



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 08:36 am GMT

PDB ID : 4DL2
Title : Human DNA polymerase eta inserting dCMPNPP opposite CG template (GG0a)
Authors : Zhao, Y.; Biertumpfel, C.; Gregory, M.; Hua, Y.; Hanaoka, F.; Yang, W.
Deposited on : 2012-02-05
Resolution : 2.15 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

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 : 1.7.2 (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

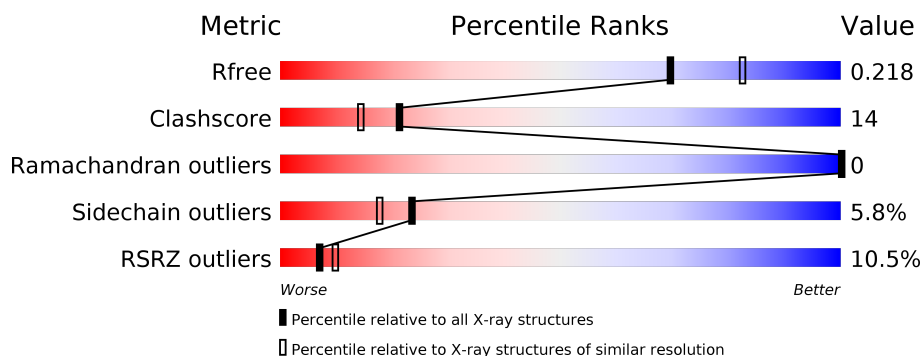
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	1170 (2.16-2.16)
Clashscore	112137	1278 (2.16-2.16)
Ramachandran outliers	110173	1256 (2.16-2.16)
Sidechain outliers	110143	1255 (2.16-2.16)
RSRZ outliers	101464	1175 (2.16-2.16)

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 ≥ 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	435	<div> <div>10%</div> <div>74%</div> <div>21%</div> <div>••</div> </div>
2	T	12	<div> <div>25%</div> <div>50%</div> <div>50%</div> </div>
3	P	9	<div> <div>44%</div> <div>56%</div> </div>

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	506	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4126 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 eta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	428	Total	C	N	O	S	0	8	0
			3380	2121	605	628	26			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP Q9Y253
A	-1	PRO	-	EXPRESSION TAG	UNP Q9Y253
A	0	HIS	-	EXPRESSION TAG	UNP Q9Y253

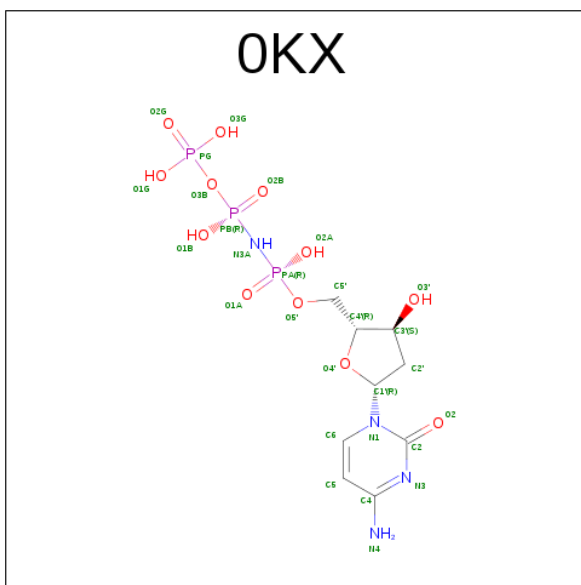
- Molecule 2 is a DNA chain called DNA (5'-D(*AP*CP*GP*GP*CP*TP*CP*AP*CP*AP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	12	Total	C	N	O	P	0	2	1
			266	125	49	79	13			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	9	Total	C	N	O	P	0	2	1
			214	100	44	60	10			

- Molecule 4 is 2'-DEOXY-5'-O-[(R)-HYDROXY{[(R)-HYDROXY(PHOSPHONOOXY)PHOSPHORYL]AMINO}PHOSPHORYL]CYTIDINE (three-letter code: 0KX) (formula: C₉H₁₇N₄O₁₂P₃).

