



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 06:29 PM UTC

PDB ID : 9V4M / pdb_00009v4m
Title : Crystal structure of the T.thermophilus transcription initiation complex bound to Ap4G
Authors : Duan, W.; Kaushik, A.; Serganov, A.
Deposited on : 2025-05-24
Resolution : 3.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

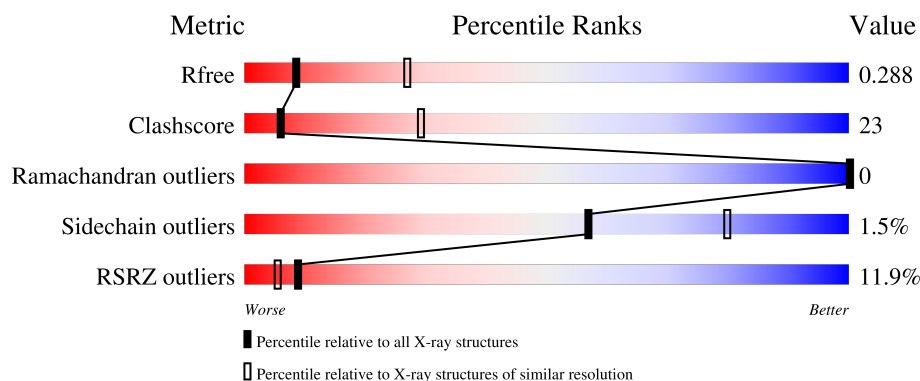
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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.



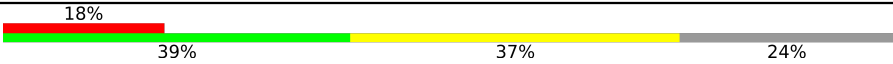
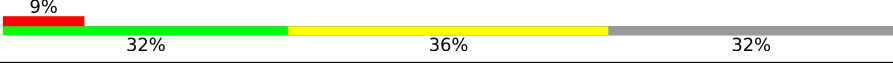
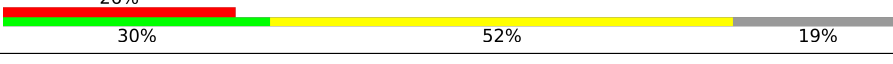
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	315	
1	B	315	
2	C	1119	
3	D	1524	
4	E	99	

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Mol	Chain	Length	Quality of chain
5	F	444	
6	G	22	
7	H	27	

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 27396 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-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	227	Total	C	N	O	S	0	1	0
			1788	1144	311	331	2			
1	B	226	Total	C	N	O	S	0	0	0
			1776	1135	307	332	2			

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	1113	Total	C	N	O	S	0	1	0
			8778	5555	1560	1639	24			

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	1357	Total	C	N	O	S	0	2	0
			10660	6738	1898	1990	34			

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	94	Total	C	N	O	S	0	0	0
			753	481	129	139	4			

- Molecule 5 is a protein called RNA polymerase sigma factor SigA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	338	Total	C	N	O	S	0	1	0
			2743	1729	498	512	4			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-20	MET	-	initiating methionine	UNP Q5SKW1
F	-19	GLY	-	expression tag	UNP Q5SKW1
F	-18	SER	-	expression tag	UNP Q5SKW1
F	-17	SER	-	expression tag	UNP Q5SKW1
F	-16	HIS	-	expression tag	UNP Q5SKW1
F	-15	HIS	-	expression tag	UNP Q5SKW1
F	-14	HIS	-	expression tag	UNP Q5SKW1
F	-13	HIS	-	expression tag	UNP Q5SKW1
F	-12	HIS	-	expression tag	UNP Q5SKW1
F	-11	HIS	-	expression tag	UNP Q5SKW1
F	-10	SER	-	expression tag	UNP Q5SKW1
F	-9	SER	-	expression tag	UNP Q5SKW1
F	-8	GLY	-	expression tag	UNP Q5SKW1
F	-7	LEU	-	expression tag	UNP Q5SKW1
F	-6	VAL	-	expression tag	UNP Q5SKW1
F	-5	PRO	-	expression tag	UNP Q5SKW1
F	-4	ARG	-	expression tag	UNP Q5SKW1
F	-3	GLY	-	expression tag	UNP Q5SKW1
F	-2	SER	-	expression tag	UNP Q5SKW1
F	-1	HIS	-	expression tag	UNP Q5SKW1
F	0	MET	-	expression tag	UNP Q5SKW1

