



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 21, 2022 – 12:57 PM EDT

PDB ID : 8E24
Title : Human DNA polymerase theta in complex with allosteric inhibitor
Authors : Mader, P.; Pau, V.P.T.; Sicheri, F.
Deposited on : 2022-08-13
Resolution : 2.34 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.29
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.29

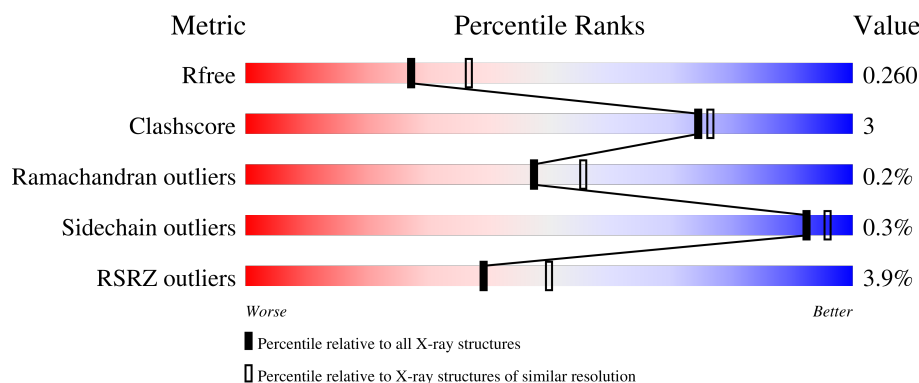
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 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}	130704	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

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	668	<div> <div>2%</div> <div>89%</div> <div>7%</div> <div>.</div> </div>
1	D	668	<div> <div>6%</div> <div>88%</div> <div>9%</div> <div>.</div> </div>
2	B	17	<div> <div>53%</div> <div>29%</div> <div>18%</div> </div>
2	E	17	<div> <div>65%</div> <div>12%</div> <div>24%</div> </div>
3	C	13	<div> <div>62%</div> <div>15%</div> <div>23%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	13	<div><div></div><div></div><div></div><div>23%</div><div>46%</div><div>31%</div></div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11101 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 theta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	644	Total	C	N	O	S	0	0	0
			4999	3181	856	932	30			
1	D	648	Total	C	N	O	S	0	0	0
			4889	3105	838	918	28			

There are 214 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1817	GLY	-	expression tag	UNP O75417
A	?	-	ALA	deletion	UNP O75417
A	?	-	THR	deletion	UNP O75417
A	?	-	ILE	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	PHE	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	GLN	deletion	UNP O75417
A	?	-	ALA	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	PRO	deletion	UNP O75417
A	?	-	GLN	deletion	UNP O75417
A	?	-	GLU	deletion	UNP O75417
A	?	-	ILE	deletion	UNP O75417
A	?	-	PRO	deletion	UNP O75417
A	?	-	ILE	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	ASP	deletion	UNP O75417
A	?	-	ASP	deletion	UNP O75417
A	?	-	GLN	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	HIS	deletion	UNP O75417

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	SER	deletion	UNP O75417
A	?	-	GLU	deletion	UNP O75417
A	?	-	ILE	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	ALA	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	LEU	deletion	UNP O75417
A	?	-	VAL	deletion	UNP O75417
A	?	-	PRO	deletion	UNP O75417
A	?	-	PRO	deletion	UNP O75417
A	?	-	MET	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	ASN	deletion	UNP O75417
A	?	-	GLN	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	THR	deletion	UNP O75417
A	?	-	LEU	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	THR	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	ILE	deletion	UNP O75417
A	?	-	ASP	deletion	UNP O75417
A	?	-	ASN	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	LEU	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	VAL	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	GLU	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	PRO	deletion	UNP O75417
A	?	-	PRO	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	GLN	deletion	UNP O75417

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	ALA	deletion	UNP O75417
A	?	-	VAL	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	LEU	deletion	UNP O75417
A	?	-	LEU	deletion	UNP O75417
A	?	-	PRO	deletion	UNP O75417
A	?	-	MET	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	TYR	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	GLY	deletion	UNP O75417
A	?	-	PHE	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	VAL	deletion	UNP O75417
A	?	-	ASN	deletion	UNP O75417
A	?	-	PRO	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	CYS	deletion	UNP O75417
A	?	-	GLN	deletion	UNP O75417
A	?	-	ALA	deletion	UNP O75417
A	?	-	GLN	deletion	UNP O75417
A	?	-	MET	deletion	UNP O75417
A	?	-	GLU	deletion	UNP O75417
A	?	-	GLU	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	ALA	deletion	UNP O75417
A	?	-	LEU	deletion	UNP O75417
A	?	-	SER	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	ARG	deletion	UNP O75417
A	?	-	LYS	deletion	UNP O75417
A	?	-	LEU	deletion	UNP O75417
A	?	-	GLN	deletion	UNP O75417
D	1817	GLY	-	expression tag	UNP O75417
D	?	-	ALA	deletion	UNP O75417

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Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	THR	deletion	UNP O75417
D	?	-	ILE	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	PHE	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	GLN	deletion	UNP O75417
D	?	-	ALA	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	PRO	deletion	UNP O75417
D	?	-	GLN	deletion	UNP O75417
D	?	-	GLU	deletion	UNP O75417
D	?	-	ILE	deletion	UNP O75417
D	?	-	PRO	deletion	UNP O75417
D	?	-	ILE	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	ASP	deletion	UNP O75417
D	?	-	ASP	deletion	UNP O75417
D	?	-	GLN	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	HIS	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	GLU	deletion	UNP O75417
D	?	-	ILE	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	ALA	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	LEU	deletion	UNP O75417
D	?	-	VAL	deletion	UNP O75417
D	?	-	PRO	deletion	UNP O75417
D	?	-	PRO	deletion	UNP O75417
D	?	-	MET	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	ASN	deletion	UNP O75417
D	?	-	GLN	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	THR	deletion	UNP O75417

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Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	LEU	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	THR	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	ILE	deletion	UNP O75417
D	?	-	ASP	deletion	UNP O75417
D	?	-	ASN	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	LEU	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	VAL	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	GLU	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	PRO	deletion	UNP O75417
D	?	-	PRO	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	GLN	deletion	UNP O75417
D	?	-	ALA	deletion	UNP O75417
D	?	-	VAL	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	LEU	deletion	UNP O75417
D	?	-	LEU	deletion	UNP O75417
D	?	-	PRO	deletion	UNP O75417
D	?	-	MET	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	TYR	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	GLY	deletion	UNP O75417
D	?	-	PHE	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417