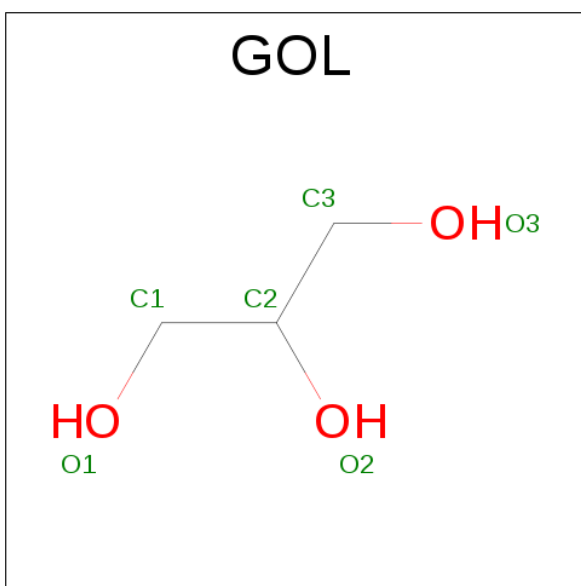


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	P	1	Total	C	N	O	P	0	0
			28	9	4	12	3		

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

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

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



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

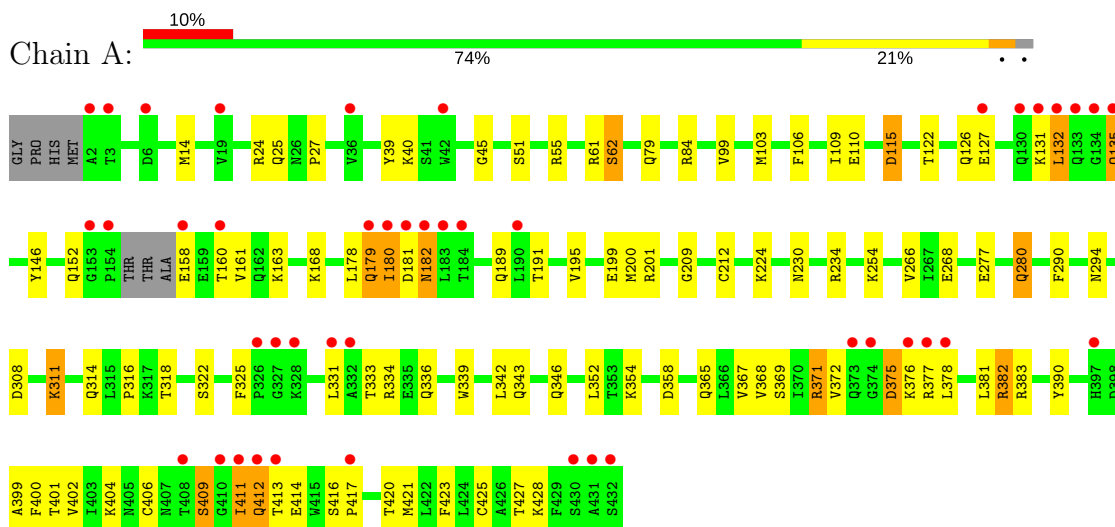
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	189	Total	O	0	0
			189	189		
7	P	15	Total	O	0	0
			15	15		
7	T	14	Total	O	0	0
			14	14		

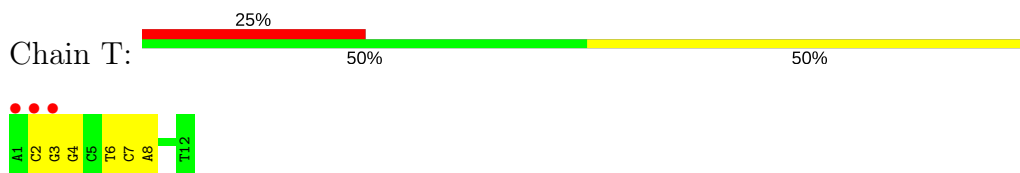
3 Residue-property plots

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.

