



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 6, 2022 – 01:09 PM EDT

PDB ID : 5X21
Title : Crystal structure of Thermus thermophilus transcription initiation complex with GpA and pseudouridimycin (PUM)
Authors : Zhang, Y.; Ebright, R.
Deposited on : 2017-01-29
Resolution : 3.32 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.31.2
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) : 2.31.2

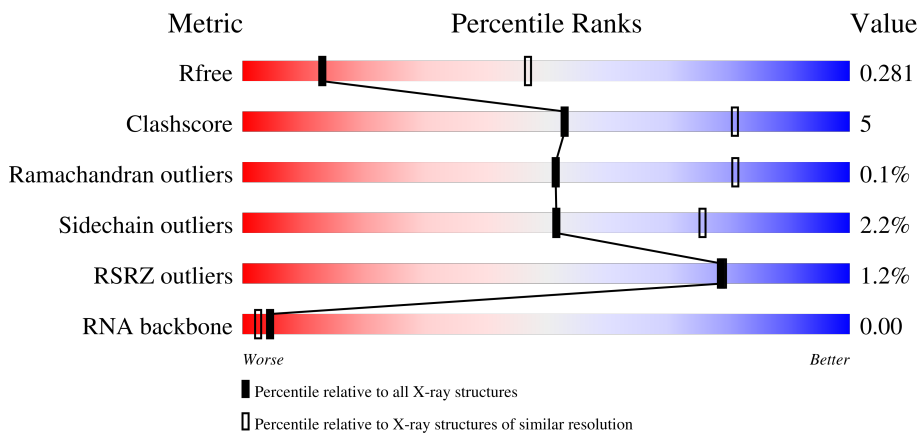
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.32 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.






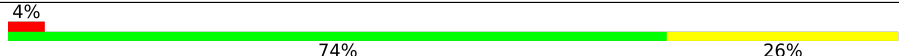
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1089 (3.36-3.28)
Clashscore	141614	1137 (3.36-3.28)
Ramachandran outliers	138981	1115 (3.36-3.28)
Sidechain outliers	138945	1114 (3.36-3.28)
RSRZ outliers	127900	1059 (3.36-3.28)
RNA backbone	3102	1125 (3.74-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	 57% 14% 28%
1	B	315	 61% 9% 30%
2	C	1119	 82% 17% ..
3	D	1524	 81% 16% ..

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Mol	Chain	Length	Quality of chain
4	E	99	 83% 12% 5%
5	F	443	 3% 71% 7% 22%
6	G	21	 10% 76% 14% 10%
7	H	27	 4% 74% 26%
8	I	2	 100%

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 28731 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
			Total	C	N	O	S			
1	A	226	Total	C	N	O	S	0	0	0
			1782	1138	310	332	2			
1	B	222	Total	C	N	O	S	0	0	0
			1750	1118	304	326	2			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	1111	Total	C	N	O	S	0	0	0
			8770	5548	1564	1634	24			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	D	1496	Total	C	N	O	S	0	0	0
			11800	7479	2078	2207	36			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	E	94	Total	C	N	O	S	0	0	0
			758	483	132	139	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	F	346	Total	C	N	O	S	0	0	0
			2807	1770	509	524	4			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-19	MET	-	initiating methionine	UNP Q72L95
F	-18	GLY	-	expression tag	UNP Q72L95
F	-17	SER	-	expression tag	UNP Q72L95
F	-16	SER	-	expression tag	UNP Q72L95
F	-15	HIS	-	expression tag	UNP Q72L95
F	-14	HIS	-	expression tag	UNP Q72L95
F	-13	HIS	-	expression tag	UNP Q72L95
F	-12	HIS	-	expression tag	UNP Q72L95
F	-11	HIS	-	expression tag	UNP Q72L95
F	-10	HIS	-	expression tag	UNP Q72L95
F	-9	SER	-	expression tag	UNP Q72L95
F	-8	SER	-	expression tag	UNP Q72L95
F	-7	GLY	-	expression tag	UNP Q72L95
F	-6	LEU	-	expression tag	UNP Q72L95
F	-5	VAL	-	expression tag	UNP Q72L95
F	-4	PRO	-	expression tag	UNP Q72L95
F	-3	ARG	-	expression tag	UNP Q72L95
F	-2	GLY	-	expression tag	UNP Q72L95
F	-1	SER	-	expression tag	UNP Q72L95
F	0	HIS	-	expression tag	UNP Q72L95

