



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:38 AM UTC

PDB ID : 5KG3 / pdb_00005kg3
Title : Human DNA polymerase eta-DNA ternary complex: reaction first with 1 mM Mn²⁺ for 1800s then with 10 mM Mn²⁺ for 60s
Authors : Gao, Y.; Yang, W.
Deposited on : 2016-06-12
Resolution : 1.70 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

<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>.

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)
Xtriage (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

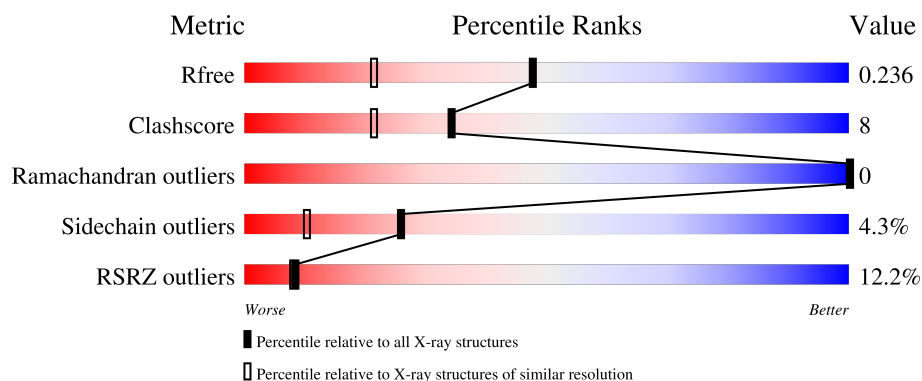
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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 ≥ 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	435	12% 78% 19% ..
2	T	12	25% 67% 33%
3	P	9	78% 22%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 4400 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	430	Total	C	N	O	S	0	16	0
			3420	2137	622	632	29			

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(*CP*AP*TP*TP*AP*TP*GP*AP*CP*GP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	12	Total	C	N	O	P	0	0	1
			226	108	39	68	11			

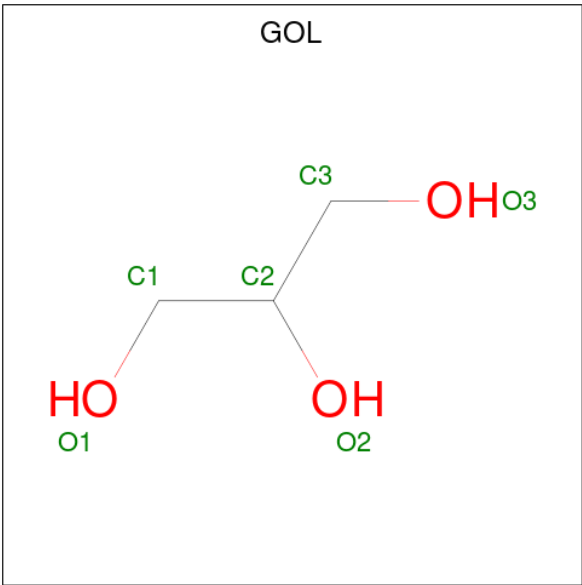
- Molecule 3 is a DNA chain called DNA (5'-D(*AP*GP*CP*GP*TP*CP*AP*TP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	9	Total	C	N	O	P	0	4	0
			228	109	39	69	11			

- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

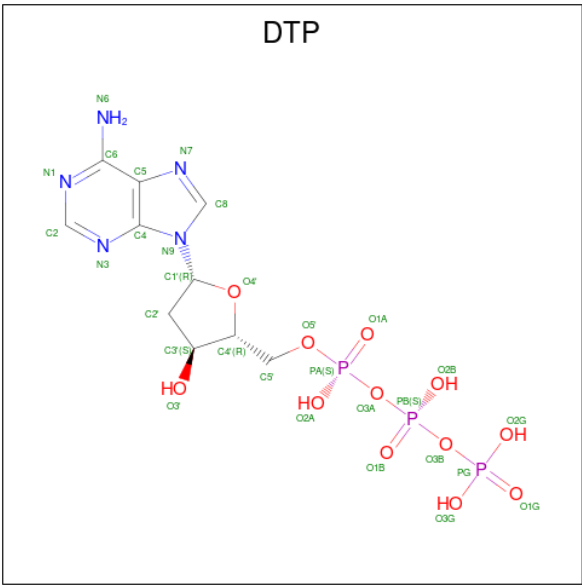
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mn	0	1
			2	2		
4	P	1	Total	Mn	0	1
			1	1		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