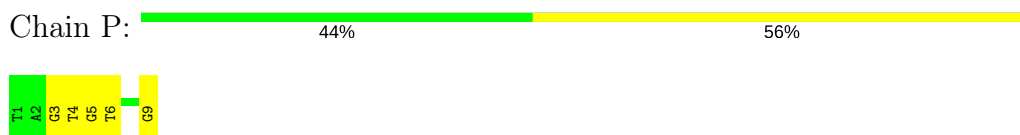
- Molecule 1: DNA polymerase eta



- Molecule 2: DNA (5'-D(*AP*CP*GP*GP*CP*TP*CP*AP*CP*AP*CP*T)-3')



- Molecule 3: DNA (5'-D(*TP*AP*GP*TP*GP*TP*GP*AP*G)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	97.86Å 97.86Å 80.53Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.15 19.44 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.6 (30.00-2.15) 99.5 (19.44-2.15)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 2.15Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.7.1_743)	Depositor
R, R_{free}	0.193 , 0.222 0.186 , 0.218	Depositor DCC
R_{free} test set	1831 reflections (7.69%)	DCC
Wilson B-factor (Å ²)	25.1	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 52.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.055 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4126	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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, OKX, MG

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	0.24	0/3461	0.42	0/4669
2	T	0.60	0/297	0.89	1/457 (0.2%)
3	P	0.40	0/240	0.86	0/369
All	All	0.30	0/3998	0.51	1/5495 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	6	DT	P-O3'-C3'	5.12	125.84	119.70

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	3380	0	3436	89	0
2	T	266	0	146	17	0
3	P	214	0	114	4	0
4	P	28	0	15	3	0
5	A	2	0	0	0	0
6	A	18	0	24	2	0
7	A	189	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	P	15	0	0	0	0
7	T	14	0	0	0	0
All	All	4126	0	3735	104	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 14.