- Molecule 6 is a DNA chain called promoter DNA template strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
6	G	19	385	184	71	112	18	0	0	0

- Molecule 7 is a DNA chain called promoter DNA nontemplate strand.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
7	H	27	560	266	109	159	26	0	0	0

- Molecule 8 is a RNA chain called RNA (5'-R(*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
8	I	2	42	20	10	11	1	0	0	0

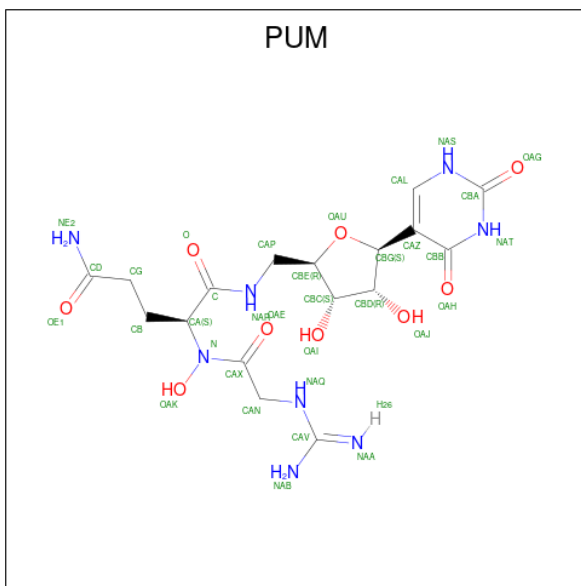
- Molecule 9 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total Mg 1 1	0	0
9	C	1	Total Mg 1 1	0	0
9	D	2	Total Mg 2 2	0	0
9	F	1	Total Mg 1 1	0	0
9	H	1	Total Mg 1 1	0	0

- Molecule 10 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 11 is (1S)-1,4-anhydro-5-[(N-carbamimidoylglycyl-N 2 -hydroxy-L-glutaminy]a mino]-5-deoxy-1-(2,4-dioxo-1,2,3,4-tetrahydropyrimidin-5-yl)-D-ribitol (three-letter code: PUM) (formula: C₁₇H₂₆N₈O₉).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	I	1	Total C N O 34 17 8 9	0	0

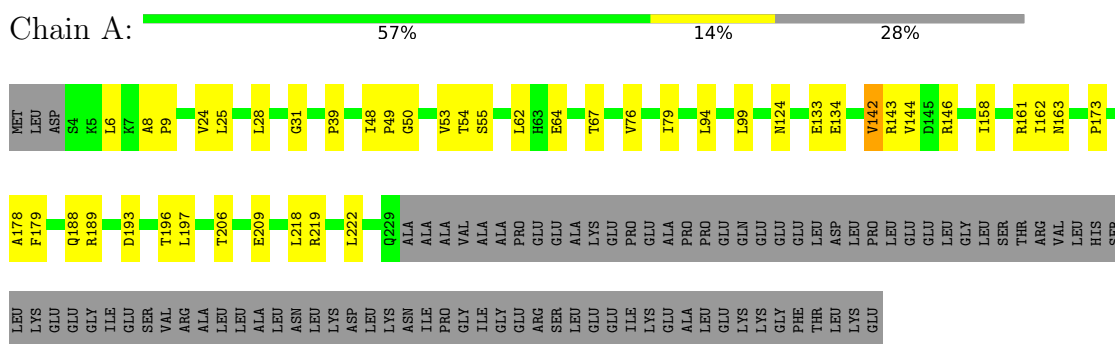
- Molecule 12 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
12	A	4	Total O 4 4	0	0
12	B	3	Total O 3 3	0	0
12	C	10	Total O 10 10	0	0
12	D	11	Total O 11 11	0	0
12	E	3	Total O 3 3	0	0
12	F	2	Total O 2 2	0	0
12	H	2	Total O 2 2	0	0