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Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	VAL	deletion	UNP O75417
D	?	-	ASN	deletion	UNP O75417
D	?	-	PRO	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	CYS	deletion	UNP O75417
D	?	-	GLN	deletion	UNP O75417
D	?	-	ALA	deletion	UNP O75417
D	?	-	GLN	deletion	UNP O75417
D	?	-	MET	deletion	UNP O75417
D	?	-	GLU	deletion	UNP O75417
D	?	-	GLU	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	ALA	deletion	UNP O75417
D	?	-	LEU	deletion	UNP O75417
D	?	-	SER	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	ARG	deletion	UNP O75417
D	?	-	LYS	deletion	UNP O75417
D	?	-	LEU	deletion	UNP O75417
D	?	-	GLN	deletion	UNP O75417

- Molecule 2 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	14	Total	C	N	O	P	0	0	0
			282	135	54	80	13			
2	E	13	Total	C	N	O	P	0	0	0
			263	126	51	74	12			

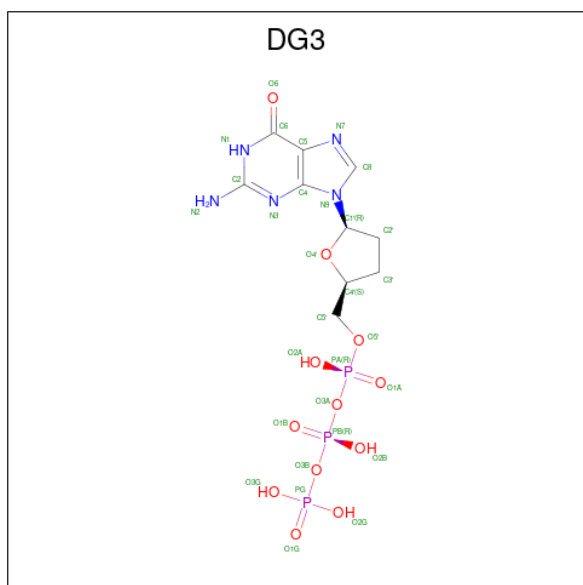
- Molecule 3 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	P	0	0	0
			204	98	34	62	10			
3	F	9	Total	C	N	O	P	0	0	0
			182	88	29	56	9			

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

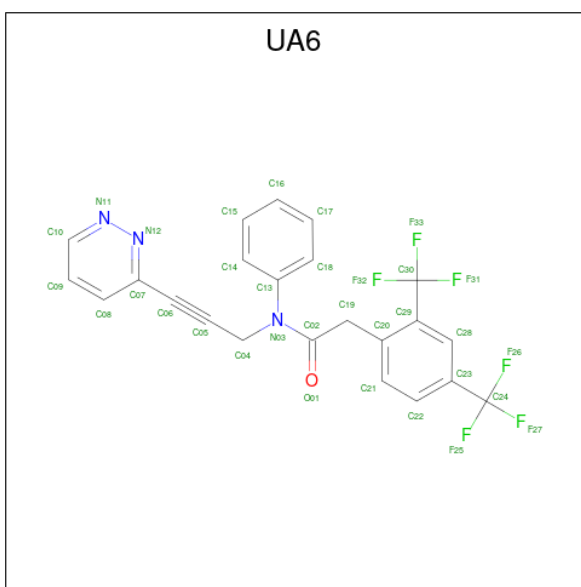
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Mg 1 1	0	0
4	D	1	Total Mg 1 1	0	0

- Molecule 5 is 2'-3'-DIDEOXYGUANOSINE-5'-TRIPHOSPHATE (three-letter code: DG3) (formula: C₁₀H₁₆N₅O₁₂P₃).



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

- Molecule 6 is 2-[2,4-bis(trifluoromethyl)phenyl]-N-phenyl-N-[3-(pyridazin-3-yl)prop-2-yn-1-yl]acetamide (three-letter code: UA6) (formula: C₂₃H₁₅F₆N₃O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	F	N	O	0	0
			33	23	6	3	1		
6	D	1	Total	C	F	N	O	0	0
			33	23	6	3	1		

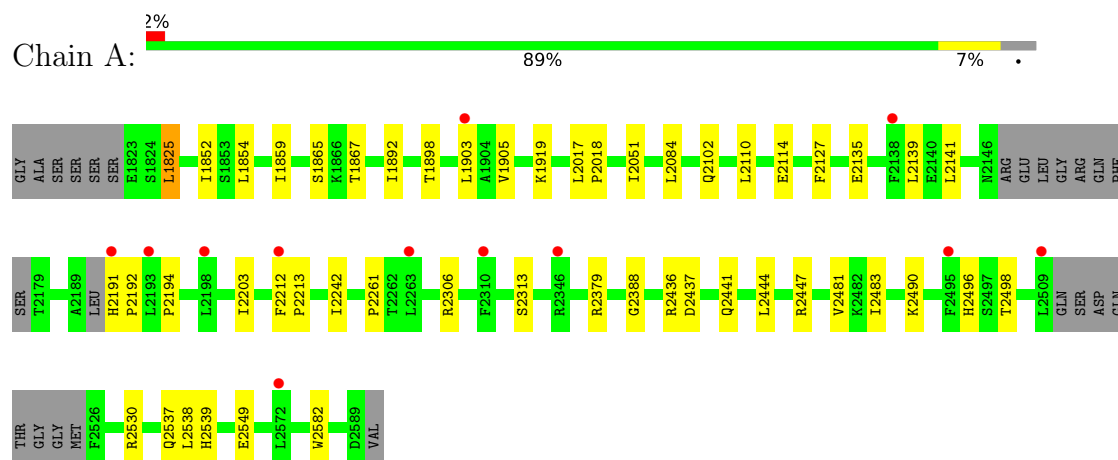
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	97	Total	O	0	0
			97	97		
7	B	3	Total	O	0	0
			3	3		
7	C	2	Total	O	0	0
			2	2		
7	D	49	Total	O	0	0
			49	49		
7	E	2	Total	O	0	0
			2	2		
7	F	1	Total	O	0	0
			1	1		