All (104) 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:62:SER:HB2	2:T:3[B]:DG:C2	1.98	0.99
1:A:382:ARG:HG2	1:A:382:ARG:HH11	1.26	0.97
1:A:382:ARG:HH11	1:A:382:ARG:CG	1.81	0.93
1:A:322:SER:HB3	1:A:423:PHE:HD1	1.33	0.92
2:T:3[B]:DG:N3	2:T:3[B]:DG:H2'	1.84	0.90
2:T:2:DC:H4'	2:T:3[B]:DG:OP1	1.71	0.88
1:A:372:VAL:HB	1:A:417:PRO:HG2	1.65	0.77
1:A:382:ARG:HG2	1:A:382:ARG:NH1	1.96	0.77
1:A:383:ARG:HB3	1:A:402:VAL:CG1	2.15	0.76
1:A:409:SER:OG	1:A:411:ILE:HD12	1.88	0.73
1:A:371:ARG:NH2	1:A:375:ASP:O	2.23	0.71
1:A:84:ARG:NH1	7:A:751:HOH:O	2.23	0.70
1:A:158:GLU:O	1:A:161:VAL:HG22	1.92	0.70
1:A:381:LEU:HD22	1:A:406:CYS:SG	2.32	0.69
1:A:122:THR:O	1:A:126:GLN:HG2	1.93	0.68
1:A:322:SER:HB3	1:A:423:PHE:CD1	2.24	0.67
2:T:2:DC:C4'	2:T:3[B]:DG:OP1	2.44	0.65
1:A:412:GLN:O	1:A:412:GLN:HG2	1.96	0.65
1:A:62:SER:HB2	2:T:3[B]:DG:N1	2.12	0.64
1:A:383:ARG:HB3	1:A:402:VAL:HG11	1.80	0.62
2:T:7:DC:H2''	2:T:8:DA:H5'	1.82	0.62
2:T:3[A]:DG:N2	2:T:4[A]:DG:O6	2.31	0.60
1:A:181:ASP:OD1	1:A:182:ASN:N	2.34	0.60
2:T:3[B]:DG:N3	2:T:3[B]:DG:C2'	2.60	0.59
1:A:61[B]:ARG:HG3	2:T:3[B]:DG:O6	2.03	0.58
1:A:99:VAL:HG12	1:A:103[A]:MET:CE	2.33	0.58
1:A:131:LYS:NZ	7:A:784:HOH:O	2.36	0.58
1:A:62:SER:HA	2:T:3[B]:DG:C5	2.38	0.58
2:T:7:DC:H2'	2:T:8:DA:C8	2.38	0.58
1:A:423:PHE:CZ	1:A:425:CYS:HB2	2.39	0.57
1:A:25:GLN:O	1:A:27:PRO:HD3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:308:ASP:OD2	1:A:311:LYS:NZ	2.38	0.56
1:A:369:SER:HB3	1:A:423:PHE:HB3	1.87	0.56
1:A:179:GLN:CG	1:A:181:ASP:OD1	2.53	0.56
2:T:2:DC:H5''	2:T:3[A]:DG:OP1	2.04	0.56
1:A:318:THR:HG22	1:A:427:THR:HG22	1.88	0.55
1:A:381:LEU:HD11	1:A:383:ARG:CZ	2.36	0.55
1:A:371:ARG:HG2	1:A:420:THR:OG1	2.06	0.55
1:A:62:SER:OG	1:A:62:SER:O	2.23	0.55
1:A:368:VAL:HG21	1:A:399:ALA:HA	1.90	0.54
1:A:39:TYR:O	1:A:45:GLY:HA2	2.08	0.54
1:A:382:ARG:HB3	3:P:4:DT:OP2	2.08	0.53
1:A:376:LYS:HG2	1:A:376:LYS:O	2.10	0.52
1:A:230:ASN:O	1:A:234:ARG:HB2	2.11	0.51
1:A:371:ARG:HB3	1:A:421:MET:HB3	1.93	0.50
1:A:371:ARG:HD3	1:A:378:LEU:O	2.11	0.50
1:A:40:LYS:HD3	1:A:79:GLN:OE1	2.11	0.50
1:A:325:PHE:CD2	1:A:331:LEU:HD21	2.47	0.50
1:A:99:VAL:HG12	1:A:103[A]:MET:HE2	1.92	0.49
1:A:371:ARG:NH2	1:A:377:ARG:O	2.45	0.49
1:A:110:GLU:HB2	1:A:311:LYS:O	2.13	0.49