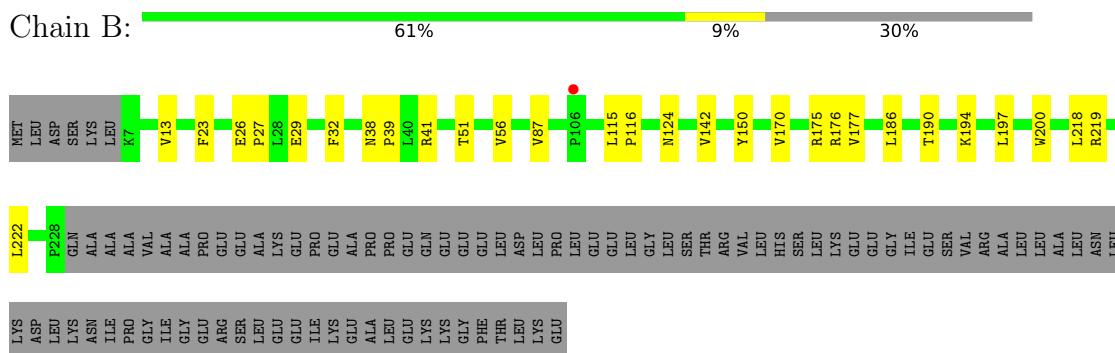
3 Residue-property plots [i](#)

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.

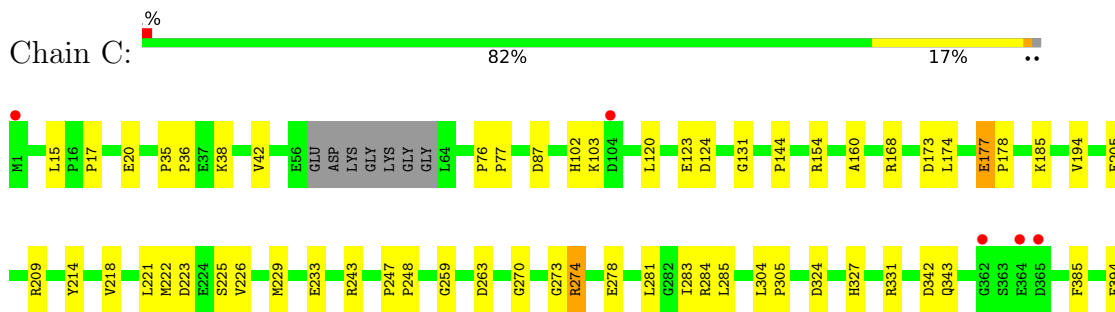
- Molecule 1: DNA-directed RNA polymerase subunit alpha