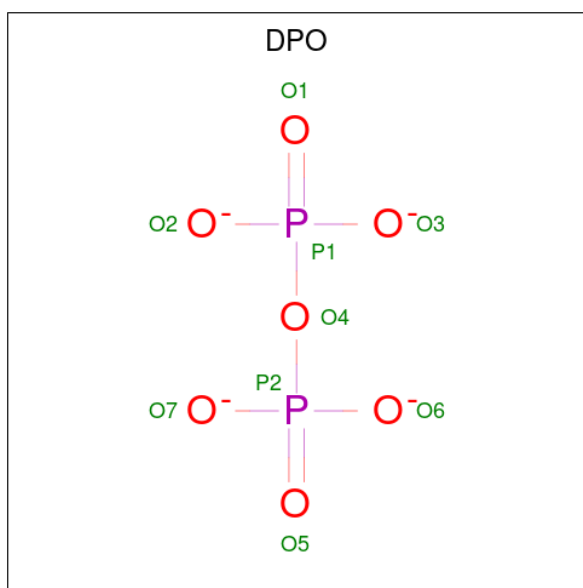
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is 2'-DEOXYADENOSINE 5'-TRIPHOSPHATE (CCD ID: DTP) (formula: C₁₀H₁₆N₅O₁₂P₃)



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	1
			30	10	5	12	3		

- Molecule 7 is DIPHOSPHATE (CCD ID: DPO) (formula: O_7P_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	P	0	1
			9	7	2		

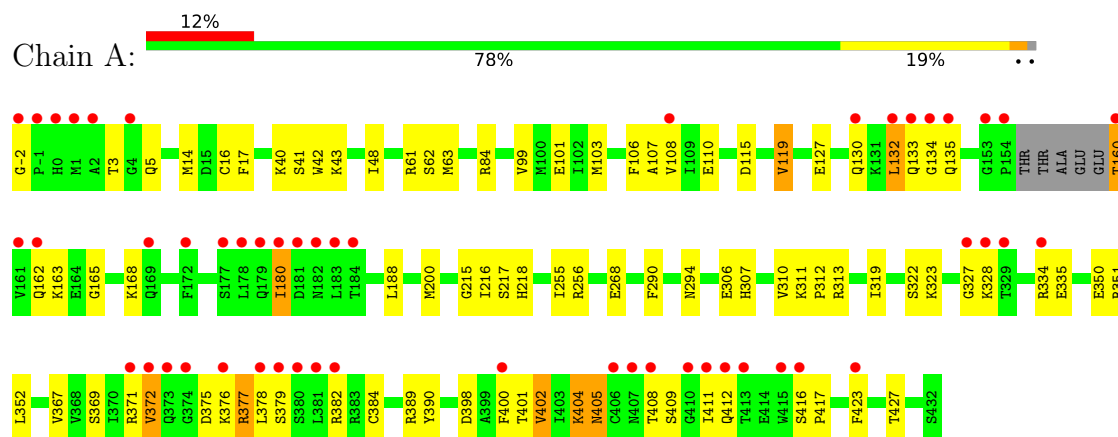
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	405	Total	O	0	4
			406	406		
8	T	32	Total	O	0	0
			32	32		
8	P	28	Total	O	0	0
			28	28		

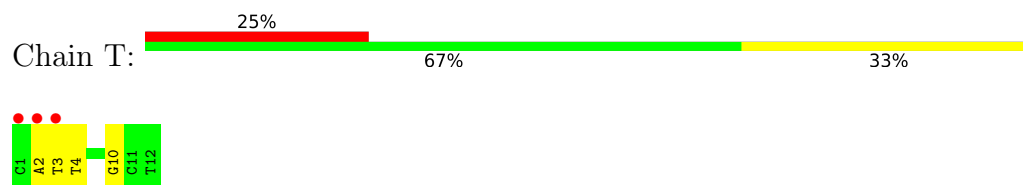
3 Residue-property plots

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.