- Molecule 6 is a DNA chain called Template DNA strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	G	15	Total	C	N	O	P	0	0	0
			301	141	57	88	15			

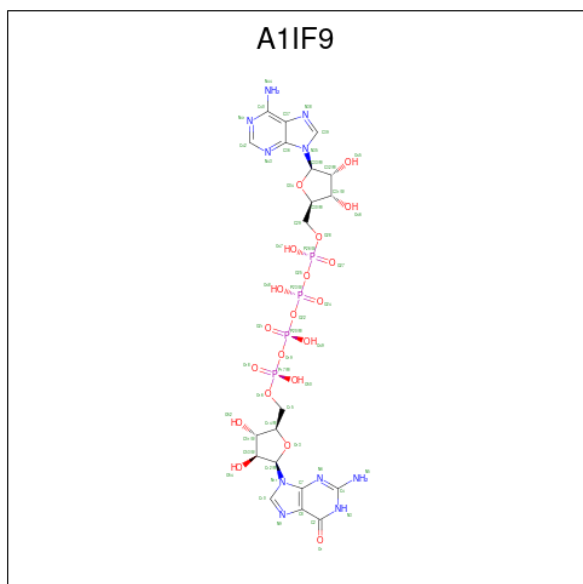
- Molecule 7 is a DNA chain called Non-template DNA Strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	H	22	Total	C	N	O	P	0	0	0
			454	217	86	130	21			

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

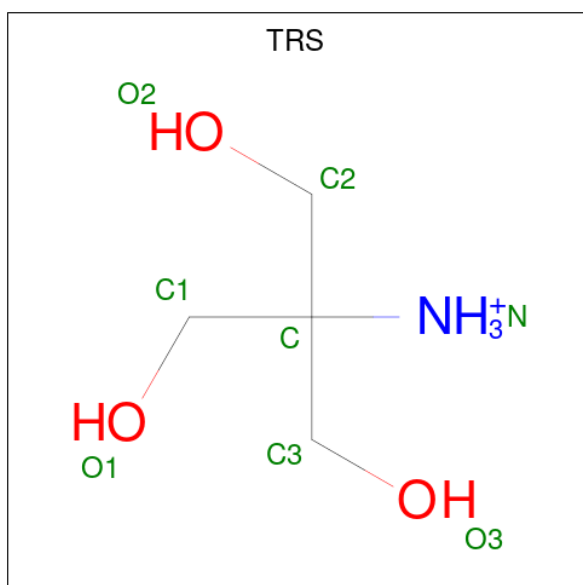
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Mg	0	0
			1	1		
8	B	1	Total	Mg	0	0
			1	1		
8	D	2	Total	Mg	0	0
			2	2		

- Molecule 9 is [[(2 {R},3 {S},4 {R},5 {R})-5-(6-aminopurin-9-yl)-3,4-bis(oxidanyl)oxolan-2-yl]methoxy-oxidanyl-phosphoryl] [[[(2 {R},3 {S},4 {R},5 {R})-5-(2-azanyl-6-oxidanylidene-1 {H}-purin-9-yl)-3,4-bis(oxidanyl)oxolan-2-yl]methoxy-oxidanyl-phosphoryl]oxy-oxidanyl-phosphoryl] hydrogen phosphate (CCD ID: A1IF9) (formula: C₂₀H₂₈N₁₀O₂₀P₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
9	C	1	54	20	10	20	4	0	0

- Molecule 10 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: C₄H₁₂NO₃).