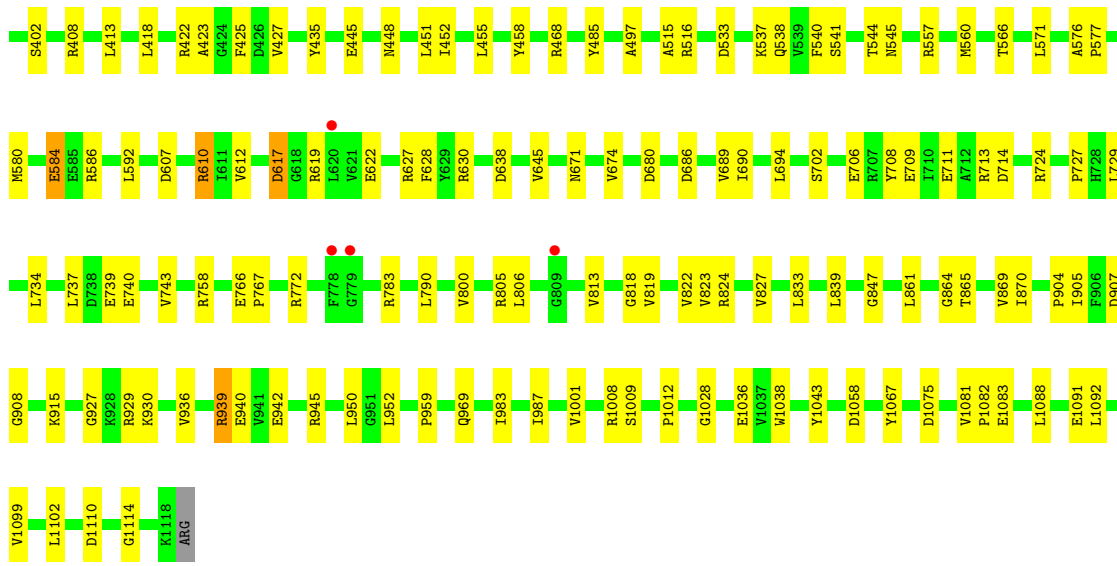


- Molecule 1: DNA-directed RNA polymerase subunit alpha

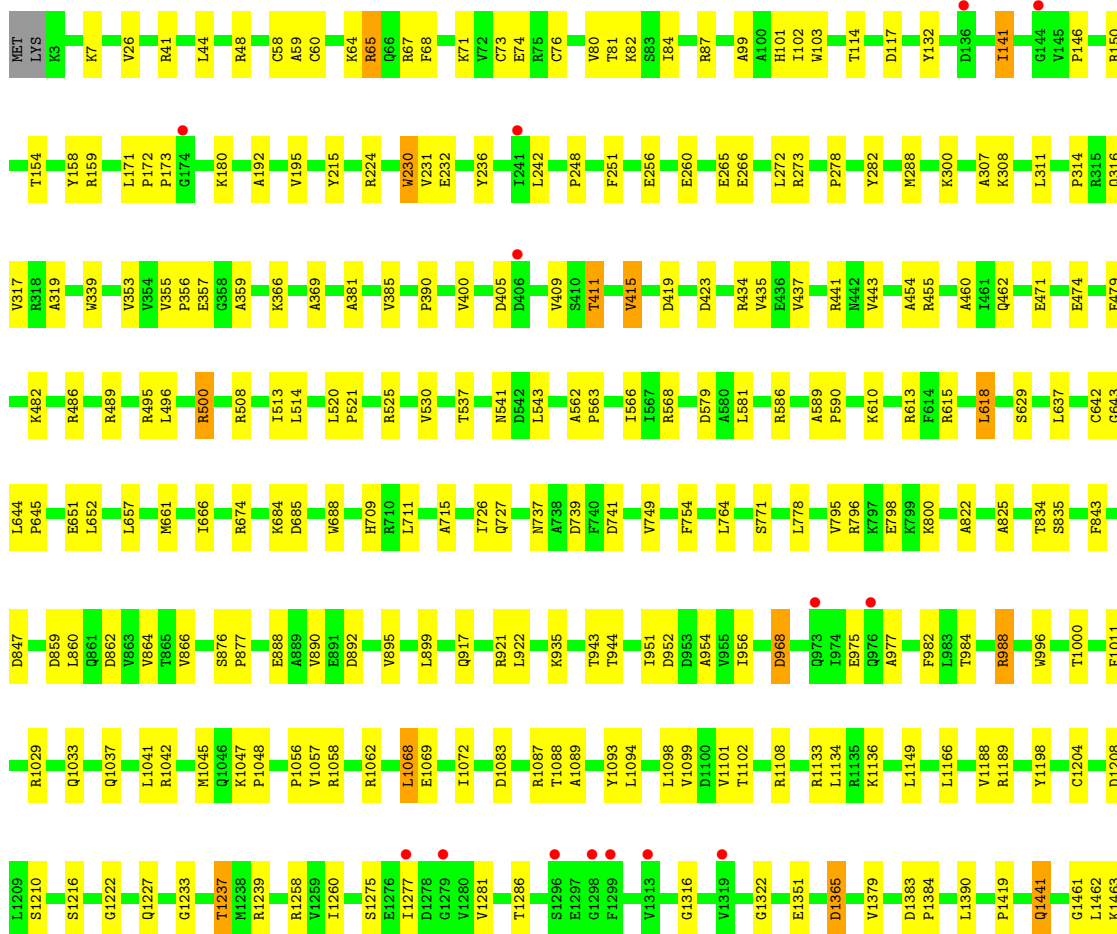
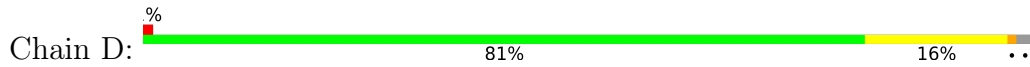