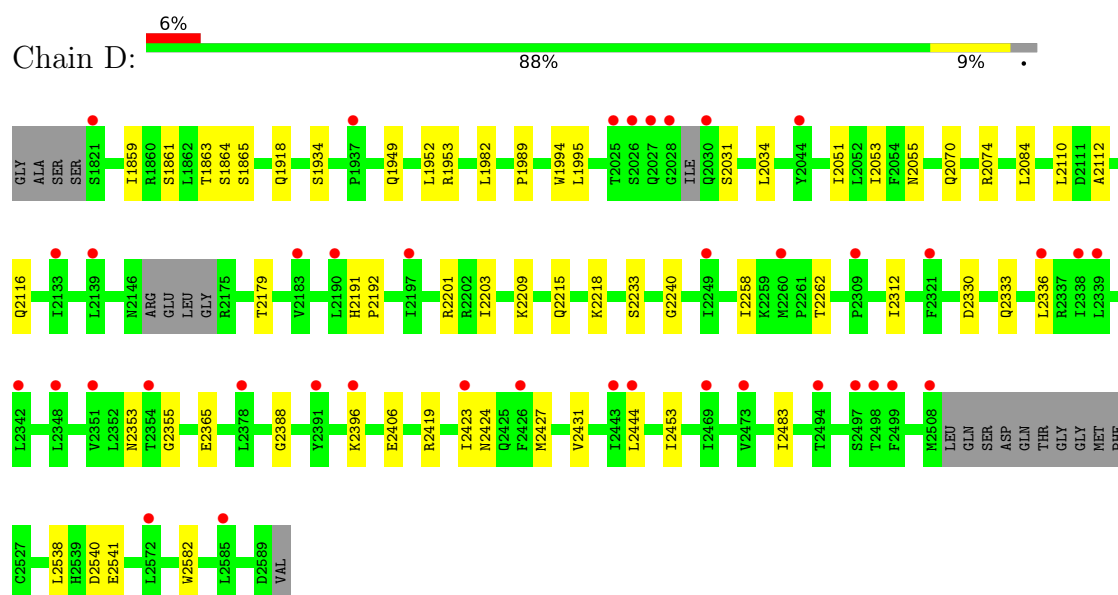
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: DNA polymerase theta

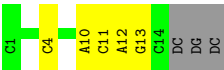


• Molecule 1: DNA polymerase theta

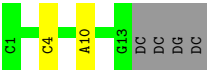


• Molecule 2: DNA





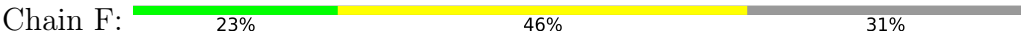
● Molecule 2: DNA



● Molecule 3: DNA



● Molecule 3: DNA



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	263.74Å 66.29Å 156.66Å 90.00° 126.14° 90.00°	Depositor
Resolution (Å)	126.52 – 2.34 126.52 – 2.34	Depositor EDS
% Data completeness (in resolution range)	80.0 (126.52-2.34) 80.0 (126.52-2.34)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.231 , 0.266 0.227 , 0.260	Depositor DCC
R_{free} test set	3686 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	59.5	Xtriage
Anisotropy	0.156	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.034 for -h-2*k,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11101	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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: UA6, MG, DG3

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	0/5096	0.44	0/6895
1	D	0.23	0/4984	0.44	0/6762
2	B	0.45	0/316	0.79	0/485
2	E	0.47	0/295	0.83	0/453
3	C	0.46	0/227	0.96	0/348
3	F	0.48	0/202	0.98	0/309
All	All	0.27	0/11120	0.50	0/15252

There are no bond length outliers.

There are no bond angle outliers.

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	4999	0	4916	27	0
1	D	4889	0	4647	34	0
2	B	282	0	158	3	0
2	E	263	0	147	2	0
3	C	204	0	113	1	0
3	F	182	0	102	3	0
4	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1	0	0	0	0
5	A	30	0	12	1	0
5	D	30	0	12	1	0
6	A	33	0	0	0	0
6	D	33	0	0	0	0
7	A	97	0	0	0	0
7	B	3	0	0	0	0
7	C	2	0	0	0	0
7	D	49	0	0	0	0
7	E	2	0	0	0	0
7	F	1	0	0	0	0
All	All	11101	0	10107	68	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 3.

