



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 10:26 AM EDT

PDB ID : 4Y4M / pdb_00004y4m
Title : Thiazole synthase Thi4 from Methanocaldococcus jannaschii
Authors : Zhang, X.; Ealick, S.E.
Deposited on : 2015-02-10
Resolution : 2.71 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

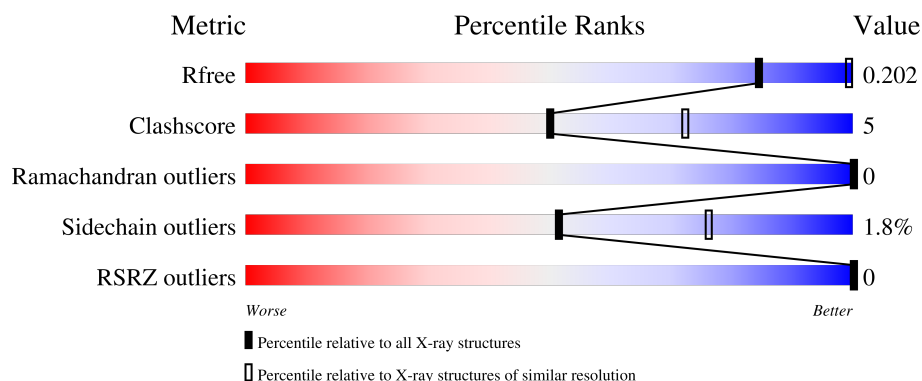
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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	290	
1	B	290	
1	C	290	
1	D	290	
1	E	290	

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Mol	Chain	Length	Quality of chain
1	F	290	 83%7%10%
1	G	290	 83%7%10%
1	H	290	 83%7%11%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15644 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 Putative ribose 1,5-bisphosphate isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	260	Total	C	N	O	S	0	0	0
			1924	1230	322	360	12			
1	B	256	Total	C	N	O	S	0	0	0
			1888	1208	316	353	11			
1	C	255	Total	C	N	O	S	0	0	0
			1887	1207	315	354	11			
1	D	255	Total	C	N	O	S	0	0	0
			1887	1207	315	354	11			
1	E	256	Total	C	N	O	S	0	0	0
			1895	1212	317	355	11			
1	F	262	Total	C	N	O	S	0	0	0
			1941	1239	325	365	12			
1	G	262	Total	C	N	O	S	0	0	0
			1940	1239	324	365	12			
1	H	259	Total	C	N	O	S	0	0	0
			1908	1219	319	359	11			

There are 184 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	initiating methionine	UNP Q58018
A	-21	GLY	-	expression tag	UNP Q58018
A	-20	SER	-	expression tag	UNP Q58018
A	-19	ASP	-	expression tag	UNP Q58018
A	-18	LYS	-	expression tag	UNP Q58018
A	-17	ILE	-	expression tag	UNP Q58018
A	-16	HIS	-	expression tag	UNP Q58018
A	-15	HIS	-	expression tag	UNP Q58018
A	-14	HIS	-	expression tag	UNP Q58018
A	-13	HIS	-	expression tag	UNP Q58018
A	-12	HIS	-	expression tag	UNP Q58018
A	-11	HIS	-	expression tag	UNP Q58018
A	-10	SER	-	expression tag	UNP Q58018