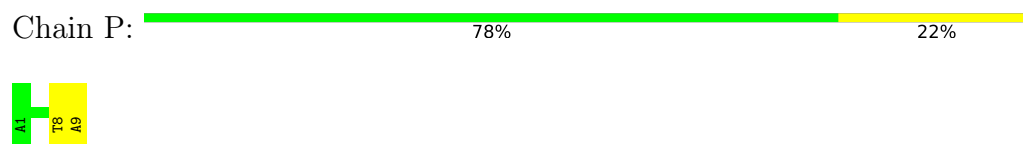
- Molecule 1: DNA polymerase eta



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	98.21Å 98.21Å 82.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.98 – 1.70 19.98 – 1.70	Depositor EDS
% Data completeness (in resolution range)	98.9 (19.98-1.70) 98.9 (19.98-1.70)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 1.70Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.197 , 0.237 0.198 , 0.236	Depositor DCC
R_{free} test set	3798 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å ²)	15.0	Xtriage
Anisotropy	0.109	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.048 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4400	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.35% 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: DPO, DTP, GOL, MN

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	0.84	15/3551 (0.4%)	0.88	10/4787 (0.2%)
2	T	0.39	0/252	0.60	0/388
3	P	0.32	0/291	0.52	0/441
All	All	0.80	15/4094 (0.4%)	0.84	10/5616 (0.2%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	217	SER	C-O	-10.45	1.16	1.24
1	A	310	VAL	C-O	-8.47	1.14	1.23
1	A	63	MET	C-O	-7.83	1.14	1.23
1	A	16	CYS	C-O	-7.76	1.14	1.23
1	A	218	HIS	C-O	-7.57	1.14	1.24
1	A	14	MET	C-O	-7.45	1.14	1.23
1	A	255	ILE	C-O	-7.37	1.14	1.24
1	A	216	ILE	C-O	-7.20	1.16	1.24
1	A	256	ARG	C-O	-6.88	1.15	1.23
1	A	379	SER	CA-C	-6.86	1.44	1.52
1	A	17	PHE	C-O	-6.78	1.16	1.24
1	A	427	THR	C-O	-6.12	1.16	1.23
1	A	312	PRO	N-CD	6.03	1.56	1.47
1	A	215	GLY	C-O	-5.23	1.16	1.23
1	A	327	GLY	C-O	-5.10	1.16	1.24

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	378	LEU	N-CA-C	9.86	122.03	111.28
1	A	377	ARG	CA-C-N	7.64	130.51	120.28
1	A	377	ARG	C-N-CA	7.64	130.51	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	379	SER	N-CA-C	-7.54	97.11	109.40
1	A	404	LYS	CB-CA-C	-7.01	99.15	110.79
1	A	134	GLY	CA-C-N	6.83	129.82	120.67
1	A	134	GLY	C-N-CA	6.83	129.82	120.67
1	A	-2	GLY	CA-C-N	6.82	126.50	119.82
1	A	-2	GLY	C-N-CA	6.82	126.50	119.82
1	A	17	PHE	N-CA-C	5.65	117.11	111.07

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	3420	0	3489	52	0
2	T	226	0	126	5	0
3	P	228	0	111	1	0
4	A	2	0	0	0	0
4	P	1	0	0	0	0
5	A	18	0	24	2	0
6	A	30	0	10	1	0
7	A	9	0	0	0	0
8	A	406	0	0	11	4
8	P	28	0	0	0	0
8	T	32	0	0	1	0
All	All	4400	0	3760	59	4

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 8.

