	464	HIS
1	A	557	HIS
1	A	735	ASN
1	A	756	GLN

### 5.3.3 RNA [i](#)

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$
3	DOC	P	913	3,2	13,19,20	0.97	1 (7%)	12,26,29	1.38	1 (8%)

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
3	DOC	P	913	3,2	-	0/3/18/19	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	913	DOC	O5'-C5'	-2.01	1.41	1.44

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	913	DOC	C3'-C2'-C1'	3.43	106.50	102.69

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	P	913	DOC	2	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	DCP	A	914	5	23,29,29	2.80	2 (8%)	25,45,45	2.13	5 (20%)

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	DCP	A	914	5	-	0/18/34/34	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	914	DCP	C5'-C4'	-12.18	1.13	1.51
4	A	914	DCP	C6-N1	3.94	1.41	1.35

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	914	DCP	C2'-C1'-N1	-3.05	107.03	114.23
4	A	914	DCP	O2A-PA-O5'	-2.93	94.31	108.14
4	A	914	DCP	O5'-PA-O1A	3.09	121.72	109.25
4	A	914	DCP	O5'-C5'-C4'	4.80	126.03	109.00
4	A	914	DCP	O4'-C1'-N1	6.48	118.70	107.78

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	914	DCP	2	0

## 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	A	781/786 (99%)	0.47	55 (7%) 17 19	15, 26, 50, 67	0
2	T	17/17 (100%)	0.90	5 (29%) 1 0	20, 32, 71, 76	0
3	P	12/13 (92%)	0.68	2 (16%) 2 2	20, 45, 65, 66	0
All	All	810/816 (99%)	0.48	62 (7%) 14 16	15, 26, 54, 76	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	311	TRP	9.1
1	A	306	SER	8.2
1	A	260	PHE	7.5
1	A	305	LYS	6.6
1	A	308	ASP	6.1
1	A	264	VAL	5.7
1	A	259	GLY	5.3
1	A	310	PRO	5.2
1	A	262	ASN	5.1
1	A	304	GLY	4.9
1	A	-1	PRO	4.8
1	A	189	SER	4.4
2	T	818	DA	4.3
1	A	0	HIS	4.3
1	A	261	LYS	4.2
1	A	190	SER	4.2
1	A	1	MET	4.1
1	A	777	THR	4.1
1	A	263	GLY	3.9
1	A	778	GLY	3.9
1	A	187	ASP	3.7
1	A	776	MET	3.7
1	A	696	GLN	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	258	HIS	3.6
1	A	61	GLN	3.6
1	A	697	ARG	3.6
1	A	59	GLY	3.4
1	A	625	ASP	3.4
1	A	433	ILE	3.3
2	T	802	DT	3.1
1	A	132	ASN	3.1
1	A	724	ASN	3.0
3	P	901	DG	3.0
1	A	771	ASN	2.9
1	A	555	GLY	2.9
1	A	397	HIS	2.9
1	A	298	GLN	2.9
1	A	427	ILE	2.8
2	T	817	DC	2.8
1	A	556	ALA	2.8
3	P	903	DG	2.7
1	A	698	ASN	2.7
1	A	265	PHE	2.6
1	A	288	PHE	2.6
2	T	815	DC	2.6
2	T	816	DA	2.5
1	A	626	LYS	2.5
1	A	774	THR	2.5
1	A	273	LEU	2.4
1	A	515	LEU	2.4
1	A	191	LEU	2.3
1	A	294	GLU	2.3
1	A	214	ASN	2.2
1	A	220	ILE	2.2
1	A	744	ARG	2.2
1	A	2	ALA	2.2
1	A	323	GLU	2.2
1	A	779	GLN	2.2
1	A	192	ASP	2.1
1	A	769	GLU	2.1
1	A	125	TRP	2.1
1	A	266	PHE	2.1



## 6.2 Non-standard residues in protein, DNA, RNA chains [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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	DOC	P	913	18/19	0.98	0.10	-	17,18,21,22	0

## 6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
5	MG	A	1001	1/1	0.98	0.12	0.37	20,20,20,20	0
4	DCP	A	914	28/28	0.97	0.11	0.05	14,17,19,22	0
5	MG	A	1002	1/1	0.99	0.05	-	21,21,21,21	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.