All (68) 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:D:2330:ASP:HB3	1:D:2541:GLU:HG2	1.53	0.89
1:A:1852:ILE:HB	1:A:1903:LEU:HD11	1.70	0.74
2:B:10:DA:H2''	2:B:11:DC:H5''	1.79	0.65
3:F:9:DC:H2'	3:F:10:DA:C8	2.33	0.64
1:A:1867:THR:HB	1:A:1919:LYS:HD3	1.84	0.59
1:A:1854:LEU:HD23	1:A:1903:LEU:HD13	1.85	0.58
3:C:4:DG:H2''	3:C:5:DC:H5''	1.85	0.57
1:A:2102:GLN:NE2	1:A:2313:SER:O	2.35	0.57
1:D:2538:LEU:HB2	1:D:2541:GLU:HB2	1.86	0.57
1:A:2135:GLU:HG3	1:A:2139:LEU:HD23	1.86	0.56
1:D:2209:LYS:HE3	2:E:10:DA:H4'	1.88	0.55
1:D:2070:GLN:HG2	1:D:2074:ARG:HE	1.72	0.54
1:A:2084:LEU:HD11	1:A:2242:ILE:HD13	1.89	0.54
1:A:2538:LEU:HD11	1:A:2582:TRP:HZ2	1.72	0.54
1:D:2215:GLN:HA	1:D:2218:LYS:HE3	1.90	0.53
1:D:1952:LEU:HB3	1:D:1982:LEU:HD23	1.91	0.53
1:D:2431:VAL:HG13	1:D:2453:ILE:HD11	1.90	0.53
1:D:2051:ILE:O	1:D:2055:ASN:ND2	2.42	0.52
1:A:2496:HIS:HB3	1:A:2498:THR:HG22	1.92	0.52
1:A:2191:HIS:O	1:A:2194:PRO:HD2	2.10	0.51
1:A:2530:ARG:CZ	1:A:2549:GLU:HG3	2.40	0.51
1:D:2538:LEU:HD11	1:D:2582:TRP:HZ2	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:2538:LEU:O	1:D:2540:ASP:N	2.41	0.49
1:A:2379:ARG:NH1	5:A:2601:DG3:O2G	2.44	0.47
1:D:1863:THR:O	1:D:1865:SER:N	2.47	0.47
1:A:2388:GLY:HA2	2:B:4:DC:C2	2.49	0.47
1:D:2424:ASN:HA	1:D:2427:MET:HE2	1.97	0.47
1:A:2212:PHE:HB3	1:A:2213:PRO:HD3	1.98	0.46
1:D:2538:LEU:HD11	1:D:2582:TRP:CZ2	2.50	0.46
1:A:2110:LEU:HD21	1:A:2203:ILE:HB	1.98	0.46
1:A:1825:LEU:HD12	1:A:2051:ILE:HD12	1.98	0.46
1:D:1859:ILE:O	1:D:1861:SER:N	2.47	0.46
1:A:2436:ARG:NH1	1:A:2437:ASP:OD1	2.49	0.45
3:F:6:DT:H2''	3:F:7:DG:C8	2.51	0.45
1:D:2110:LEU:HD21	1:D:2203:ILE:HB	1.99	0.45
1:A:2261:PRO:HA	1:A:2306:ARG:HA	1.98	0.45
1:A:2490:LYS:HE3	1:A:2490:LYS:HB2	1.75	0.45
1:D:2365:GLU:OE2	1:D:2419:ARG:NH2	2.50	0.45
1:D:2396:LYS:HA	1:D:2406:GLU:HG2	1.98	0.45
1:D:1994:TRP:HZ2	1:D:2240:GLY:HA2	1.81	0.45
1:A:2444:LEU:HD11	1:A:2483:ILE:HD11	1.99	0.44
1:A:2441:GLN:HG2	1:A:2447:ARG:HG2	1.99	0.44
1:A:2017:LEU:N	1:A:2018:PRO:HD2	2.33	0.44
1:D:2112:ALA:O	1:D:2116:GLN:HG2	2.18	0.44
1:D:2258:ILE:HD13	1:D:2312:ILE:HD11	2.00	0.44
1:A:1859:ILE:HG23	1:A:1865:SER:OG	2.18	0.43
1:D:1918:GLN:HE21	1:D:1934:SER:HA	1.84	0.43
1:D:2423:ILE:O	1:D:2427:MET:HG3	2.18	0.43
1:A:1852:ILE:HG22	1:A:1905:VAL:HG22	2.00	0.43
1:D:2031:SER:HB3	1:D:2034:LEU:HG	2.00	0.43
3:F:11:DT:H2'	3:F:12:DT:C6	2.54	0.43
1:D:2179:THR:HG22	1:D:2201:ARG:HH22	1.84	0.43
1:A:1892:ILE:HD12	1:A:1898:THR:HG22	1.99	0.43
1:A:2481:VAL:HG21	1:A:2539:HIS:O	2.19	0.42
1:D:2084:LEU:HD13	1:D:2233:SER:HB2	2.00	0.42
1:D:1995:LEU:HD23	1:D:1995:LEU:HA	1.92	0.42
1:D:2191:HIS:CG	1:D:2192:PRO:HD2	2.55	0.42
1:A:2141:LEU:HD13	1:A:2191:HIS:CE1	2.54	0.42
1:D:2388:GLY:HA2	2:E:4:DC:C2	2.55	0.42
1:D:1949:GLN:NE2	1:D:1953:ARG:HD2	2.35	0.42
1:D:2353:ASN:C	1:D:2355:GLY:H	2.24	0.42
1:A:2114:GLU:HG2	1:A:2127:PHE:CE1	2.54	0.41
1:D:2444:LEU:HD11	1:D:2483:ILE:HD11	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:12:DA:H2''	2:B:13:DG:N7	2.36	0.41
5:D:2601:DG3:H8	5:D:2601:DG3:O5'	2.21	0.41
1:D:1994:TRP:CZ2	1:D:2240:GLY:HA2	2.55	0.41
1:D:1989:PRO:HG3	1:D:2053:ILE:HG23	2.02	0.41
1:D:2333:GLN:HB3	1:D:2336:LEU:HD12	2.02	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	636/668 (95%)	617 (97%)	18 (3%)	1 (0%)	47	55
1	D	640/668 (96%)	621 (97%)	18 (3%)	1 (0%)	47	55
All	All	1276/1336 (96%)	1238 (97%)	36 (3%)	2 (0%)	47	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	1864	SER
1	A	2192	PRO

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	535/586 (91%)	533 (100%)	2 (0%)	91	95
1	D	498/586 (85%)	497 (100%)	1 (0%)	93	96
All	All	1033/1172 (88%)	1030 (100%)	3 (0%)	92	96

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1825	LEU
1	A	2537	GLN
1	D	2262	THR

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

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	UA6	A	2602	-	35,35,35	1.76	7 (20%)	48,50,50	1.91	12 (25%)
5	DG3	D	2601	4	25,32,32	0.87	1 (4%)	28,50,50	1.58	5 (17%)
6	UA6	D	2602	-	35,35,35	1.85	7 (20%)	48,50,50	1.99	13 (27%)
5	DG3	A	2601	4	25,32,32	0.86	1 (4%)	28,50,50	1.50	6 (21%)

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	UA6	A	2602	-	-	1/28/30/30	0/3/3/3
5	DG3	D	2601	4	-	8/18/31/31	0/3/3/3
6	UA6	D	2602	-	-	1/28/30/30	0/3/3/3
5	DG3	A	2601	4	-	7/18/31/31	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	2602	UA6	N11-N12	-5.93	1.27	1.35
6	D	2602	UA6	N11-N12	-5.72	1.27	1.35
6	D	2602	UA6	C02-N03	3.60	1.42	1.36
6	D	2602	UA6	C19-C02	3.52	1.56	1.52
6	D	2602	UA6	C13-N03	3.49	1.50	1.43
6	A	2602	UA6	C13-N03	3.33	1.50	1.43
6	A	2602	UA6	C02-N03	3.17	1.41	1.36
6	A	2602	UA6	C19-C02	2.84	1.55	1.52
6	D	2602	UA6	C07-N12	2.57	1.38	1.34
6	A	2602	UA6	C07-N12	2.49	1.38	1.34
6	D	2602	UA6	C10-N11	2.43	1.38	1.32
6	A	2602	UA6	C10-N11	2.39	1.38	1.32
5	A	2601	DG3	C6-N1	-2.30	1.34	1.37
5	D	2601	DG3	C6-N1	-2.27	1.34	1.37
6	A	2602	UA6	C30-C29	2.25	1.55	1.50
6	D	2602	UA6	C30-C29	2.22	1.55	1.50