All (59) 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:401:THR:HA	1:A:404:LYS:HD2	1.51	0.92
1:A:382:ARG:NH2	8:A:601:HOH:O	2.13	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:ARG:HH21	1:A:412:GLN:HB3	1.51	0.76
1:A:382:ARG:NH1	8:A:605:HOH:O	2.20	0.75
1:A:99:VAL:HG12	1:A:103[A]:MET:HE2	1.74	0.69
2:T:10:DG:N7	8:T:101:HOH:O	2.26	0.68
1:A:334:ARG:NH2	1:A:412:GLN:HB3	2.09	0.67
1:A:375:ASP:OD1	1:A:377:ARG:HG3	2.01	0.61
1:A:3:THR:HB	1:A:5:GLN:HE21	1.64	0.61
1:A:375:ASP:OD1	1:A:377:ARG:HB2	2.01	0.61
1:A:101:GLU:OE1	8:A:603:HOH:O	2.16	0.60
1:A:372:VAL:HG11	1:A:417:PRO:HG2	1.84	0.59
1:A:132:LEU:O	1:A:135:GLN:HB2	2.09	0.53
1:A:61[A]:ARG:NH1	8:A:607[A]:HOH:O	2.31	0.53
1:A:416:SER:HA	1:A:417:PRO:C	2.34	0.52
5:A:503:GOL:H32	8:A:635:HOH:O	2.09	0.51
1:A:61[B]:ARG:NH1	2:T:3:DT:O4	2.44	0.51
1:A:130:GLN:O	1:A:133:GLN:NE2	2.43	0.51
1:A:389:ARG:NH1	8:A:619:HOH:O	2.44	0.51
1:A:268:GLU:OE1	8:A:604:HOH:O	2.19	0.50
1:A:372:VAL:CG1	1:A:417:PRO:HG2	2.42	0.49
5:A:505:GOL:O2	8:A:602:HOH:O	2.16	0.49
1:A:62:SER:HB2	2:T:3:DT:C5	2.48	0.49
1:A:127:GLU:O	1:A:130:GLN:HG2	2.13	0.49
1:A:162:GLN:HG3	1:A:165:GLY:H	1.77	0.49
1:A:352:LEU:HB3	1:A:390:TYR:CE1	2.48	0.48
1:A:323:LYS:NZ	8:A:621:HOH:O	2.46	0.48
2:T:3:DT:H2'	2:T:4:DT:H71	1.96	0.48
1:A:160:THR:HG23	1:A:162:GLN:HG2	1.94	0.48
3:P:8[B]:DT:H2'	3:P:9[B]:DA:C8	2.48	0.48
1:A:375:ASP:OD1	1:A:377:ARG:CB	2.61	0.48
1:A:375:ASP:OD1	1:A:377:ARG:CG	2.62	0.47
1:A:42:TRP:CD1	1:A:43:LYS:HG3	2.49	0.47
1:A:398:ASP:O	1:A:402:VAL:HG13	2.13	0.47
1:A:180:ILE:H	1:A:180:ILE:HG13	1.53	0.47
1:A:322[A]:SER:HB3	1:A:423:PHE:CD2	2.50	0.47
1:A:371:ARG:NE	1:A:377:ARG:O	2.48	0.46
1:A:405:ASN:OD1	1:A:405:ASN:N	2.40	0.46
1:A:400:PHE:O	1:A:404:LYS:HG3	2.17	0.44
1:A:367:VAL:HG22	1:A:384[B]:CYS:SG	2.57	0.44
1:A:48:ILE:HA	1:A:61[B]:ARG:HG3	1.99	0.44
1:A:290:PHE:O	1:A:294:ASN:HB2	2.18	0.44
1:A:401:THR:CA	1:A:404:LYS:HD2	2.37	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:GLU:HG2	1:A:307:HIS:N	2.33	0.43
1:A:369:SER:HB3	1:A:423:PHE:HB3	2.00	0.43
1:A:180:ILE:HA	1:A:188:LEU:HD22	2.00	0.43
1:A:133:GLN:OE1	1:A:133:GLN:HA	2.19	0.43
6:A:506[A]:DTP:H8	6:A:506[A]:DTP:O5'	2.19	0.43
1:A:319:ILE:HG23	1:A:351[B]:ARG:CZ	2.48	0.43
1:A:311:LYS:CD	1:A:313[A]:ARG:HH22	2.30	0.43
1:A:409:SER:OG	1:A:411:ILE:O	2.37	0.43
1:A:132:LEU:HD12	1:A:132:LEU:HA	1.78	0.43
1:A:106:PHE:CG	1:A:200[B]:MET:HG2	2.54	0.42
1:A:3:THR:HB	1:A:5:GLN:NE2	2.31	0.42
1:A:40:LYS:HE2	8:A:864:HOH:O	2.18	0.42
8:A:694:HOH:O	2:T:2:DA:H2''	2.20	0.42
1:A:110:GLU:HB2	1:A:311:LYS:O	2.20	0.41
1:A:107:ALA:HB3	1:A:119:VAL:CG1	2.50	0.41
1:A:84:ARG:NH2	1:A:350:GLU:OE1	2.54	0.41