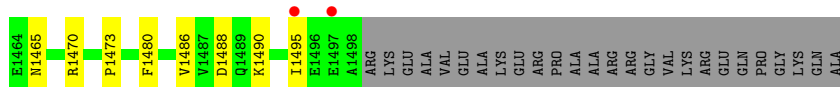
- Molecule 2: DNA-directed RNA polymerase subunit beta





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





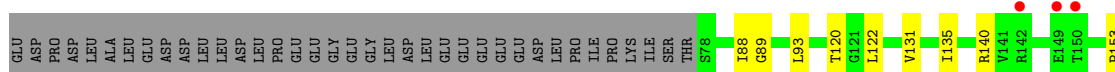
- Molecule 4: DNA-directed RNA polymerase subunit omega

Chain E: 83% 12% 5%



- Molecule 5: RNA polymerase sigma factor SigA

Chain F: 3% 71% 7% 22%



- Molecule 6: promoter DNA template strand

Chain G: 10% 76% 14% 10%



- Molecule 7: promoter DNA nontemplate strand

Chain H: 4% 74% 26%



- Molecule 8: RNA (5'-R(*GP*A)-3')

Chain I: 100%

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	186.79Å 103.09Å 296.23Å 90.00° 98.71° 90.00°	Depositor
Resolution (Å)	39.67 – 3.32 39.67 – 3.32	Depositor EDS
% Data completeness (in resolution range)	85.1 (39.67-3.32) 85.1 (39.67-3.32)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.10 (at 3.32Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.232 , 0.280 0.234 , 0.281	Depositor DCC
R_{free} test set	1565 reflections (2.23%)	wwPDB-VP
Wilson B-factor (Å ²)	60.4	Xtrriage
Anisotropy	0.035	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 20.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	28731	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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: MG, PUM, ZN

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	0.21	0/1814	0.40	0/2466
1	B	0.20	0/1782	0.40	0/2424
2	C	0.20	0/8937	0.38	0/12087
3	D	0.20	0/12009	0.38	0/16240
4	E	0.20	0/772	0.36	0/1040
5	F	0.20	0/2852	0.36	0/3837
6	G	0.51	0/431	0.86	0/663
7	H	0.50	0/630	0.90	0/973
8	I	0.13	0/47	0.52	0/72
All	All	0.22	0/29274	0.41	0/39802

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	1782	0	1834	27	0
1	B	1750	0	1797	17	0
2	C	8770	0	8874	109	0
3	D	11800	0	12024	157	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	758	0	770	8	0
5	F	2807	0	2882	20	0
6	G	385	0	215	2	0
7	H	560	0	305	7	0
8	I	42	0	23	0	0
9	A	1	0	0	0	0
9	C	1	0	0	0	0
9	D	2	0	0	0	0
9	F	1	0	0	0	0
9	H	1	0	0	0	0
10	D	2	0	0	0	0
11	I	34	0	0	1	0
12	A	4	0	0	1	0
12	B	3	0	0	0	0
12	C	10	0	0	0	0
12	D	11	0	0	0	0
12	E	3	0	0	0	0
12	F	2	0	0	0	0
12	H	2	0	0	0	0
All	All	28731	0	28724	311	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 5.