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	2602	UA6	C20-C19-C02	6.62	122.95	112.46
6	A	2602	UA6	C20-C19-C02	5.83	121.68	112.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	2602	UA6	C05-C04-N03	4.35	122.35	111.28
6	A	2602	UA6	C10-N11-N12	4.33	122.49	118.99
6	D	2602	UA6	C10-N11-N12	4.33	122.49	118.99
6	A	2602	UA6	C05-C04-N03	4.04	121.56	111.28
5	D	2601	DG3	PB-O3B-PG	-3.80	119.79	132.83
6	A	2602	UA6	C04-N03-C02	3.65	123.22	117.34
6	D	2602	UA6	C04-N03-C02	3.61	123.16	117.34
5	D	2601	DG3	PA-O3A-PB	-3.40	121.15	132.83
5	A	2601	DG3	PB-O3B-PG	-3.32	121.44	132.83
6	A	2602	UA6	F33-C30-C29	3.14	118.17	112.70
6	A	2602	UA6	O01-C02-N03	-2.99	117.31	121.66
6	D	2602	UA6	O01-C02-N03	-2.99	117.31	121.66
6	D	2602	UA6	C08-C07-N12	-2.89	119.93	123.25
5	A	2601	DG3	PA-O3A-PB	-2.84	123.08	132.83
6	A	2602	UA6	C08-C07-N12	-2.81	120.02	123.25
6	D	2602	UA6	F33-C30-C29	2.72	117.44	112.70
6	A	2602	UA6	C23-C28-C29	-2.61	119.75	123.80
6	D	2602	UA6	C18-C13-N03	2.53	123.87	120.18
6	D	2602	UA6	C23-C28-C29	-2.53	119.88	123.80
5	D	2601	DG3	C3'-C2'-C1'	2.51	105.68	102.78
5	A	2601	DG3	C3'-C2'-C1'	2.45	105.61	102.78
6	A	2602	UA6	C18-C13-N03	2.42	123.72	120.18
6	D	2602	UA6	C21-C20-C29	-2.40	116.44	118.76
6	A	2602	UA6	O01-C02-C19	-2.39	117.71	121.60
5	A	2601	DG3	C8-N7-C5	2.32	107.42	102.99
6	D	2602	UA6	O01-C02-C19	-2.31	117.84	121.60
5	D	2601	DG3	C8-N7-C5	2.25	107.27	102.99
5	A	2601	DG3	C5-C6-N1	2.25	117.92	113.95
6	D	2602	UA6	C17-C18-C13	2.24	122.60	119.68
5	D	2601	DG3	C5-C6-N1	2.23	117.89	113.95
6	A	2602	UA6	C17-C18-C13	2.23	122.59	119.68
6	D	2602	UA6	C28-C23-C24	2.15	122.48	119.58
6	A	2602	UA6	C21-C20-C29	-2.07	116.76	118.76
5	A	2601	DG3	C2'-C1'-N9	-2.01	108.70	112.48

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	2601	DG3	C5'-O5'-PA-O2A
5	D	2601	DG3	C5'-O5'-PA-O1A
5	D	2601	DG3	C5'-O5'-PA-O2A

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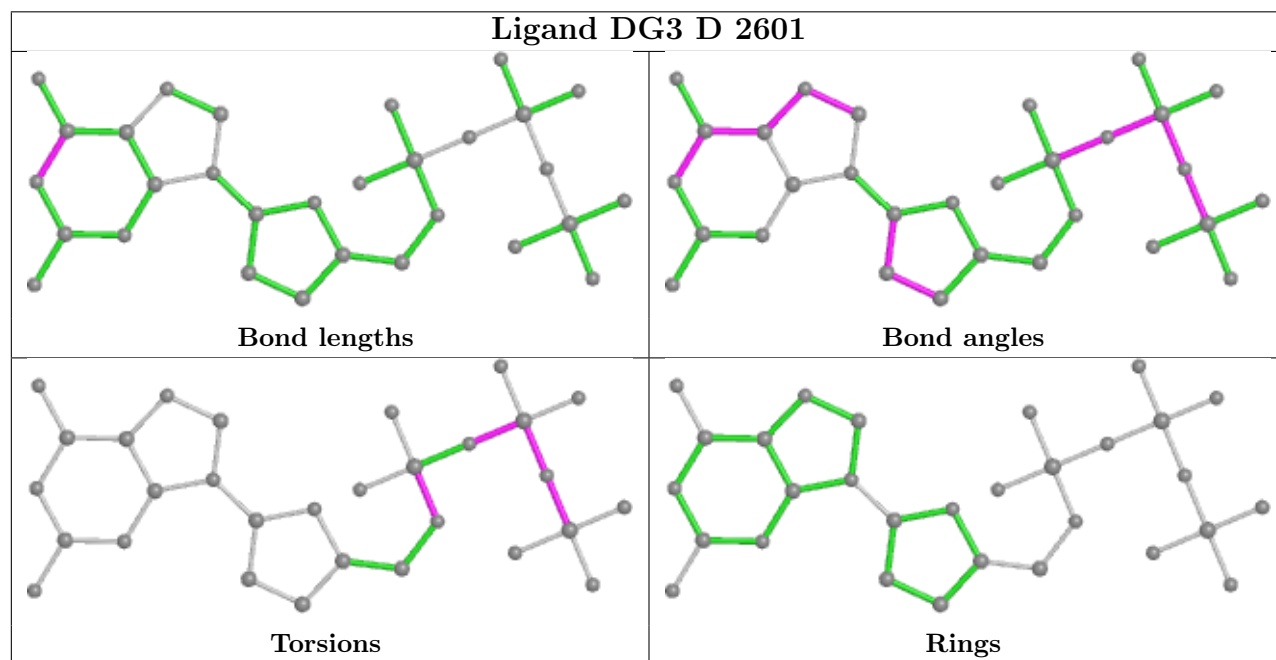
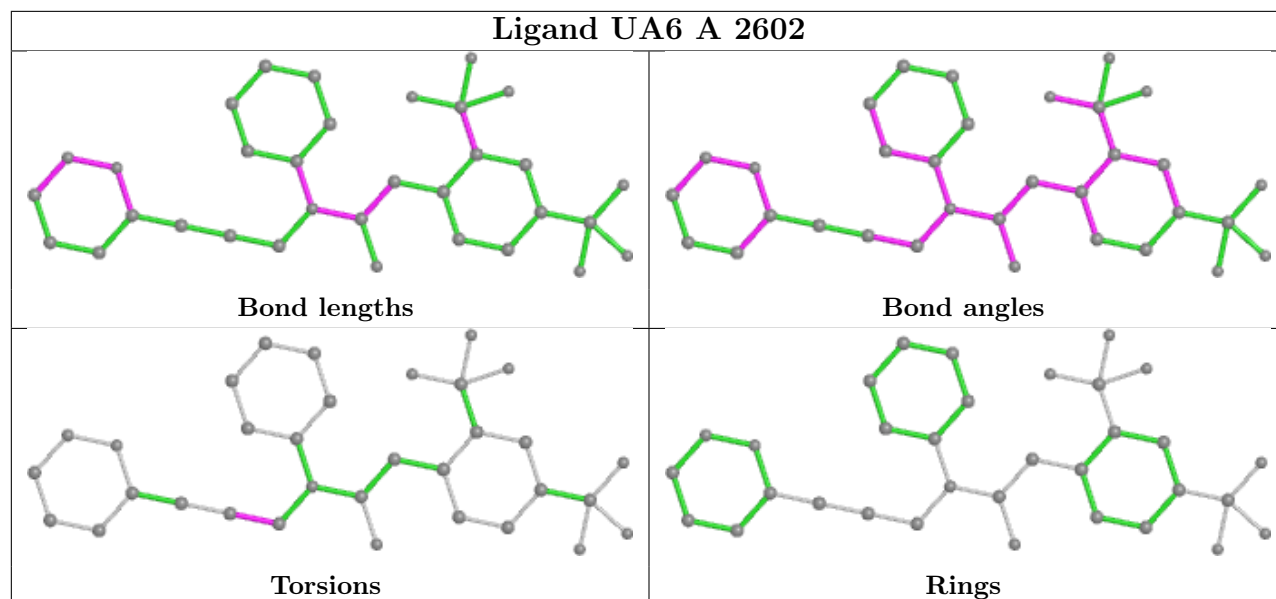
Mol	Chain	Res	Type	Atoms
6	A	2602	UA6	N03-C04-C05-C06
6	D	2602	UA6	C05-C04-N03-C13
5	D	2601	DG3	PG-O3B-PB-O1B
5	D	2601	DG3	PB-O3B-PG-O2G
5	A	2601	DG3	C5'-O5'-PA-O3A
5	D	2601	DG3	PG-O3B-PB-O2B
5	D	2601	DG3	PA-O3A-PB-O2B
5	A	2601	DG3	O4'-C4'-C5'-O5'
5	A	2601	DG3	PG-O3B-PB-O2B
5	A	2601	DG3	PB-O3A-PA-O2A
5	D	2601	DG3	C5'-O5'-PA-O3A
5	A	2601	DG3	PG-O3B-PB-O1B
5	D	2601	DG3	PA-O3A-PB-O1B
5	A	2601	DG3	C5'-O5'-PA-O1A

