



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:23 PM UTC

PDB ID : 4EBC / pdb\_00004ebc  
Title : Conformationally Restrained North-methanocarba-2'-deoxyadenosine Corrects the Error-Prone Nature of Human DNA Polymerase Iota  
Authors : Eoff, R.L.; Ketkar, A.; Banerjee, S.  
Deposited on : 2012-03-23  
Resolution : 2.90 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 : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtrriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

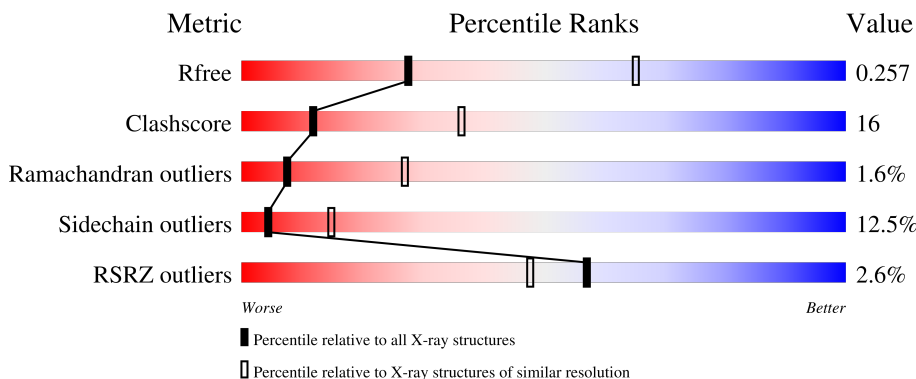
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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))
$R_{free}$	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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 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	A	420	 2% 60% 25% 10%
2	P	7	 43% 57%
3	T	9	 22% 78%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	GOL	A	505	-	-	X	-

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3365 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 iota.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	376	2922	1852	509	539	22	3	0	0

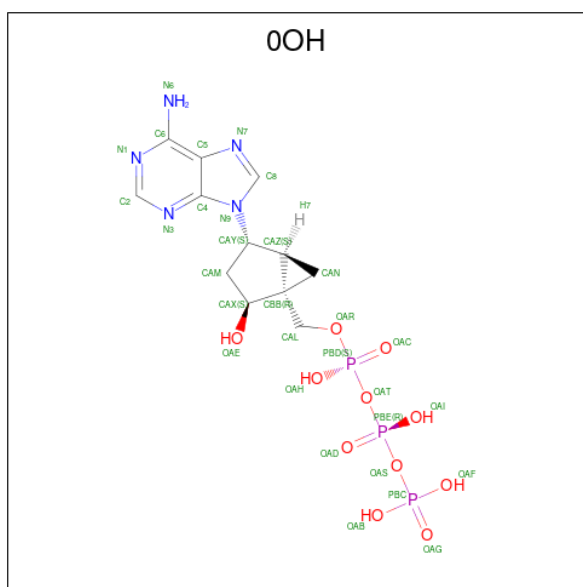
- Molecule 2 is a DNA chain called 5'-D(P\*AP\*GP\*GP\*AP\*CP\*CP\*(DOC))-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	P	7	142	67	29	39	7	0	0	0

- Molecule 3 is a DNA chain called 5'-D(P\*CP\*TP\*GP\*GP\*GP\*TP\*CP\*CP\*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	T	9	183	87	30	57	9	0	0	0

- Molecule 4 is North-methanocarpa-2'-deoxyadenosine triphosphate (CCD ID: 0OH) (formula: C<sub>12</sub>H<sub>18</sub>N<sub>5</sub>O<sub>11</sub>P<sub>3</sub>).