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	SER	-	expression tag	UNP Q58018
A	-8	GLY	-	expression tag	UNP Q58018
A	-7	GLU	-	expression tag	UNP Q58018
A	-6	ASN	-	expression tag	UNP Q58018
A	-5	LEU	-	expression tag	UNP Q58018
A	-4	TYR	-	expression tag	UNP Q58018
A	-3	PHE	-	expression tag	UNP Q58018
A	-2	GLN	-	expression tag	UNP Q58018
A	-1	GLY	-	expression tag	UNP Q58018
A	0	HIS	-	expression tag	UNP Q58018
B	-22	MET	-	initiating methionine	UNP Q58018
B	-21	GLY	-	expression tag	UNP Q58018
B	-20	SER	-	expression tag	UNP Q58018
B	-19	ASP	-	expression tag	UNP Q58018
B	-18	LYS	-	expression tag	UNP Q58018
B	-17	ILE	-	expression tag	UNP Q58018
B	-16	HIS	-	expression tag	UNP Q58018
B	-15	HIS	-	expression tag	UNP Q58018
B	-14	HIS	-	expression tag	UNP Q58018
B	-13	HIS	-	expression tag	UNP Q58018
B	-12	HIS	-	expression tag	UNP Q58018
B	-11	HIS	-	expression tag	UNP Q58018
B	-10	SER	-	expression tag	UNP Q58018
B	-9	SER	-	expression tag	UNP Q58018
B	-8	GLY	-	expression tag	UNP Q58018
B	-7	GLU	-	expression tag	UNP Q58018
B	-6	ASN	-	expression tag	UNP Q58018
B	-5	LEU	-	expression tag	UNP Q58018
B	-4	TYR	-	expression tag	UNP Q58018
B	-3	PHE	-	expression tag	UNP Q58018
B	-2	GLN	-	expression tag	UNP Q58018
B	-1	GLY	-	expression tag	UNP Q58018
B	0	HIS	-	expression tag	UNP Q58018
C	-22	MET	-	initiating methionine	UNP Q58018
C	-21	GLY	-	expression tag	UNP Q58018
C	-20	SER	-	expression tag	UNP Q58018
C	-19	ASP	-	expression tag	UNP Q58018
C	-18	LYS	-	expression tag	UNP Q58018
C	-17	ILE	-	expression tag	UNP Q58018
C	-16	HIS	-	expression tag	UNP Q58018
C	-15	HIS	-	expression tag	UNP Q58018
C	-14	HIS	-	expression tag	UNP Q58018

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-13	HIS	-	expression tag	UNP Q58018
C	-12	HIS	-	expression tag	UNP Q58018
C	-11	HIS	-	expression tag	UNP Q58018
C	-10	SER	-	expression tag	UNP Q58018
C	-9	SER	-	expression tag	UNP Q58018
C	-8	GLY	-	expression tag	UNP Q58018
C	-7	GLU	-	expression tag	UNP Q58018
C	-6	ASN	-	expression tag	UNP Q58018
C	-5	LEU	-	expression tag	UNP Q58018
C	-4	TYR	-	expression tag	UNP Q58018
C	-3	PHE	-	expression tag	UNP Q58018
C	-2	GLN	-	expression tag	UNP Q58018
C	-1	GLY	-	expression tag	UNP Q58018
C	0	HIS	-	expression tag	UNP Q58018
D	-22	MET	-	initiating methionine	UNP Q58018
D	-21	GLY	-	expression tag	UNP Q58018
D	-20	SER	-	expression tag	UNP Q58018
D	-19	ASP	-	expression tag	UNP Q58018
D	-18	LYS	-	expression tag	UNP Q58018
D	-17	ILE	-	expression tag	UNP Q58018
D	-16	HIS	-	expression tag	UNP Q58018
D	-15	HIS	-	expression tag	UNP Q58018
D	-14	HIS	-	expression tag	UNP Q58018
D	-13	HIS	-	expression tag	UNP Q58018
D	-12	HIS	-	expression tag	UNP Q58018
D	-11	HIS	-	expression tag	UNP Q58018
D	-10	SER	-	expression tag	UNP Q58018
D	-9	SER	-	expression tag	UNP Q58018
D	-8	GLY	-	expression tag	UNP Q58018
D	-7	GLU	-	expression tag	UNP Q58018
D	-6	ASN	-	expression tag	UNP Q58018
D	-5	LEU	-	expression tag	UNP Q58018
D	-4	TYR	-	expression tag	UNP Q58018
D	-3	PHE	-	expression tag	UNP Q58018
D	-2	GLN	-	expression tag	UNP Q58018
D	-1	GLY	-	expression tag	UNP Q58018
D	0	HIS	-	expression tag	UNP Q58018
E	-22	MET	-	initiating methionine	UNP Q58018
E	-21	GLY	-	expression tag	UNP Q58018
E	-20	SER	-	expression tag	UNP Q58018
E	-19	ASP	-	expression tag	UNP Q58018
E	-18	LYS	-	expression tag	UNP Q58018