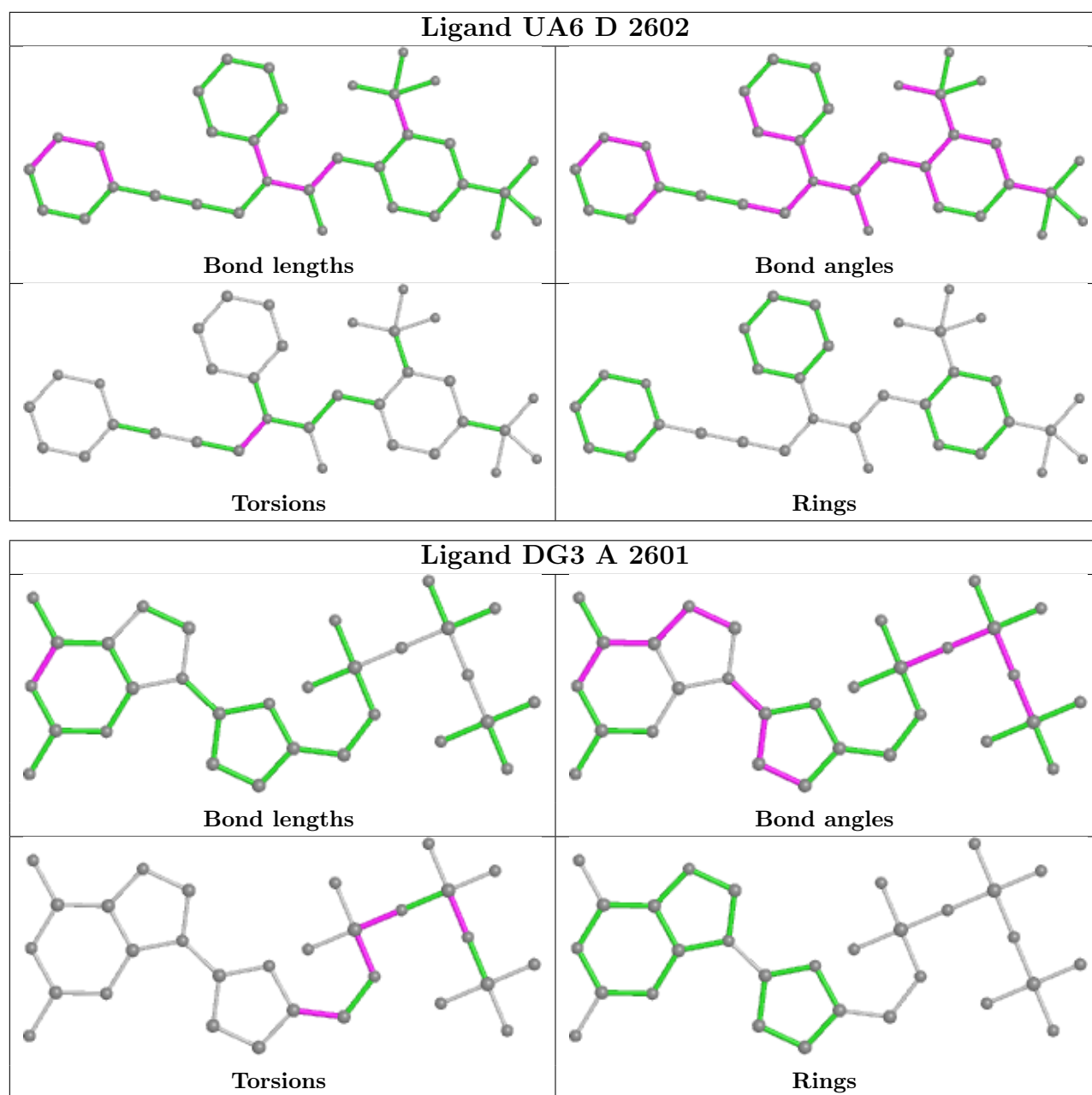
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	2601	DG3	1	0
5	A	2601	DG3	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	644/668 (96%)	0.25	12 (1%) 66 75	43, 67, 107, 138	0
1	D	648/668 (97%)	0.36	40 (6%) 20 28	55, 81, 122, 177	0
2	B	14/17 (82%)	-0.25	0 100 100	54, 76, 180, 221	0
2	E	13/17 (76%)	-0.45	0 100 100	64, 84, 121, 141	0
3	C	10/13 (76%)	-0.41	0 100 100	45, 101, 173, 197	0
3	F	9/13 (69%)	-0.38	0 100 100	66, 78, 139, 157	0
All	All	1338/1396 (95%)	0.28	52 (3%) 39 50	43, 75, 119, 221	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2495	PHE	4.9
1	D	2498	THR	4.3
1	D	2025	THR	4.3
1	D	2497	SER	4.0
1	A	2509	LEU	3.9
1	D	2030	GLN	3.5
1	D	2339	LEU	3.4
1	D	2469	ILE	3.4
1	D	2026	SER	3.1
1	D	2572	LEU	3.0
1	D	2027	GLN	3.0
1	D	2396	LYS	2.9
1	D	2190	LEU	2.9
1	D	2499	PHE	2.8
1	D	2028	GLY	2.8
1	D	2342	LEU	2.8
1	D	2426	PHE	2.8
1	D	2473	VAL	2.7
1	D	2133	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	1903	LEU	2.7
1	D	2348	LEU	2.7
1	A	2310	PHE	2.7
1	D	2494	THR	2.6
1	D	2351	VAL	2.6
1	A	2572	LEU	2.6
1	D	2354	THR	2.6
1	D	2585	LEU	2.5
1	D	2508	MET	2.5
1	D	2378	LEU	2.5
1	D	2249	ILE	2.5
1	D	2391	TYR	2.5
1	A	2191	HIS	2.4
1	D	2336	LEU	2.4
1	A	2212	PHE	2.4
1	D	1821	SER	2.4
1	D	2044	TYR	2.3
1	D	2444	LEU	2.3
1	D	2338	ILE	2.3
1	D	1937	PRO	2.2
1	D	2183	VAL	2.2
1	A	2193	LEU	2.2
1	D	2309	PRO	2.2
1	D	2423	ILE	2.2
1	D	2260	MET	2.2
1	A	2138	PHE	2.1
1	A	2198	LEU	2.1
1	A	2263	LEU	2.0
1	D	2139	LEU	2.0
1	D	2197	ILE	2.0
1	A	2346	ARG	2.0
1	D	2321	PHE	2.0
1	D	2443	ILE	2.0