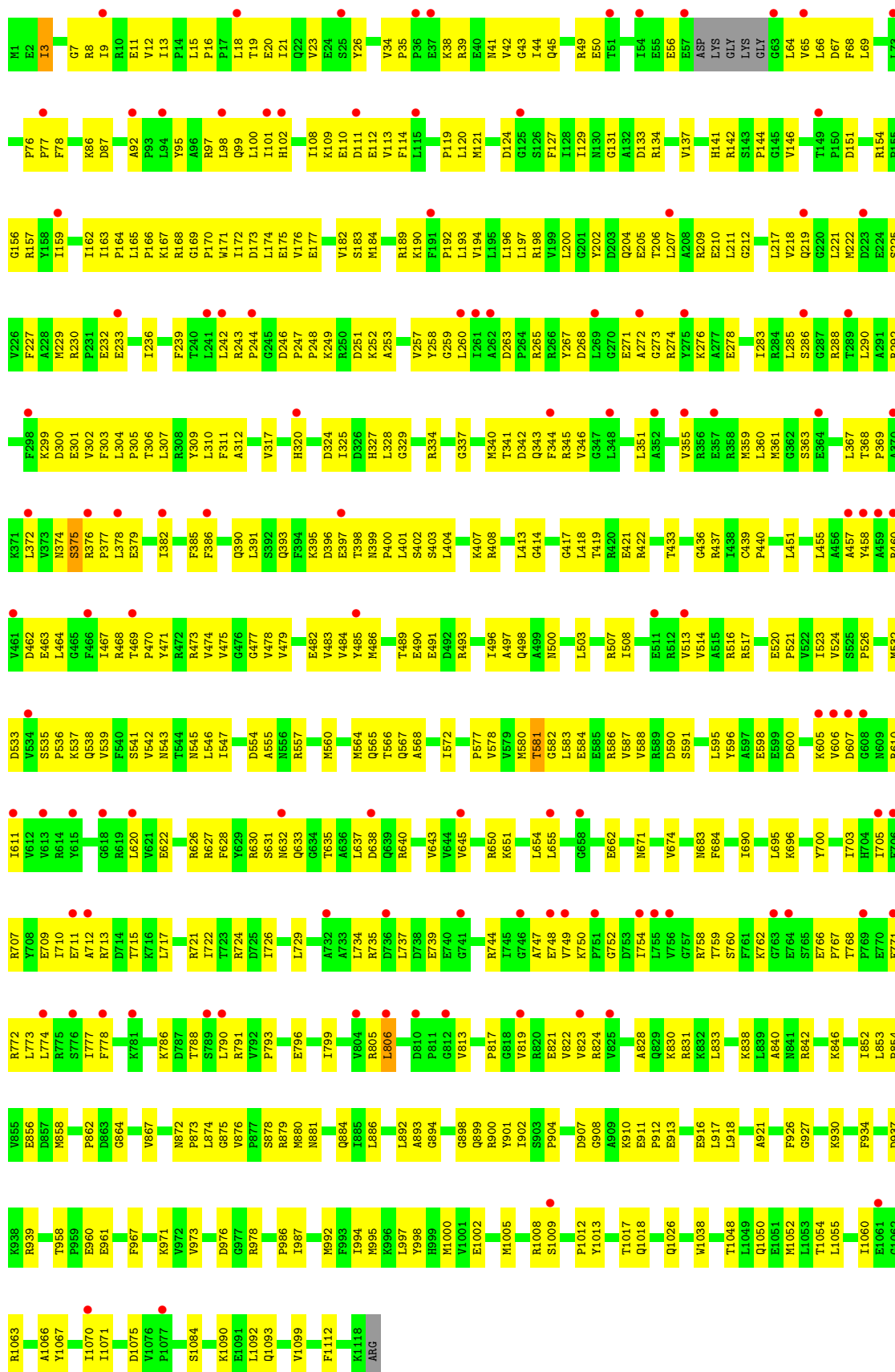
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	C	1	Total	C	N	O	0	0
			8	4	1	3		
10	D	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 11 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	D	2	Total	Zn	0	0
			2	2		