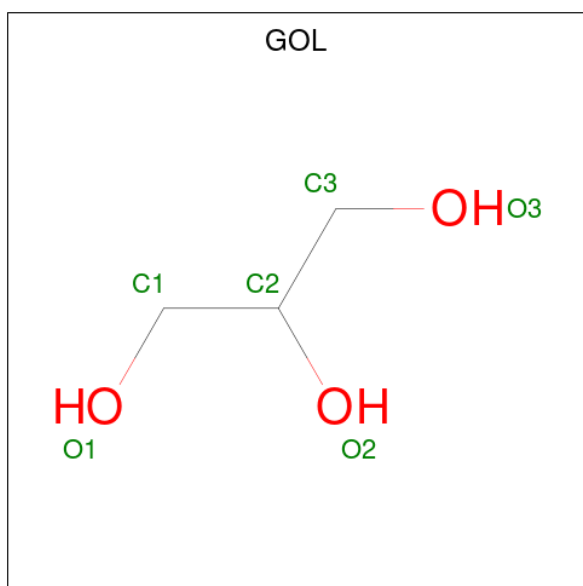


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	A	1	31	12	5	11	3	0	0

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
5	A	3	3	3	0	0

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	61	Total	O	0	0
			61	61		
7	P	8	Total	O	0	0
			8	8		
7	T	9	Total	O	0	0
			9	9		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	97.80Å 97.80Å 203.63Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.90 – 2.90 48.90 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.1 (48.90-2.90) 93.0 (48.90-2.90)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.00 (at 2.91Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, $R_{free}$	0.206 , 0.253 0.212 , 0.257	Depositor DCC
$R_{free}$ test set	1278 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	60.5	Xtrriage
Anisotropy	0.265	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 49.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3365	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	81.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.77% 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: CA, DOC, GOL, 0OH

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	0.60	0/2964	0.94	5/4003 (0.1%)
2	P	0.37	0/139	1.03	0/212
3	T	0.54	0/203	1.20	0/311
All	All	0.59	0/3306	0.97	5/4526 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	364	ILE	CB-CA-C	-7.21	102.58	110.95
1	A	298	PRO	O-C-N	5.94	124.04	121.31
1	A	149	SER	N-CA-C	-5.93	107.28	114.75
1	A	386	ILE	CB-CA-C	-5.63	104.77	111.97
1	A	151	VAL	N-CA-C	5.43	116.42	109.30

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	A	2922	0	2972	92	0
2	P	142	0	78	6	3
3	T	183	0	103	8	3

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	31	0	18	3	0
5	A	3	0	0	0	0
6	A	6	0	8	5	0
7	A	61	0	0	23	0
7	P	8	0	0	1	0
7	T	9	0	0	3	0
All	All	3365	0	3179	102	3