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

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

6.3 Carbohydrates ⓘ

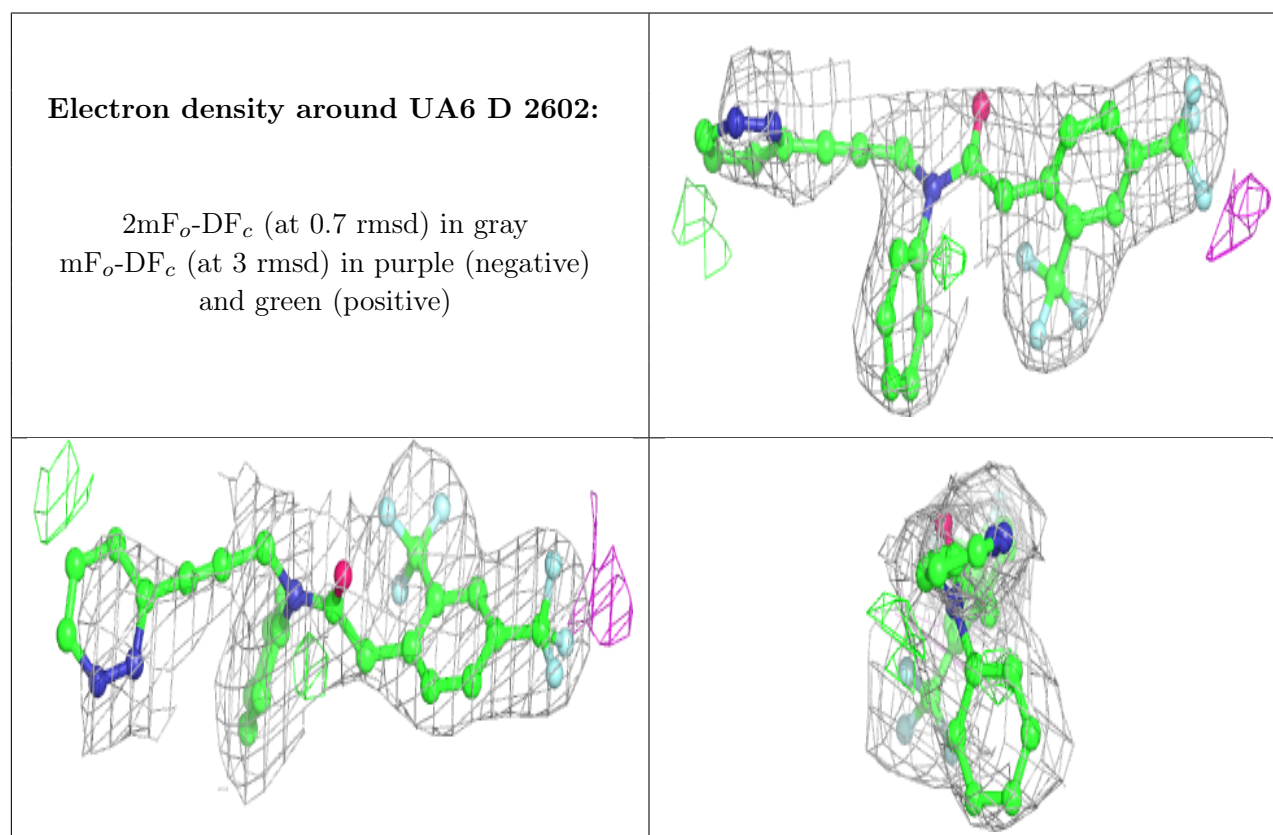
There are no monosaccharides in this entry.

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.