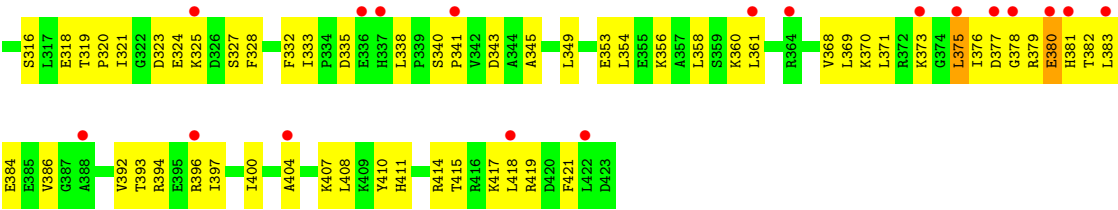
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	3	Total	O	0	0
			3	3		
12	B	4	Total	O	0	0
			4	4		
12	C	19	Total	O	0	0
			19	19		
12	D	27	Total	O	0	0
			27	27		
12	E	3	Total	O	0	0
			3	3		
12	F	6	Total	O	0	0
			6	6		
12	G	4	Total	O	0	0
			4	4		
12	H	1	Total	O	0	0
			1	1		

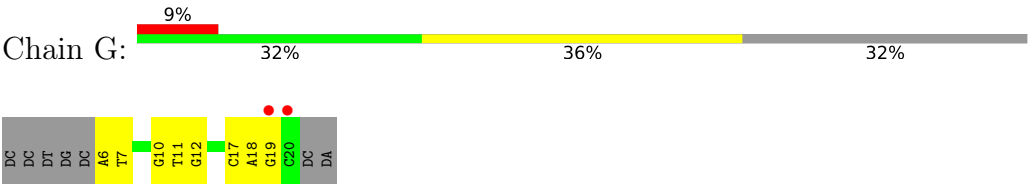


• Molecule 3: DNA-directed RNA polymerase subunit beta'

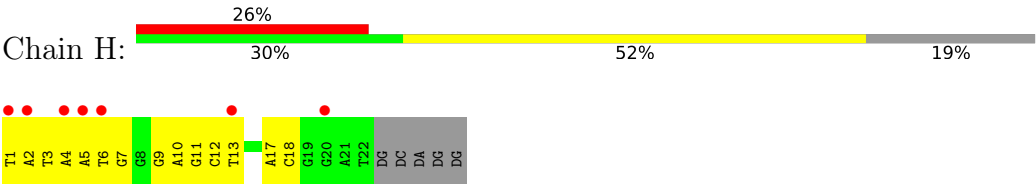




● Molecule 6: Template DNA strand



● Molecule 7: Non-template DNA Strand



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	185.06Å 103.21Å 296.51Å 90.00° 98.89° 90.00°	Depositor
Resolution (Å)	29.96 – 3.00 29.96 – 3.00	Depositor EDS
% Data completeness (in resolution range)	78.4 (29.96-3.00) 78.4 (29.96-3.00)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 3.00Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.279 , 0.288 0.279 , 0.288	Depositor DCC
R_{free} test set	2024 reflections (1.82%)	wwPDB-VP
Wilson B-factor (Å ²)	59.9	Xtriage
Anisotropy	0.217	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 67.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	27396	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.03% 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: A1IF9, ZN, TRS, 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	0.47	0/1823	0.72	0/2479
1	B	0.39	0/1808	0.68	0/2459
2	C	0.28	0/8948	0.50	0/12103
3	D	0.28	0/10848	0.48	0/14660
4	E	0.29	0/767	0.55	0/1035
5	F	0.37	0/2788	0.63	0/3749
6	G	0.45	0/337	0.69	0/518
7	H	0.46	0/510	0.72	0/787
All	All	0.32	0/27829	0.55	0/37790