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 (102) 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)
6:A:505:GOL:H12	3:T:6:DG:OP2	1.42	1.20
1:A:335:ASP:C	1:A:337:ARG:N	2.08	1.11
1:A:177:SER:HB2	7:A:651:HOH:O	1.62	0.99
1:A:214:LYS:HD3	7:A:645:HOH:O	1.69	0.92
1:A:62:LEU:HG	7:T:107:HOH:O	1.71	0.90
1:A:302:PHE:HZ	1:A:332:VAL:HG23	1.46	0.80
1:A:44:MET:HE2	1:A:51:LYS:HA	1.64	0.78
3:T:7:DG:H1'	7:T:109:HOH:O	1.83	0.78
1:A:335:ASP:CB	1:A:337:ARG:N	2.50	0.75
1:A:44:MET:HE3	1:A:67:ASN:HD22	1.52	0.74
1:A:103:ARG:NH1	7:A:613:HOH:O	2.21	0.73
1:A:218:GLN:HG3	7:A:637:HOH:O	1.88	0.72
1:A:214:LYS:CD	7:A:645:HOH:O	2.34	0.71
6:A:505:GOL:C1	3:T:6:DG:OP2	2.30	0.71
1:A:72:LYS:HG3	1:A:73:LEU:HD23	1.72	0.70
1:A:144:GLN:O	1:A:146:ASP:N	2.25	0.69
1:A:136:VAL:HG22	1:A:175:VAL:HG12	1.77	0.67
1:A:119:VAL:HB	7:A:650:HOH:O	1.93	0.67
1:A:292:PRO:HA	7:A:659:HOH:O	1.94	0.67
1:A:302:PHE:CZ	1:A:332:VAL:HG23	2.30	0.66
1:A:67:ASN:HB2	7:A:622:HOH:O	1.95	0.66
1:A:343:ARG:HB3	1:A:409:CYS:HB3	1.78	0.65
1:A:101:ARG:N	7:A:623:HOH:O	2.28	0.65
1:A:259:ARG:HA	1:A:262:GLN:HB2	1.79	0.64
1:A:335:ASP:HB2	1:A:337:ARG:N	2.12	0.62
1:A:177:SER:CB	7:A:651:HOH:O	2.32	0.62
1:A:325:LEU:HD22	1:A:380:MET:HE3	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:GLU:HG3	7:A:634:HOH:O	1.99	0.61
1:A:44:MET:CE	1:A:67:ASN:HD22	2.12	0.61
1:A:304:GLU:OE2	6:A:505:GOL:C1	2.49	0.61
1:A:167:ASP:HB3	1:A:170:HIS:HB2	1.83	0.59
1:A:302:PHE:CD2	7:A:648:HOH:O	2.51	0.59
6:A:505:GOL:H12	3:T:6:DG:P	2.43	0.58
1:A:335:ASP:CA	1:A:337:ARG:N	2.66	0.58
1:A:69:GLU:HG3	7:A:625:HOH:O	2.04	0.57
1:A:357:ARG:NH1	7:A:646:HOH:O	2.37	0.56
1:A:361:GLN:N	2:P:1:DA:OP1	2.38	0.56
1:A:204:LEU:HD22	1:A:284:LEU:HD23	1.89	0.53
1:A:103:ARG:NH2	1:A:331:ARG:HD3	2.24	0.53
1:A:267:LYS:H	1:A:267:LYS:HD2	1.75	0.52
1:A:348:ARG:NH1	1:A:358:GLU:OE1	2.43	0.52
1:A:39:TYR:CE2	4:A:501:OOH:H12	2.46	0.50
1:A:303:SER:O	1:A:331:ARG:NH2	2.43	0.50
1:A:41:GLN:O	1:A:45:ILE:HG13	2.12	0.50
1:A:359:SER:O	1:A:360:ARG:HD2	2.12	0.49
1:A:335:ASP:CG	1:A:337:ARG:N	2.70	0.49
1:A:277:VAL:O	1:A:281:ILE:HG12	2.13	0.49
1:A:201:SER:HB3	7:A:649:HOH:O	2.12	0.48
1:A:365:PRO:HB2	1:A:368:VAL:HG23	1.95	0.48
1:A:39:TYR:CD2	4:A:501:OOH:H12	2.49	0.47
1:A:72:LYS:HD2	7:A:625:HOH:O	2.14	0.47
1:A:329:LEU:O	1:A:332:VAL:HG12	2.13	0.47
1:A:335:ASP:N	1:A:335:ASP:OD1	2.48	0.47
1:A:61:TYR:HE2	7:A:655:HOH:O	1.97	0.47
1:A:331:ARG:NH1	7:A:628:HOH:O	2.47	0.47
1:A:413:LEU:HD22	1:A:413:LEU:H	1.79	0.47
3:T:9:DC:H2''	3:T:10:DC:O5'	2.14	0.46
1:A:220:VAL:HB	7:A:651:HOH:O	2.14	0.46
1:A:120:VAL:HG22	1:A:130:VAL:HG22	1.98	0.46
1:A:214:LYS:NZ	7:A:644:HOH:O	2.49	0.46
1:A:365:PRO:O	1:A:368:VAL:HB	2.16	0.46
1:A:159:ASN:N	7:A:632:HOH:O	2.31	0.45
1:A:36:ASP:O	1:A:37:CYS:C	2.58	0.45
1:A:313:SER:OG	1:A:316:GLU:HB2	2.17	0.45
1:A:239:ILE:HB	1:A:242:ILE:HD12	1.98	0.44
1:A:184:ARG:HA	1:A:187:MET:HE3	1.99	0.44
1:A:273:LEU:HB3	1:A:277:VAL:HG22	1.99	0.44
1:A:387:LEU:HD23	1:A:387:LEU:HA	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:MET:O	1:A:139:ARG:HG3	2.18	0.44
1:A:106:SER:OG	1:A:122:ARG:NH1	2.45	0.43
1:A:267:LYS:H	1:A:267:LYS:CD	2.31	0.43
1:A:305:GLU:HB2	3:T:5:DG:H2''	1.99	0.43
1:A:224:GLU:CD	1:A:224:GLU:H	2.27	0.43
1:A:273:LEU:O	1:A:277:VAL:HG13	2.19	0.43
1:A:365:PRO:HB2	1:A:368:VAL:CG2	2.49	0.42
1:A:178:GLN:NE2	7:A:653:HOH:O	2.51	0.42
2:P:4:DA:H2'	2:P:5:DC:C6	2.54	0.42
1:A:305:GLU:HB2	3:T:5:DG:C2'	2.49	0.42
1:A:88:CYS:SG	1:A:91:LEU:HB2	2.59	0.42
1:A:119:VAL:HG12	1:A:293:VAL:HG12	2.02	0.42
1:A:135:MET:HE2	1:A:135:MET:HB3	1.76	0.42
1:A:144:GLN:HG2	1:A:147:GLU:OE1	2.19	0.42
3:T:4:DT:H2'	3:T:5:DG:C8	2.54	0.42
1:A:304:GLU:OE2	6:A:505:GOL:H12	2.19	0.42
2:P:4:DA:H2''	2:P:5:DC:O4'	2.20	0.41
1:A:60:LYS:HE3	1:A:97:GLU:OE2	2.20	0.41
1:A:364:ILE:H	1:A:364:ILE:HG12	1.71	0.41
1:A:388:MET:HA	1:A:388:MET:HE2	2.02	0.41
1:A:100:THR:O	1:A:104:GLU:HG3	2.19	0.41
1:A:140:LEU:HD23	1:A:140:LEU:HA	1.84	0.41
1:A:334:GLN:HB2	1:A:335:ASP:H	1.78	0.41
1:A:39:TYR:HB2	1:A:65:THR:CG2	2.50	0.41
1:A:207:LYS:NZ	2:P:7:DOC:OP1	2.33	0.41
1:A:265:SER:HA	1:A:266:PRO:HD3	1.89	0.41
1:A:283:LYS:NZ	1:A:288:GLU:OE1	2.44	0.41
2:P:7:DOC:OP2	7:P:102:HOH:O	2.21	0.41
1:A:62:LEU:CD2	7:T:107:HOH:O	2.69	0.40
4:A:501:OOH:H5	2:P:7:DOC:H2''	2.03	0.40
1:A:39:TYR:HB2	1:A:65:THR:HG21	2.01	0.40
1:A:236:ILE:H	1:A:236:ILE:HG12	1.51	0.40
1:A:201:SER:CB	7:A:649:HOH:O	2.68	0.40
1:A:405:LEU:HD23	1:A:406:LEU:N	2.37	0.40