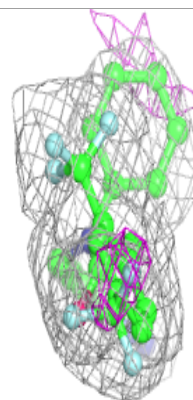
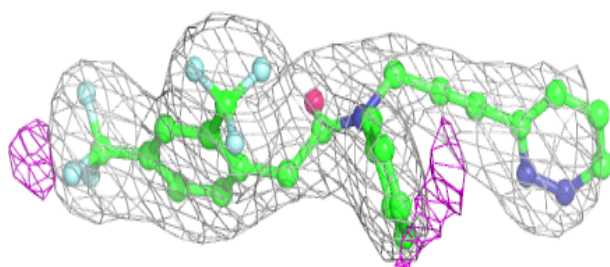
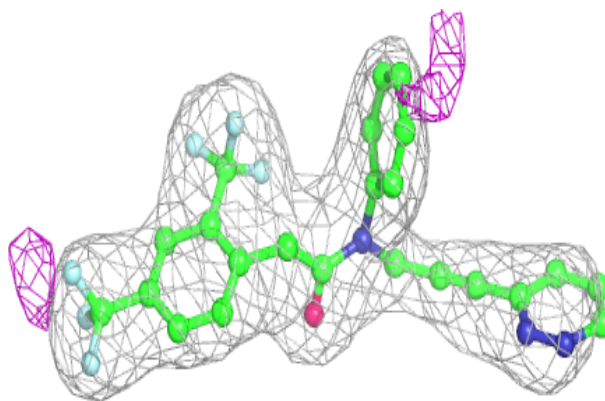
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	MG	D	2600	1/1	0.82	0.16	92,92,92,92	0
6	UA6	D	2602	33/33	0.90	0.20	91,104,135,141	0
6	UA6	A	2602	33/33	0.95	0.16	44,65,94,105	0
5	DG3	D	2601	30/30	0.97	0.14	60,75,83,89	0
5	DG3	A	2601	30/30	0.98	0.18	35,50,60,66	0
4	MG	A	2600	1/1	0.99	0.17	49,49,49,49	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.

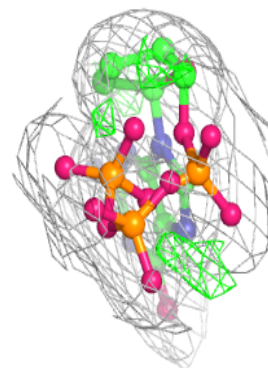
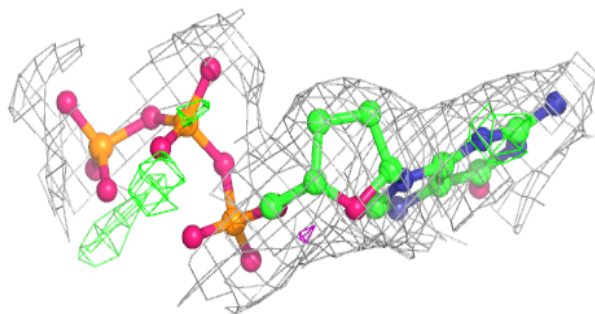
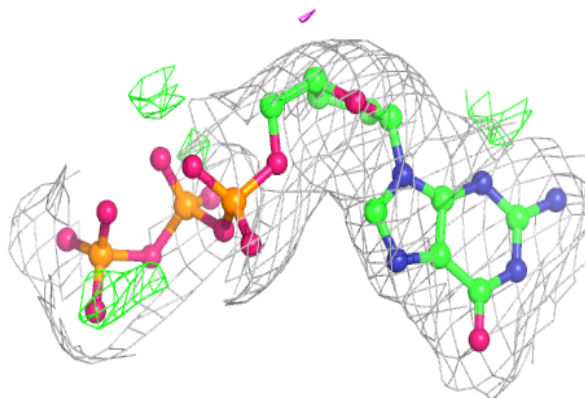


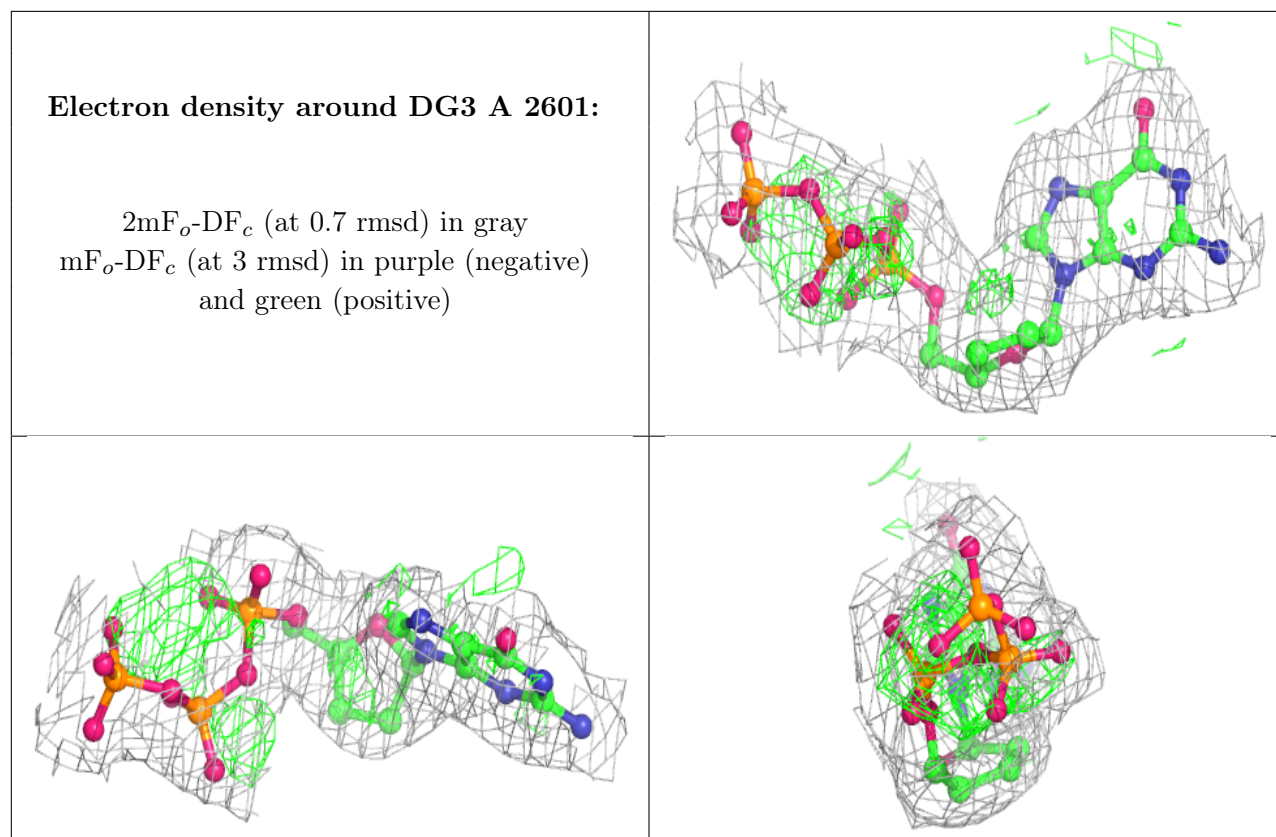
Electron density around UA6 A 2602:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around DG3 D 2601:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.