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	1788	0	1846	119	0
1	B	1776	0	1823	85	0
2	C	8778	0	8868	468	0
3	D	10660	0	10807	460	0
4	E	753	0	760	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	F	2743	0	2812	157	0
6	G	301	0	164	8	0
7	H	454	0	250	23	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
8	D	2	0	0	0	0
9	C	54	0	0	2	0
10	C	8	0	12	0	0
10	D	8	0	12	1	0
11	D	2	0	0	0	0
12	A	3	0	0	0	0
12	B	4	0	0	0	0
12	C	19	0	0	0	0
12	D	27	0	0	0	0
12	E	3	0	0	1	0
12	F	6	0	0	0	0
12	G	4	0	0	0	0
12	H	1	0	0	0	0
All	All	27396	0	27354	1246	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 23.

The worst 5 of 1246 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:53:VAL:HG11	1:A:82:LEU:CD2	1.77	1.15
1:A:53:VAL:HG11	1:A:82:LEU:HD21	1.29	1.14
2:C:109:LYS:HG2	2:C:368:THR:HG22	1.46	0.97
3:D:45:PHE:HZ	3:D:541:ASN:HD21	1.13	0.94
2:C:458:TYR:HE2	2:C:535:SER:HG	1.15	0.93

There are no symmetry-related clashes.

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	226/315 (72%)	202 (89%)	24 (11%)	0	100	100
1	B	224/315 (71%)	216 (96%)	8 (4%)	0	100	100
2	C	1110/1119 (99%)	1037 (93%)	73 (7%)	0	100	100
3	D	1353/1524 (89%)	1287 (95%)	66 (5%)	0	100	100
4	E	92/99 (93%)	86 (94%)	6 (6%)	0	100	100
5	F	335/444 (76%)	318 (95%)	17 (5%)	0	100	100
All	All	3340/3816 (88%)	3146 (94%)	194 (6%)	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	199/273 (73%)	193 (97%)	6 (3%)	36	69
1	B	198/273 (72%)	196 (99%)	2 (1%)	68	84
2	C	935/941 (99%)	928 (99%)	7 (1%)	76	86
3	D	1128/1279 (88%)	1111 (98%)	17 (2%)	57	80
4	E	81/88 (92%)	81 (100%)	0	100	100
5	F	293/389 (75%)	283 (97%)	10 (3%)	32	66
All	All	2834/3243 (87%)	2792 (98%)	42 (2%)	57	80

5 of 42 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	D	1282	ARG
5	F	333	ILE
3	D	1304	LYS
5	F	115	LYS
5	F	338	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
5	F	214	GLN
5	F	280	GLN
2	C	962	GLN
2	C	884	GLN
5	F	337	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 6 are monoatomic - leaving 3 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	$\# Z > 2$
9	A1IF9	C	1201	8	59,59,59	0.84	2 (3%)	86,93,93	1.19	8 (9%)
10	TRS	C	1202	-	7,7,7	0.15	0	9,9,9	0.24	0
10	TRS	D	1605	-	7,7,7	0.14	0	9,9,9	0.30	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
9	A1IF9	C	1201	8	-	10/38/70/70	0/6/6/6
10	TRS	C	1202	-	-	0/9/9/9	-
10	TRS	D	1605	-	-	0/9/9/9	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	C	1201	A1IF9	C36-N35	-3.43	1.30	1.37
9	C	1201	A1IF9	O46-C31	2.08	1.48	1.43

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	C	1201	A1IF9	C37-C36-N43	-4.97	119.88	126.72
9	C	1201	A1IF9	N35-C39-N38	-4.27	107.88	113.94
9	C	1201	A1IF9	C36-N35-C39	3.16	109.06	105.74
9	C	1201	A1IF9	N43-C36-N35	2.69	131.74	127.17
9	C	1201	A1IF9	C42-N43-C36	2.68	118.37	111.83

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	C	1201	A1IF9	C29-O28-P26-O25
9	C	1201	A1IF9	C29-O28-P26-O27
9	C	1201	A1IF9	C29-O28-P26-O47
9	C	1201	A1IF9	C30-C29-O28-P26
9	C	1201	A1IF9	P20-O22-P23-O24