All (3) 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)
2:P:1:DA:P	3:T:11:DT:O3'[10_557]	1.20	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:1:DA:OP2	3:T:11:DT:O3'[10_557]	2.04	0.16
2:P:1:DA:OP1	3:T:11:DT:O3'[10_557]	2.16	0.04

### 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	368/420 (88%)	332 (90%)	30 (8%)	6 (2%)	<b>7</b> <b>27</b>

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	145	SER
1	A	333	CYS
1	A	37	CYS
1	A	253	LEU
1	A	334	GLN
1	A	369	ILE

#### 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	329/376 (88%)	288 (88%)	41 (12%)	<b>4</b> <b>15</b>

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	73	LEU
1	A	80	ASN
1	A	114	GLU
1	A	132	LEU
1	A	142	GLN
1	A	143	LEU
1	A	148	LEU
1	A	149	SER
1	A	151	VAL
1	A	152	THR
1	A	154	SER
1	A	175	VAL
1	A	185	GLU
1	A	209	VAL
1	A	217	GLN
1	A	230	ILE
1	A	236	ILE
1	A	253	LEU
1	A	261	LEU
1	A	267	LYS
1	A	275	ILE
1	A	277	VAL
1	A	301	SER
1	A	308	PHE
1	A	329	LEU
1	A	332	VAL
1	A	333	CYS
1	A	334	GLN
1	A	335	ASP
1	A	342	VAL
1	A	345	ILE
1	A	364	ILE
1	A	366	SER
1	A	379	VAL
1	A	380	MET
1	A	388	MET
1	A	395	VAL
1	A	411	CYS
1	A	413	LEU
1	A	414	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	ASN
1	A	128	ASN
1	A	217	GLN
1	A	396	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](#)

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
2	DOC	P	7	3,2	16,19,20	0.45	0	20,26,29	0.80	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
2	DOC	P	7	3,2	-	0/7/18/19	0/2/2/2

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.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	P	7	DOC	3	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 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$
6	GOL	A	505	-	5,5,5	0.25	0	5,5,5	0.31	0
4	OOH	A	501	5	32,34,34	1.51	4 (12%)	43,56,56	1.96	14 (32%)

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
6	GOL	A	505	-	-	4/4/4/4	-
4	OOH	A	501	5	-	10/23/45/45	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	501	OOH	C6-N6	4.09	1.44	1.34
4	A	501	OOH	OAE-CAX	-4.07	1.36	1.43
4	A	501	OOH	CAM-CAY	-2.90	1.48	1.53
4	A	501	OOH	CAM-CAX	-2.10	1.48	1.53