The worst 5 of 311 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)
2:C:628:PHE:H	2:C:638:ASP:HB3	1.47	0.77
2:C:194:VAL:HG22	2:C:221:LEU:HD12	1.70	0.74
3:D:60:CYS:SG	3:D:76:CYS:HB3	2.28	0.74
3:D:117:ASP:HB3	3:D:150:ARG:HH22	1.53	0.71
2:C:422:ARG:HH22	7:H:13:DT:H5'	1.56	0.69

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	224/315 (71%)	218 (97%)	5 (2%)	1 (0%)	34	66
1	B	220/315 (70%)	215 (98%)	5 (2%)	0	100	100
2	C	1107/1119 (99%)	1069 (97%)	37 (3%)	1 (0%)	51	81
3	D	1494/1524 (98%)	1455 (97%)	38 (2%)	1 (0%)	51	81
4	E	92/99 (93%)	91 (99%)	1 (1%)	0	100	100
5	F	344/443 (78%)	340 (99%)	4 (1%)	0	100	100
All	All	3481/3815 (91%)	3388 (97%)	90 (3%)	3 (0%)	51	81

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	162	ILE
2	C	423	ALA
3	D	530	VAL

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	199/273 (73%)	196 (98%)	3 (2%)	65	81
1	B	195/273 (71%)	193 (99%)	2 (1%)	76	86
2	C	936/941 (100%)	915 (98%)	21 (2%)	52	76
3	D	1258/1279 (98%)	1227 (98%)	31 (2%)	47	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
4	E	82/88 (93%)	80 (98%)	2 (2%)	49 74
5	F	301/387 (78%)	296 (98%)	5 (2%)	60 79
All	All	2971/3241 (92%)	2907 (98%)	64 (2%)	52 76

5 of 64 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	E	31	LEU
5	F	88	ILE
2	C	1001	VAL
2	C	939	ARG
5	F	172	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	163	ASN
1	B	128	HIS
3	D	794	GLN
3	D	1195	GLN
3	D	1353	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
8	I	1/2 (50%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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

Of 9 ligands modelled in this entry, 8 are monoatomic - leaving 1 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
11	PUM	I	101	-	34,35,35	2.32	8 (23%)	39,49,49	1.63	6 (15%)

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
11	PUM	I	101	-	-	6/31/47/47	0/2/2/2

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	I	101	PUM	OAK-N	-7.05	1.31	1.40
11	I	101	PUM	CAL-NAS	5.80	1.45	1.36
11	I	101	PUM	CBA-NAS	5.60	1.44	1.36
11	I	101	PUM	CBD-CBG	-4.09	1.48	1.53
11	I	101	PUM	CBA-NAT	3.22	1.43	1.37

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	I	101	PUM	CBB-NAT-CBA	-4.24	120.23	126.34
11	I	101	PUM	NAS-CBA-NAT	4.09	119.76	115.13
11	I	101	PUM	OAU-CBG-CBD	3.87	110.60	105.14
11	I	101	PUM	CAL-NAS-CBA	-2.57	120.06	122.68
11	I	101	PUM	OAG-CBA-NAS	-2.47	120.07	122.79