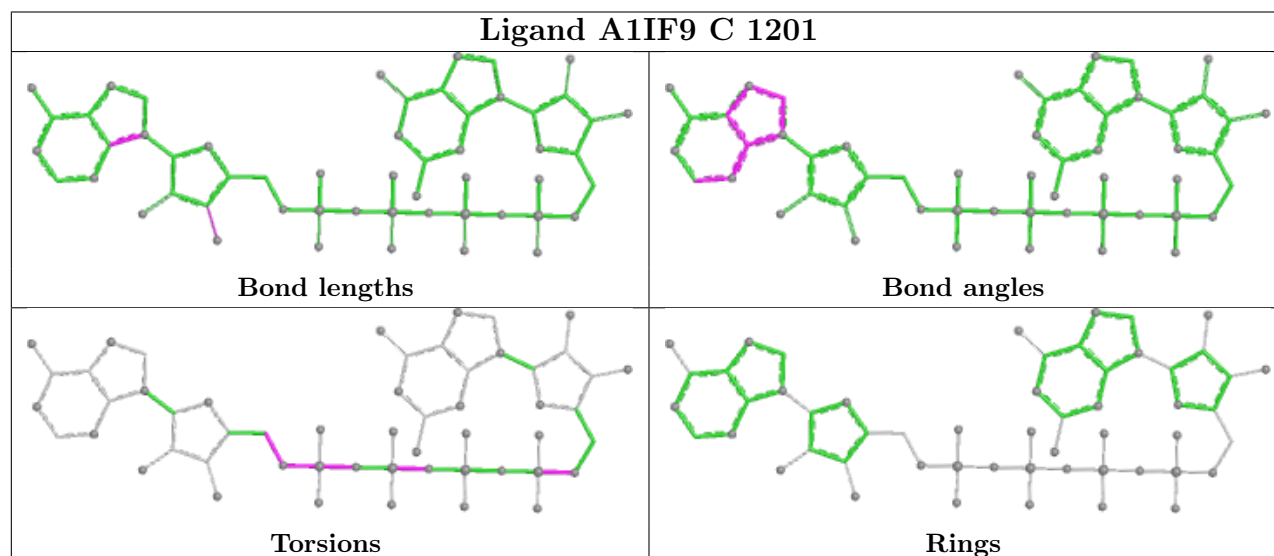
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	C	1201	A1IF9	2	0
10	D	1605	TRS	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 [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, 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	227/315 (72%)	0.84	23 (10%) 12 7	37, 71, 91, 105	1 (0%)
1	B	226/315 (71%)	0.26	0 100 100	33, 54, 79, 93	0
2	C	1113/1119 (99%)	0.86	113 (10%) 12 6	20, 75, 119, 146	1 (0%)
3	D	1357/1524 (89%)	0.69	168 (12%) 8 5	19, 56, 174, 222	3 (0%)
4	E	94/99 (94%)	0.94	9 (9%) 13 7	39, 75, 104, 109	0
5	F	338/444 (76%)	1.49	80 (23%) 2 1	43, 89, 114, 143	1 (0%)
6	G	15/22 (68%)	0.95	2 (13%) 7 4	43, 69, 127, 130	0
7	H	22/27 (81%)	1.83	7 (31%) 1 1	88, 109, 144, 157	0
All	All	3392/3865 (87%)	0.82	402 (11%) 9 5	19, 70, 136, 222	6 (0%)

The worst 5 of 402 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	216	VAL	6.6
3	D	409	VAL	5.4
3	D	365	ASP	5.3
2	C	63	GLY	5.0
3	D	207	PHE	4.9

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](#)