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	501	0OH	N1-C2-N3	-5.54	120.20	128.58
4	A	501	0OH	C5-C4-N3	-4.05	121.13	126.72
4	A	501	0OH	C4-C5-N7	-4.02	105.99	110.58
4	A	501	0OH	CBB-CAZ-CAY	-3.49	105.67	108.39
4	A	501	0OH	C5-N7-C8	3.48	108.92	103.45
4	A	501	0OH	C2-N3-C4	3.48	120.33	111.83
4	A	501	0OH	N9-C8-N7	-3.07	109.58	113.94
4	A	501	0OH	C6-C5-N7	2.71	137.31	132.09
4	A	501	0OH	CAX-CAM-CAY	2.52	108.07	104.28
4	A	501	0OH	C2-N1-C6	2.41	122.70	118.73
4	A	501	0OH	OAI-PBE-OAT	2.36	113.64	107.27
4	A	501	0OH	C5-C4-N9	2.22	108.23	105.81
4	A	501	0OH	OAR-CAL-CBB	2.06	111.52	108.96
4	A	501	0OH	N3-C4-N9	2.04	130.63	127.17

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	501	0OH	CAL-OAR-PBD-OAT
4	A	501	0OH	CAL-OAR-PBD-OAC
4	A	501	0OH	OAR-CAL-CBB-CAX
6	A	505	GOL	O1-C1-C2-O2
6	A	505	GOL	O1-C1-C2-C3
6	A	505	GOL	C1-C2-C3-O3
4	A	501	0OH	PBE-OAT-PBD-OAR
6	A	505	GOL	O2-C2-C3-O3
4	A	501	0OH	CAL-OAR-PBD-OAH
4	A	501	0OH	OAR-CAL-CBB-CAZ
4	A	501	0OH	PBC-OAS-PBE-OAI
4	A	501	0OH	PBE-OAT-PBD-OAH
4	A	501	0OH	PBE-OAT-PBD-OAC
4	A	501	0OH	CAM-CAY-N9-C4

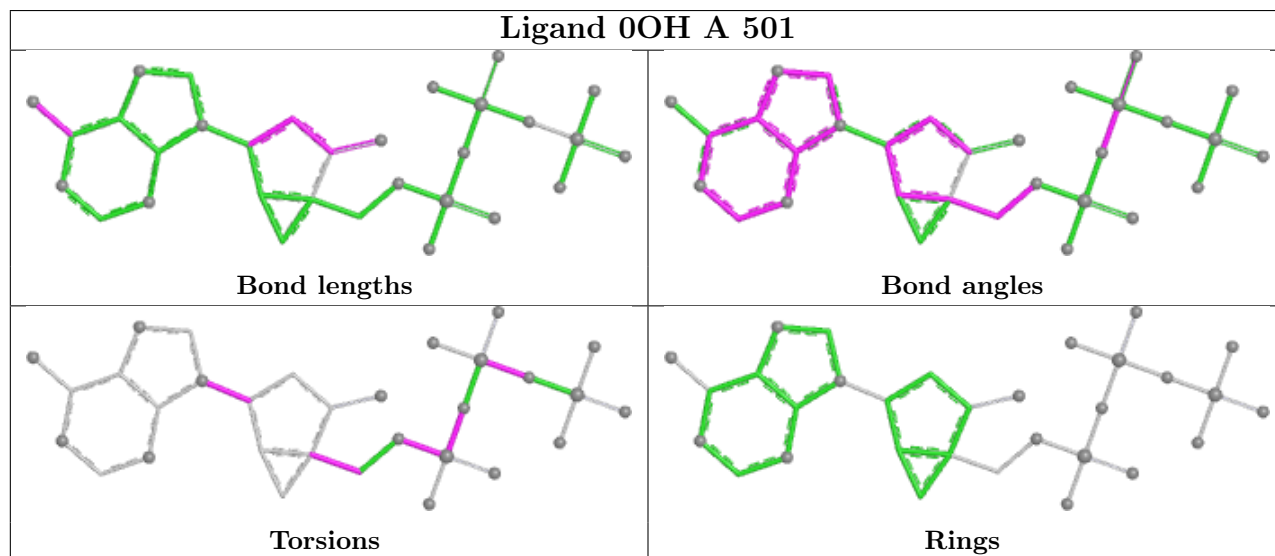
There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	505	GOL	5	0
4	A	501	0OH	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