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

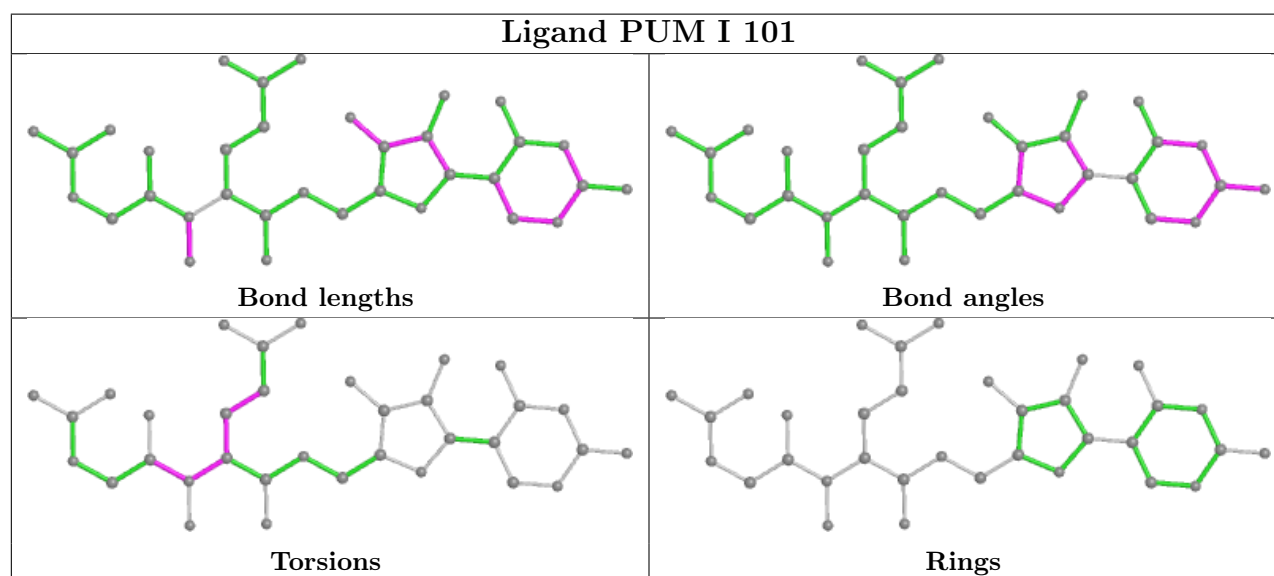
Mol	Chain	Res	Type	Atoms
11	I	101	PUM	C-CA-CB-CG
11	I	101	PUM	N-CA-CB-CG
11	I	101	PUM	C-CA-N-OAK
11	I	101	PUM	CB-CA-N-OAK
11	I	101	PUM	CA-CB-CG-CD

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	I	101	PUM	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

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	226/315 (71%)	-0.19	0 100 100	34, 60, 83, 99	0
1	B	222/315 (70%)	-0.07	1 (0%) 91 91	33, 70, 100, 122	0
2	C	1111/1119 (99%)	-0.17	9 (0%) 86 87	15, 55, 109, 132	0
3	D	1496/1524 (98%)	-0.10	16 (1%) 80 81	14, 52, 110, 144	1 (0%)
4	E	94/99 (94%)	-0.01	0 100 100	35, 68, 101, 107	0
5	F	346/443 (78%)	0.04	12 (3%) 44 42	22, 60, 126, 140	0
6	G	19/21 (90%)	0.16	2 (10%) 6 6	20, 48, 153, 153	0
7	H	27/27 (100%)	0.05	1 (3%) 41 40	49, 67, 141, 158	0
8	I	2/2 (100%)	-0.51	0 100 100	28, 28, 28, 30	0
All	All	3543/3865 (91%)	-0.11	41 (1%) 79 79	14, 57, 112, 158	1 (0%)

The worst 5 of 41 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	365	ASP	4.8
5	F	149	GLU	4.3
5	F	388	ALA	4.3
5	F	415	THR	4.2
2	C	362	GLY	3.8

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](#)