All (4) 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)
8:A:619:HOH:O	8:A:882:HOH:O[4_654]	2.10	0.10
8:A:700:HOH:O	8:A:748:HOH:O[5_664]	2.13	0.07
8:A:731:HOH:O	8:A:904:HOH:O[5_554]	2.15	0.05
8:A:751:HOH:O	8:A:973:HOH:O[4_655]	2.19	0.01

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	442/435 (102%)	436 (99%)	6 (1%)	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	384/372 (103%)	368 (96%)	16 (4%)	26 11

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	SER
1	A	108	VAL
1	A	115	ASP
1	A	119	VAL
1	A	132	LEU
1	A	160	THR
1	A	163	LYS
1	A	168	LYS
1	A	180	ILE
1	A	328	LYS
1	A	335	GLU
1	A	372	VAL
1	A	376	LYS
1	A	402	VAL
1	A	405	ASN
1	A	408	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	126	GLN
1	A	245	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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	$\# Z > 2$
5	GOL	A	504	-	5,5,5	0.37	0	5,5,5	0.32	0
6	DTP	A	506[A]	4	31,32,32	1.45	5 (16%)	46,50,50	1.73	9 (19%)
5	GOL	A	505	-	5,5,5	0.37	0	5,5,5	0.24	0
7	DPO	A	507[B]	4	6,8,8	0.81	0	12,13,13	0.65	0
5	GOL	A	503	-	5,5,5	0.60	0	5,5,5	0.51	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
5	GOL	A	504	-	-	0/4/4/4	-
6	DTP	A	506[A]	4	-	5/22/34/34	0/3/3/3
5	GOL	A	505	-	-	2/4/4/4	-
7	DPO	A	507[B]	4	-	2/6/6/6	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	A	503	-	-	2/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	506[A]	DTP	C6-N6	3.65	1.43	1.34
6	A	506[A]	DTP	C3'-C4'	-3.06	1.44	1.53
6	A	506[A]	DTP	C1'-N9	-2.48	1.41	1.46
6	A	506[A]	DTP	C5-N7	-2.46	1.34	1.39
6	A	506[A]	DTP	C2'-C3'	-2.17	1.47	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	506[A]	DTP	N3-C2-N1	-5.79	119.81	128.58
6	A	506[A]	DTP	C5-C4-N3	-4.66	120.30	126.72
6	A	506[A]	DTP	C5-N7-C8	3.52	108.97	103.45
6	A	506[A]	DTP	N3-C4-N9	3.49	133.10	127.17
6	A	506[A]	DTP	C2-N3-C4	3.44	120.24	111.83
6	A	506[A]	DTP	C4-C5-N7	-2.92	107.25	110.58
6	A	506[A]	DTP	N9-C8-N7	-2.76	110.02	113.94
6	A	506[A]	DTP	C2'-C3'-C4'	2.25	107.36	102.80
6	A	506[A]	DTP	C2-N1-C6	2.03	122.06	118.73

There are no chirality outliers.