1:A:61[B]:ARG:HH12	4:P:10:0KX:H5	1.61	0.49
1:A:334:ARG:HG3	1:A:400:PHE:CZ	2.48	0.49
1:A:383:ARG:CB	1:A:402:VAL:CG1	2.89	0.49
1:A:209:GLY:O	6:A:505:GOL:H31	2.13	0.49
1:A:24:ARG:HB2	6:A:505:GOL:O2	2.12	0.49
1:A:106:PHE:CG	1:A:200[B]:MET:HG2	2.48	0.49
1:A:106:PHE:CD1	1:A:200[B]:MET:HG2	2.47	0.48
1:A:189:GLN:OE1	7:A:619:HOH:O	2.20	0.48
1:A:109:ILE:HD11	1:A:314:GLN:HG2	1.95	0.48
1:A:99:VAL:HG12	1:A:103[A]:MET:HE1	1.95	0.48
1:A:381:LEU:HD11	1:A:383:ARG:NH1	2.28	0.48
4:P:10:0KX:H15	4:P:10:0KX:O5'	2.14	0.48
1:A:14:MET:HG2	1:A:115:ASP:O	2.13	0.48
1:A:179:GLN:HG3	1:A:181:ASP:OD1	2.15	0.47
1:A:352:LEU:HB3	1:A:390:TYR:CE1	2.49	0.47
1:A:354:LYS:NZ	1:A:358:ASP:OD2	2.41	0.47
4:P:10:0KX:H7	4:P:10:0KX:O2B	2.15	0.47
1:A:277:GLU:O	1:A:280:GLN:HG3	2.15	0.47
1:A:382:ARG:NH1	1:A:382:ARG:CG	2.51	0.47
1:A:266:VAL:HA	1:A:290:PHE:CZ	2.51	0.46
3:P:5:DG:C8	3:P:6:DT:H72	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:ARG:HG2	1:A:212:CYS:SG	2.56	0.45
1:A:339:TRP:O	1:A:343:GLN:HG2	2.16	0.45
1:A:365:GLN:OE1	1:A:428:LYS:HE2	2.16	0.45
2:T:2:DC:H4'	2:T:3[A]:DG:OP1	2.16	0.45
1:A:195:VAL:O	1:A:199:GLU:HG3	2.17	0.45
1:A:336:GLN:HG2	7:A:699:HOH:O	2.16	0.45
1:A:416:SER:HA	1:A:417:PRO:C	2.38	0.44
1:A:290:PHE:O	1:A:294:ASN:HB2	2.18	0.44
1:A:51:SER:O	1:A:55:ARG:HG3	2.16	0.44
1:A:411:ILE:HD13	1:A:414:GLU:HB2	2.00	0.44
1:A:342:LEU:O	1:A:346:GLN:HG3	2.18	0.44
1:A:103[B]:MET:HG2	1:A:200[B]:MET:CE	2.48	0.44
1:A:180:ILE:H	1:A:180:ILE:HG13	1.46	0.43
1:A:224:LYS:HE2	3:P:9[B]:DG:OP1	2.18	0.43
1:A:411:ILE:O	1:A:412:GLN:HB3	2.17	0.43
1:A:178:LEU:HD11	1:A:191:THR:OG1	2.18	0.43
2:T:2:DC:C4'	2:T:3[A]:DG:OP1	2.66	0.43
3:P:3:DG:H2''	3:P:4:DT:O5'	2.18	0.43
1:A:333:THR:OG1	1:A:336:GLN:HB2	2.19	0.43
1:A:146:TYR:OH	1:A:163:LYS:HG2	2.18	0.42
1:A:401:THR:O	1:A:404:LYS:HG2	2.19	0.42
1:A:371:ARG:HD2	1:A:377:ARG:O	2.20	0.42
1:A:254:LYS:HA	1:A:254:LYS:HD3	1.78	0.42
1:A:409:SER:CB	1:A:416:SER:OG	2.67	0.42
1:A:200[A]:MET:HE2	1:A:200[A]:MET:HB2	1.86	0.42
1:A:381:LEU:CD2	1:A:406:CYS:SG	3.05	0.42
2:T:2:DC:C5'	2:T:3[A]:DG:OP1	2.67	0.42
1:A:316:PRO:HA	2:T:8:DA:OP1	2.20	0.41
1:A:62:SER:CB	2:T:3[B]:DG:N1	2.82	0.41
1:A:132:LEU:HB2	1:A:135:GLN:HB3	2.02	0.41
1:A:135:GLN:O	1:A:180:ILE:HG21	2.21	0.41
1:A:367:VAL:HG11	1:A:382:ARG:NE	2.36	0.41