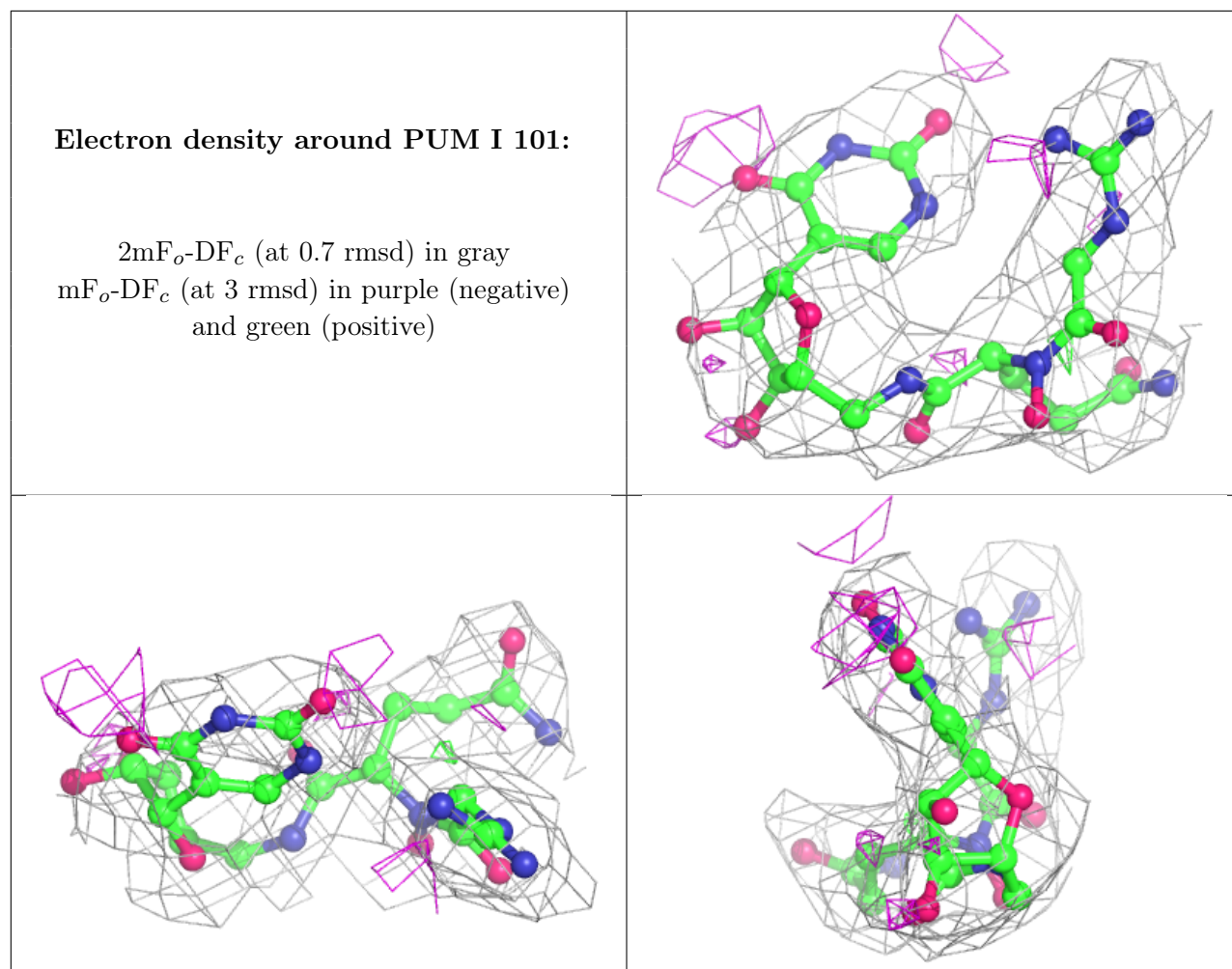
There are no monosaccharides in this entry.

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
9	MG	A	2001	1/1	0.65	0.25	43,43,43,43	0
9	MG	H	2001	1/1	0.82	0.20	50,50,50,50	0
9	MG	F	2001	1/1	0.84	0.08	45,45,45,45	0
9	MG	C	2000	1/1	0.86	0.15	31,31,31,31	0
9	MG	D	1604	1/1	0.86	0.10	21,21,21,21	0
11	PUM	I	101	34/34	0.94	0.20	20,27,36,40	0
9	MG	D	1601	1/1	0.97	0.41	39,39,39,39	0
10	ZN	D	1602	1/1	0.98	0.09	36,36,36,36	0
10	ZN	D	1603	1/1	0.99	0.06	80,80,80,80	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.