All (11) torsion outliers are listed below:

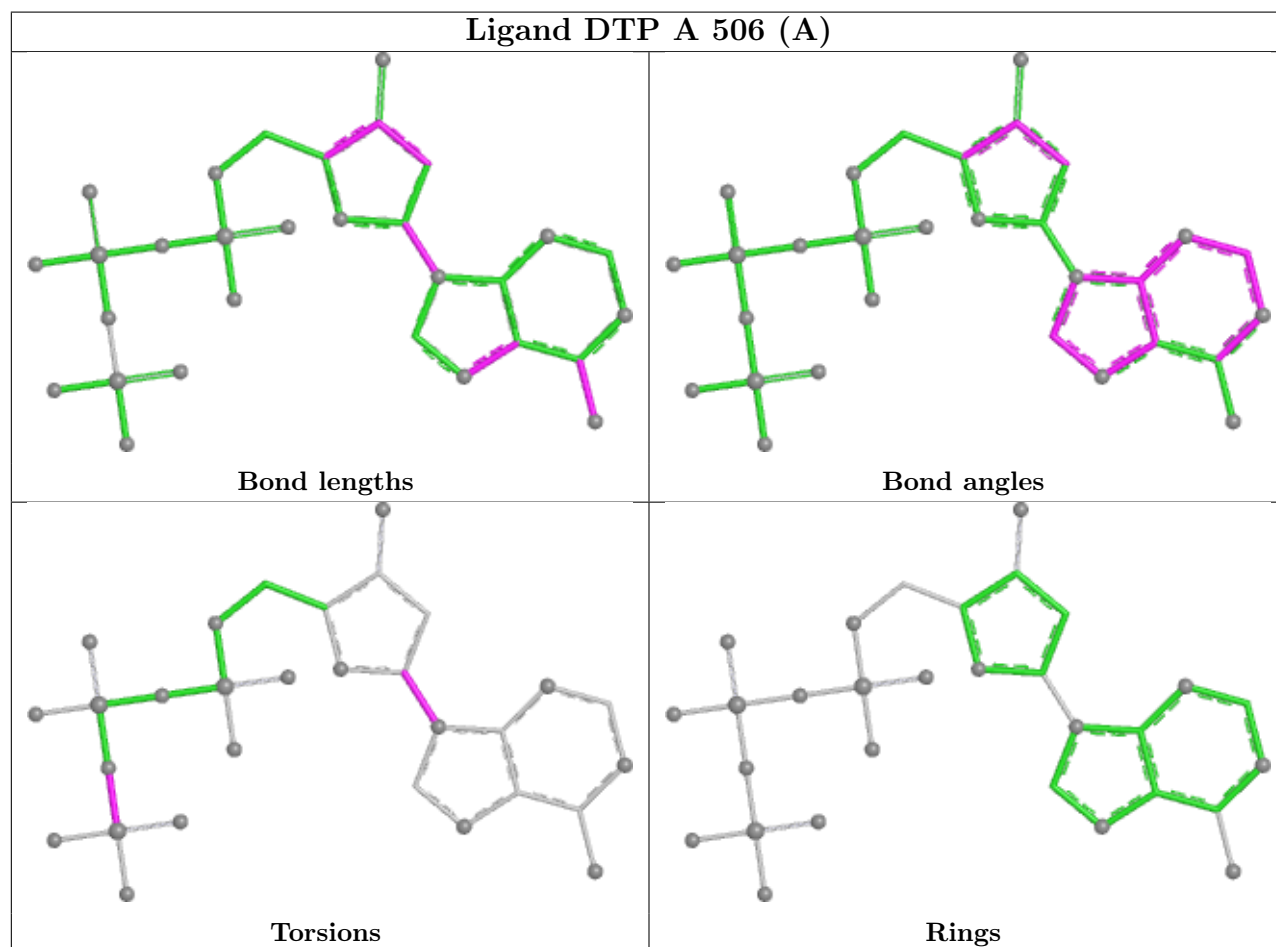
Mol	Chain	Res	Type	Atoms
5	A	503	GOL	O1-C1-C2-C3
5	A	505	GOL	O1-C1-C2-C3
5	A	505	GOL	O1-C1-C2-O2
5	A	503	GOL	O1-C1-C2-O2
6	A	506[A]	DTP	PB-O3B-PG-O1G
6	A	506[A]	DTP	C2'-C1'-N9-C4
7	A	507[B]	DPO	P1-O4-P2-O5
6	A	506[A]	DTP	PB-O3B-PG-O2G
6	A	506[A]	DTP	PB-O3B-PG-O3G
7	A	507[B]	DPO	P1-O4-P2-O7
6	A	506[A]	DTP	O4'-C1'-N9-C4