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-17	ILE	-	expression tag	UNP Q58018
E	-16	HIS	-	expression tag	UNP Q58018
E	-15	HIS	-	expression tag	UNP Q58018
E	-14	HIS	-	expression tag	UNP Q58018
E	-13	HIS	-	expression tag	UNP Q58018
E	-12	HIS	-	expression tag	UNP Q58018
E	-11	HIS	-	expression tag	UNP Q58018
E	-10	SER	-	expression tag	UNP Q58018
E	-9	SER	-	expression tag	UNP Q58018
E	-8	GLY	-	expression tag	UNP Q58018
E	-7	GLU	-	expression tag	UNP Q58018
E	-6	ASN	-	expression tag	UNP Q58018
E	-5	LEU	-	expression tag	UNP Q58018
E	-4	TYR	-	expression tag	UNP Q58018
E	-3	PHE	-	expression tag	UNP Q58018
E	-2	GLN	-	expression tag	UNP Q58018
E	-1	GLY	-	expression tag	UNP Q58018
E	0	HIS	-	expression tag	UNP Q58018
F	-22	MET	-	initiating methionine	UNP Q58018
F	-21	GLY	-	expression tag	UNP Q58018
F	-20	SER	-	expression tag	UNP Q58018
F	-19	ASP	-	expression tag	UNP Q58018
F	-18	LYS	-	expression tag	UNP Q58018
F	-17	ILE	-	expression tag	UNP Q58018
F	-16	HIS	-	expression tag	UNP Q58018
F	-15	HIS	-	expression tag	UNP Q58018
F	-14	HIS	-	expression tag	UNP Q58018
F	-13	HIS	-	expression tag	UNP Q58018
F	-12	HIS	-	expression tag	UNP Q58018
F	-11	HIS	-	expression tag	UNP Q58018
F	-10	SER	-	expression tag	UNP Q58018
F	-9	SER	-	expression tag	UNP Q58018
F	-8	GLY	-	expression tag	UNP Q58018
F	-7	GLU	-	expression tag	UNP Q58018
F	-6	ASN	-	expression tag	UNP Q58018
F	-5	LEU	-	expression tag	UNP Q58018
F	-4	TYR	-	expression tag	UNP Q58018
F	-3	PHE	-	expression tag	UNP Q58018
F	-2	GLN	-	expression tag	UNP Q58018
F	-1	GLY	-	expression tag	UNP Q58018
F	0	HIS	-	expression tag	UNP Q58018
G	-22	MET	-	initiating methionine	UNP Q58018

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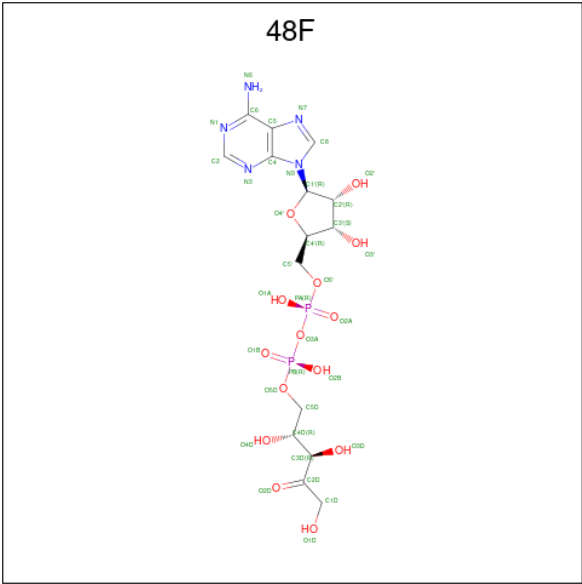
Chain	Residue	Modelled	Actual	Comment	Reference
G	-21	GLY	-	expression tag	UNP Q58018
G	-20	SER	-	expression tag	UNP Q58018
G	-19	ASP	-	expression tag	UNP Q58018
G	-18	LYS	-	expression tag	UNP Q58018
G	-17	ILE	-	expression tag	UNP Q58018
G	-16	HIS	-	expression tag	UNP Q58018
G	-15	HIS	-	expression tag	UNP Q58018
G	-14	HIS	-	expression tag	UNP Q58018
G	-13	HIS	-	expression tag	UNP Q58018
G	-12	HIS	-	expression tag	UNP Q58018
G	-11	HIS	-	expression tag	UNP Q58018
G	-10	SER	-	expression tag	UNP Q58018
G	-9	SER	-	expression tag	UNP Q58018
G	-8	GLY	-	expression tag	UNP Q58018
G	-7	GLU	-	expression tag	UNP Q58018
G	-6	ASN	-	expression tag	UNP Q58018
G	-5	LEU	-	expression tag	UNP Q58018
G	-4	TYR	-	expression tag	UNP Q58018
G	-3	PHE	-	expression tag	UNP Q58018
G	-2	GLN	-	expression tag	UNP Q58018
G	-1	GLY	-	expression tag	UNP Q58018
G	0	HIS	-	expression tag	UNP Q58018
H	-22	MET	-	initiating methionine	UNP Q58018
H	-21	GLY	-	expression tag	UNP Q58018
H	-20	SER	-	expression tag	UNP Q58018
H	-19	ASP	-	expression tag	UNP Q58018
H	-18	LYS	-	expression tag	UNP Q58018
H	-17	ILE	-	expression tag	UNP Q58018
H	-16	HIS	-	expression tag	UNP Q58018
H	-15	HIS	-	expression tag	UNP Q58018
H	-14	HIS	-	expression tag	UNP Q58018
H	-13	HIS	-	expression tag	UNP Q58018
H	-12	HIS	-	expression tag	UNP Q58018
H	-11	HIS	-	expression tag	UNP Q58018
H	-10	SER	-	expression tag	UNP Q58018
H	-9	SER	-	expression tag	UNP Q58018
H	-8	GLY	-	expression tag	UNP Q58018
H	-7	GLU	-	expression tag	UNP Q58018
H	-6	ASN	-	expression tag	UNP Q58018
H	-5	LEU	-	expression tag	UNP Q58018
H	-4	TYR	-	expression tag	UNP Q58018
H	-3	PHE	-	expression tag	UNP Q58018