## 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](#)

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	376/420 (89%)	-0.11	10 (2%) 56 47	35, 77, 136, 180	9 (2%)
2	P	6/7 (85%)	-0.33	0 100 100	63, 91, 96, 99	0
3	T	9/9 (100%)	-0.43	0 100 100	61, 66, 105, 163	0
All	All	391/436 (89%)	-0.12	10 (2%) 57 48	35, 77, 136, 180	9 (2%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	408	VAL	3.4
1	A	343	ARG	2.9
1	A	78	LEU	2.7
1	A	407	SER	2.4
1	A	245	LYS	2.3
1	A	344	LEU	2.2
1	A	364	ILE	2.2
1	A	406	LEU	2.2
1	A	302	PHE	2.1
1	A	237	LYS	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. 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
2	DOC	P	7	18/19	0.96	0.07	56,61,63,65	0

### 6.3 Carbohydrates [i](#)

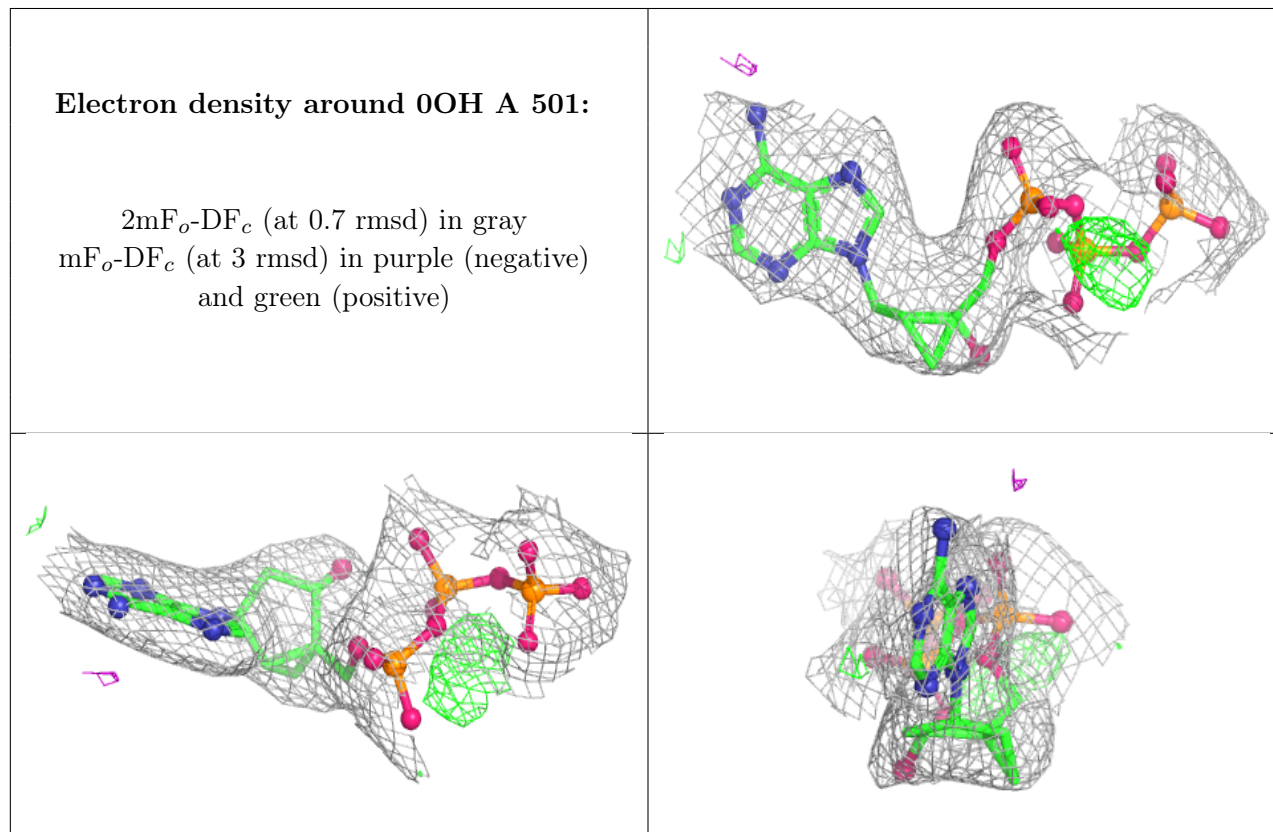
There are no oligosaccharides in this entry.

### 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
5	CA	A	504	1/1	0.72	0.13	103,103,103,103	0
6	GOL	A	505	6/6	0.75	0.15	59,66,71,72	0
5	CA	A	503	1/1	0.85	0.12	83,83,83,83	0
5	CA	A	502	1/1	0.97	0.07	49,49,49,49	0
4	OOH	A	501	31/31	0.97	0.07	48,58,67,70	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



## 6.5 Other polymers [i](#)

There are no such residues in this entry.