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	506[A]	DTP	1	0
5	A	505	GOL	1	0
5	A	503	GOL	1	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

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	430/435 (98%)	0.55	52 (12%) 8 8	7, 16, 50, 83	16 (3%)
2	T	12/12 (100%)	0.65	3 (25%) 2 1	15, 20, 57, 68	0
3	P	9/9 (100%)	0.31	0 100 100	3, 17, 29, 30	3 (33%)
All	All	451/456 (98%)	0.55	55 (12%) 8 8	3, 17, 50, 83	19 (4%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	378	LEU	6.1
1	A	161	VAL	5.7
1	A	180	ILE	5.2
1	A	411	ILE	4.8
1	A	408	THR	4.6
1	A	413	THR	4.5
1	A	154	PRO	4.2
1	A	374	GLY	4.0
1	A	134	GLY	3.8
1	A	406	CYS	3.7
2	T	3	DT	3.6
1	A	182	ASN	3.4
1	A	1	MET	3.4
1	A	183	LEU	3.3
1	A	2	ALA	3.3
1	A	372	VAL	3.2
1	A	160	THR	3.1
1	A	-2	GLY	3.1
1	A	181	ASP	2.9
1	A	108	VAL	2.9
1	A	327	GLY	2.8
1	A	334	ARG	2.8
1	A	376	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	415	TRP	2.7
1	A	162	GLN	2.7
1	A	0	HIS	2.7
1	A	380	SER	2.7
1	A	153	GLY	2.7
1	A	172	PHE	2.6
1	A	184	THR	2.6
1	A	132	LEU	2.6
1	A	412	GLN	2.6
1	A	381	LEU	2.5
1	A	373	GLN	2.5
1	A	130	GLN	2.5
1	A	133	GLN	2.5
1	A	178	LEU	2.5
1	A	416	SER	2.5
1	A	179	GLN	2.4
1	A	423	PHE	2.4
1	A	169	GLN	2.3
1	A	379	SER	2.2
1	A	400	PHE	2.2
1	A	135	GLN	2.2
2	T	1	DC	2.2
1	A	-1	PRO	2.2
1	A	4	GLY	2.1
1	A	382	ARG	2.1
1	A	410	GLY	2.1
1	A	329	THR	2.1
2	T	2	DA	2.1
1	A	407	ASN	2.0
1	A	328	LYS	2.0
1	A	371	ARG	2.0
1	A	177	SER	2.0

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](#)