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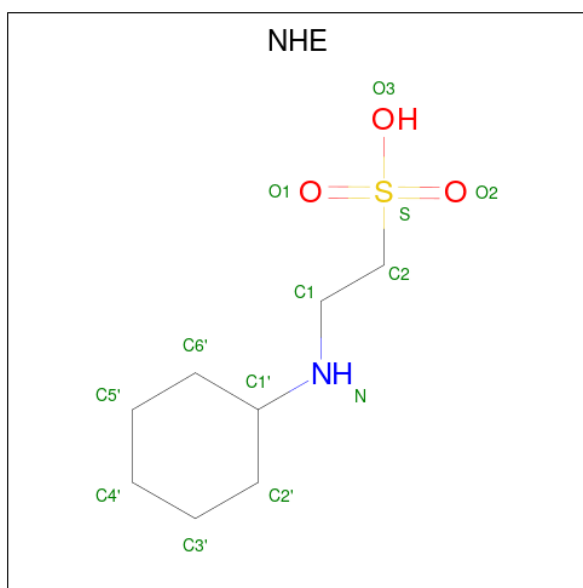
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Chain	Residue	Modelled	Actual	Comment	Reference
H	-2	GLN	-	expression tag	UNP Q58018
H	-1	GLY	-	expression tag	UNP Q58018
H	0	HIS	-	expression tag	UNP Q58018

- Molecule 2 is [[(2R,3S,4R,5R)-5-(6-aminopurin-9-yl)-3,4-bis(oxidanyl)oxolan-2-yl]methoxy-oxidanyl-phosphoryl] [(2R,3R)-2,3,5-tris(oxidanyl)-4-oxidanylidene-pentyl] hydrogen phosphate (CCD ID: 48F) (formula: C₁₅H₂₃N₅O₁₄P₂).



(formula: C₈H₁₇NO₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			13	8	1	3	1		
3	C	1	Total	C	N	O	S	0	0
			13	8	1	3	1		
3	F	1	Total	C	N	O	S	0	0
			13	8	1	3	1		
3	G	1	Total	C	N	O	S	0	0
			13	8	1	3	1		
3	H	1	Total	C	N	O	S	0	0
			13	8	1	3	1		

- Molecule 4 is water.