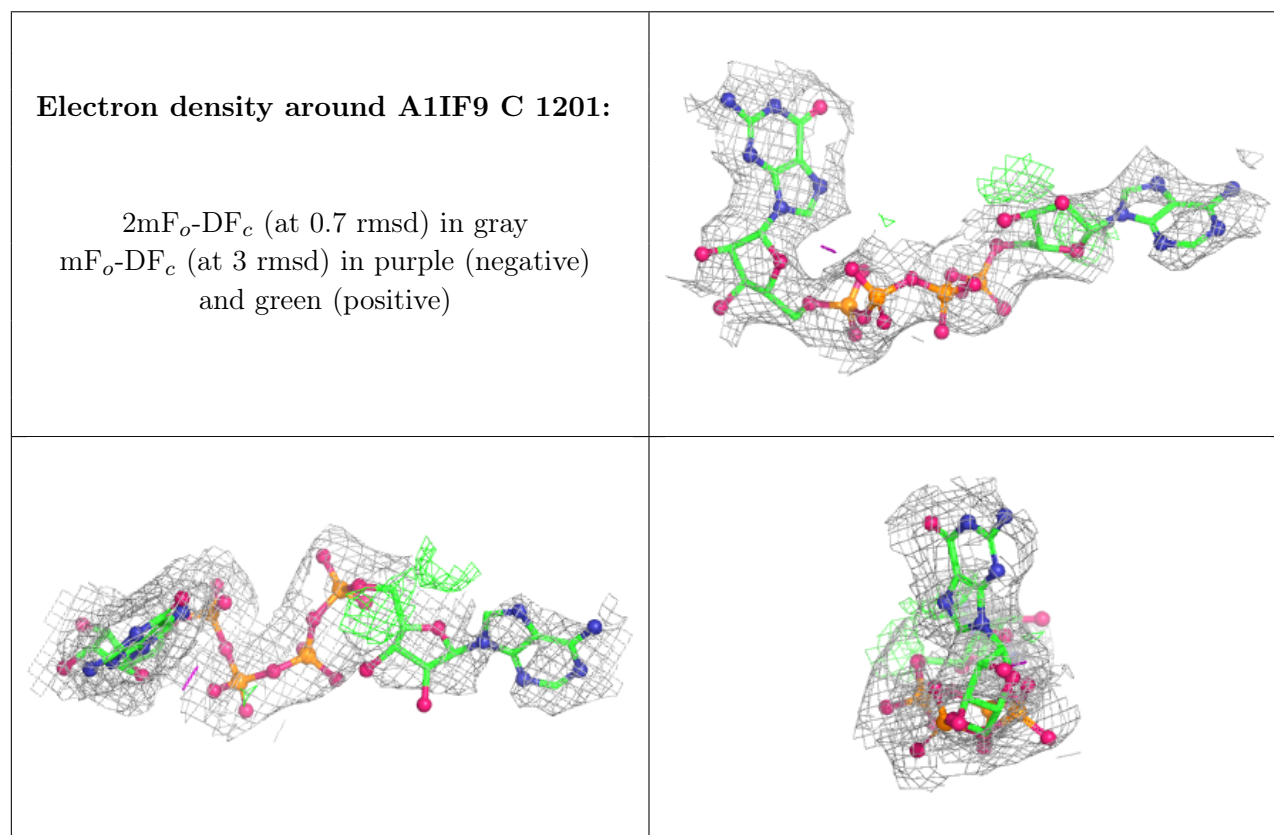
There are no oligosaccharides in this entry.

6.4 Ligands [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, 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
10	TRS	D	1605	8/8	0.61	0.21	31,37,45,47	0
10	TRS	C	1202	8/8	0.64	0.21	41,51,56,56	0
11	ZN	D	1602	1/1	0.80	0.12	174,174,174,174	0
8	MG	A	401	1/1	0.83	0.24	60,60,60,60	0
9	A1IF9	C	1201	54/54	0.89	0.11	37,47,66,68	21
8	MG	B	401	1/1	0.90	0.20	32,32,32,32	0
8	MG	D	1604	1/1	0.93	0.08	36,36,36,36	0
11	ZN	D	1601	1/1	0.97	0.04	36,36,36,36	0
8	MG	D	1603	1/1	0.98	0.03	25,25,25,25	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.