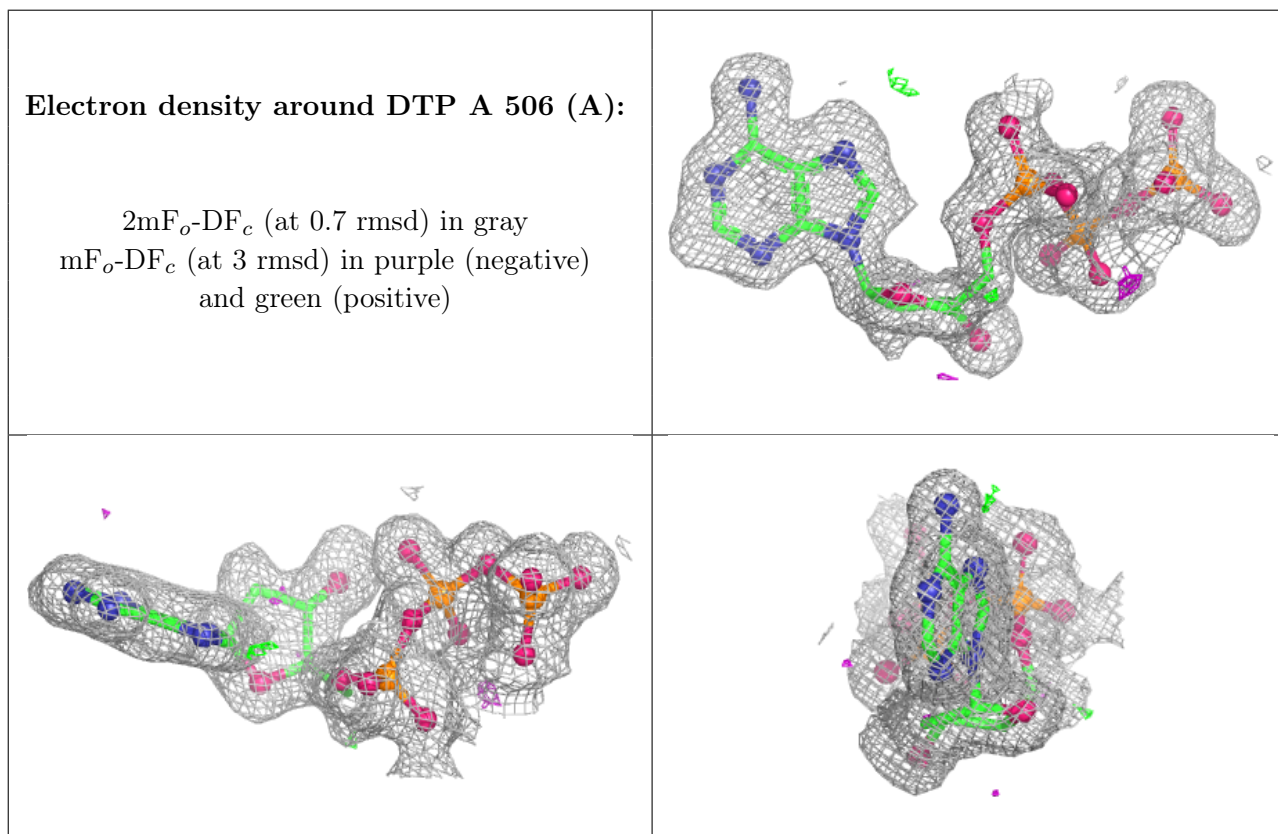
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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, 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	B-factors(Å ²)	Q<0.9
5	GOL	A	505	6/6	0.70	0.22	28,31,34,39	0
5	GOL	A	504	6/6	0.87	0.15	17,31,36,40	0
5	GOL	A	503	6/6	0.90	0.17	15,19,27,31	0
4	MN	P	101[B]	1/1	0.95	0.07	18,18,18,18	1
7	DPO	A	507[B]	9/9	0.96	0.06	7,10,10,10	9
6	DTP	A	506[A]	30/30	0.98	0.05	7,10,12,13	30
4	MN	A	501	1/1	0.99	0.02	10,10,10,10	0
4	MN	A	502[B]	1/1	0.99	0.03	9,9,9,9	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.