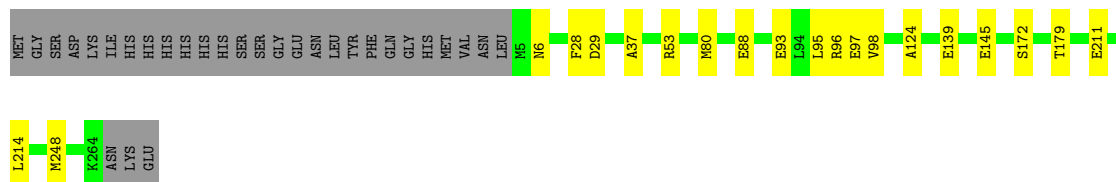
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	3	Total	O	0	0
			3	3		
4	C	2	Total	O	0	0
			2	2		
4	D	2	Total	O	0	0
			2	2		
4	E	1	Total	O	0	0
			1	1		
4	F	5	Total	O	0	0
			5	5		
4	G	8	Total	O	0	0
			8	8		

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.

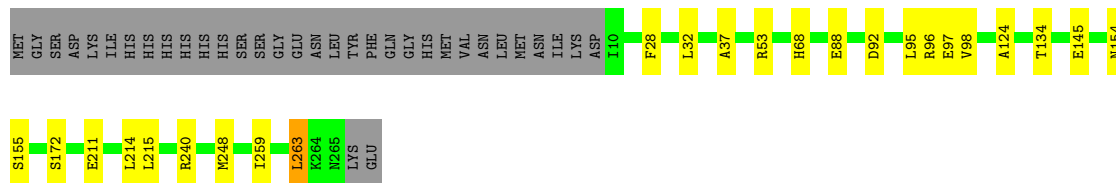
- Molecule 1: Putative ribose 1,5-bisphosphate isomerase

Chain A: 



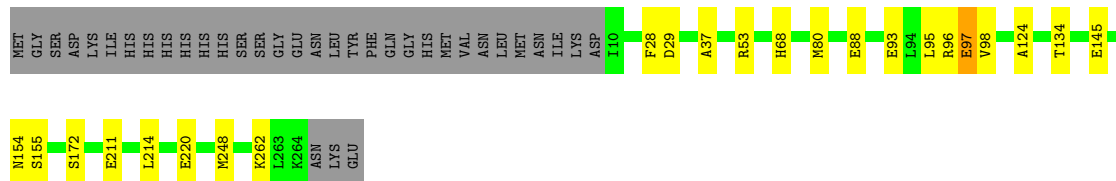
- Molecule 1: Putative ribose 1,5-bisphosphate isomerase

Chain B: 



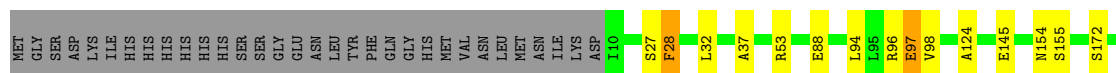
- Molecule 1: Putative ribose 1,5-bisphosphate isomerase

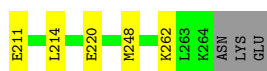
Chain C: 



- Molecule 1: Putative ribose 1,5-bisphosphate isomerase

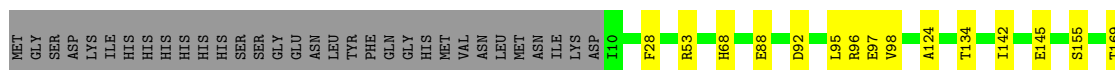
Chain D: 





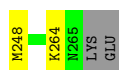
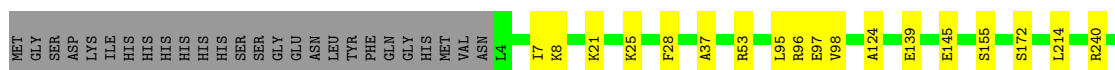
- Molecule 1: Putative ribose 1,5-bisphosphate isomerase

Chain E: 81% 8% 12%



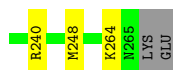
- Molecule 1: Putative ribose 1,5-bisphosphate isomerase

Chain F: 83% 7% 10%



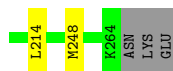
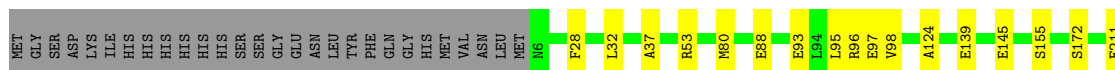
- Molecule 1: Putative ribose 1,5-bisphosphate isomerase