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	432/435 (99%)	424 (98%)	8 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	373/372 (100%)	352 (94%)	21 (6%)	25	19

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	SER
1	A	115	ASP
1	A	127	GLU
1	A	132	LEU
1	A	135	GLN
1	A	152	GLN
1	A	160	THR
1	A	168	LYS
1	A	179	GLN
1	A	180	ILE
1	A	182	ASN
1	A	268	GLU
1	A	280	GLN
1	A	311	LYS
1	A	371	ARG
1	A	375	ASP
1	A	382	ARG
1	A	409	SER
1	A	411	ILE

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Mol	Chain	Res	Type
1	A	412	GLN
1	A	413	THR

Some 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 ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	504	-	5,5,5	0.34	0	5,5,5	0.24	0
6	GOL	A	505	-	5,5,5	0.34	0	5,5,5	0.28	0
6	GOL	A	506	-	5,5,5	0.35	0	5,5,5	0.25	0
4	OKX	P	10	5	26,29,29	3.32	8 (30%)	31,45,45	1.50	7 (22%)

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	504	-	-	0/4/4/4	0/0/0/0
6	GOL	A	505	-	-	0/4/4/4	0/0/0/0
6	GOL	A	506	-	-	0/4/4/4	0/0/0/0
4	OKX	P	10	5	-	0/14/34/34	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	P	10	OKX	O3'-C3'	-4.23	1.34	1.43
4	P	10	OKX	C5'-C4'	-2.50	1.43	1.51
4	P	10	OKX	PG-O3G	-2.26	1.45	1.54
4	P	10	OKX	C2-N3	-2.18	1.33	1.38
4	P	10	OKX	C2'-C3'	-2.12	1.47	1.52
4	P	10	OKX	C3'-C4'	-2.10	1.47	1.53
4	P	10	OKX	PA-N3A	2.39	1.69	1.63
4	P	10	OKX	PA-O1A	14.80	1.63	1.46

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	P	10	OKX	C5-C4-N3	-2.72	118.44	121.68
4	P	10	OKX	C2'-C1'-N1	-2.52	108.27	114.23
4	P	10	OKX	O2A-PA-O1A	-2.46	104.76	109.87
4	P	10	OKX	O5'-PA-N3A	2.02	111.52	107.27
4	P	10	OKX	O2A-PA-O5'	2.41	113.28	106.75
4	P	10	OKX	N4-C4-N3	2.71	121.20	116.64
4	P	10	OKX	O4'-C1'-N1	3.71	114.03	107.78

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	505	GOL	2	0
4	P	10	OKX	3	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, 95th 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 ²)	Q<0.9
1	A	428/435 (98%)	0.56	44 (10%)  	12, 24, 54, 83	0
2	T	12/12 (100%)	1.51	3 (25%)  	27, 30, 68, 91	0
3	P	9/9 (100%)	-0.14	0  	22, 27, 36, 48	0
All	All	449/456 (98%)	0.57	47 (10%)  	12, 24, 54, 91	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	T	3[A]	DG	7.8
2	T	1	DA	7.4
1	A	412	GLN	6.1
1	A	133	GLN	5.9
1	A	411	ILE	5.7
1	A	181	ASP	5.3
1	A	131	LYS	5.1
1	A	179	GLN	4.9
1	A	2	ALA	4.9
2	T	2	DC	4.7
1	A	134	GLY	4.2
1	A	410	GLY	4.1
1	A	182	ASN	3.9
1	A	328	LYS	3.9
1	A	408	THR	3.7
1	A	327	GLY	3.6
1	A	431	ALA	3.5
1	A	36	VAL	3.3
1	A	432	SER	3.3
1	A	180	ILE	3.3
1	A	184	THR	3.2
1	A	3	THR	3.2
1	A	160	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	376	LYS	3.0
1	A	130	GLN	3.0
1	A	158	GLU	2.9
1	A	378	LEU	2.9
1	A	135	GLN	2.9
1	A	19	VAL	2.8
1	A	417	PRO	2.8
1	A	183	LEU	2.6
1	A	373	GLN	2.6
1	A	413	THR	2.6
1	A	153	GLY	2.6
1	A	132	LEU	2.5
1	A	190	LEU	2.5
1	A	374	GLY	2.5
1	A	42	TRP	2.5
1	A	397	HIS	2.4
1	A	331	LEU	2.3
1	A	332	ALA	2.3
1	A	326	PRO	2.3
1	A	154	PRO	2.3
1	A	127	GLU	2.3
1	A	430	SER	2.2
1	A	6	ASP	2.2
1	A	377	ARG	2.2

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

There are no non-standard protein/DNA/RNA residues in this entry.

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, 95th 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(\AA^2)	Q<0.9
6	GOL	A	506	6/6	0.89	0.24	3.46	32,34,39,40	0
6	GOL	A	505	6/6	0.89	0.20	1.56	28,31,33,34	0
6	GOL	A	504	6/6	0.94	0.12	0.42	15,18,23,23	0
4	OKX	P	10	28/28	0.95	0.12	-0.50	12,16,20,21	0
5	MG	A	503	1/1	0.88	0.09	-1.31	18,18,18,18	0
5	MG	A	502	1/1	0.96	0.08	-2.57	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.