Chain G: 83% 7% 10%



- Molecule 1: Putative ribose 1,5-bisphosphate isomerase

Chain H: 83% 7% 11%



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	180.76Å 180.76Å 73.33Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.00 – 2.71 40.00 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.0 (40.00-2.71) 99.3 (40.00-2.71)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.177 , 0.203 0.180 , 0.202	Depositor DCC
R_{free} test set	3646 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	55.6	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 28.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l 0.477 for h,-h-k,-l 0.017 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	15644	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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: NHE, 48F

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	0.67	0/1954	0.83	1/2647 (0.0%)
1	B	0.73	0/1918	0.85	0/2600
1	C	0.66	0/1917	0.81	1/2598 (0.0%)
1	D	0.68	0/1917	0.83	3/2598 (0.1%)
1	E	0.75	0/1925	0.83	0/2608
1	F	0.68	0/1971	0.83	1/2670 (0.0%)
1	G	0.69	0/1970	0.83	1/2670 (0.0%)
1	H	0.67	0/1938	0.82	0/2629
All	All	0.69	0/15510	0.83	7/21020 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	264	LYS	N-CA-C	-6.97	104.75	113.18
1	A	29	ASP	N-CA-C	-6.69	104.07	111.36
1	C	29	ASP	N-CA-C	-6.27	104.53	111.36
1	G	264	LYS	N-CA-C	-6.11	105.79	113.18
1	D	27	SER	CA-C-N	5.32	127.41	120.28
1	D	27	SER	C-N-CA	5.32	127.41	120.28
1	D	28	PHE	N-CA-CB	-5.25	102.40	110.12

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	1924	0	1946	37	0
1	B	1888	0	1905	42	0
1	C	1887	0	1907	40	0
1	D	1887	0	1907	42	0
1	E	1895	0	1918	15	0
1	F	1941	0	1958	23	0
1	G	1940	0	1956	24	0
1	H	1908	0	1915	20	0
2	A	36	0	0	1	0
2	B	36	0	0	0	0
2	C	36	0	0	0	0
2	D	36	0	0	0	0
2	E	36	0	0	0	0
2	F	36	0	0	0	0
2	G	36	0	0	0	0
2	H	36	0	0	0	0
3	A	13	0	16	0	0
3	C	13	0	17	0	0
3	F	13	0	16	0	0
3	G	13	0	16	0	0
3	H	13	0	16	0	0
4	B	3	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
4	E	1	0	0	0	0
4	F	5	0	0	0	0
4	G	8	0	0	2	0
All	All	15644	0	15493	165	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.

All (165) 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:28:PHE:CE1	1:B:28:PHE:HB2	1.08	1.61
1:A:28:PHE:CD2	1:B:28:PHE:CD1	1.91	1.56
1:C:28:PHE:CD2	1:D:28:PHE:CD1	1.96	1.53
1:A:28:PHE:CZ	1:B:28:PHE:HA	1.42	1.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:PHE:CZ	1:B:28:PHE:CA	1.97	1.44
1:A:28:PHE:CE1	1:B:28:PHE:CB	2.02	1.42
1:C:28:PHE:CE2	1:D:28:PHE:HA	1.61	1.33
1:A:28:PHE:CZ	1:B:28:PHE:HB2	1.64	1.32
1:A:28:PHE:CZ	1:B:28:PHE:CB	2.10	1.31
1:C:28:PHE:CD2	1:D:28:PHE:CG	2.16	1.31
1:A:28:PHE:CD1	1:B:28:PHE:HB2	1.65	1.31
1:C:28:PHE:CE2	1:D:28:PHE:CD1	2.21	1.28
1:C:28:PHE:CZ	1:D:28:PHE:HA	1.67	1.28
1:C:28:PHE:CZ	1:D:28:PHE:CA	2.25	1.19
1:A:28:PHE:CE2	1:B:28:PHE:HA	1.76	1.18
1:C:28:PHE:CE2	1:D:28:PHE:CA	2.25	1.17
1:F:28:PHE:CD2	1:G:28:PHE:CD1	2.34	1.16
1:C:28:PHE:CE2	1:D:28:PHE:CG	2.34	1.16
1:A:28:PHE:CE2	1:B:28:PHE:CD1	2.35	1.14
1:C:28:PHE:HD2	1:D:28:PHE:CE1	1.72	1.07
1:A:28:PHE:CG	1:B:28:PHE:CG	2.47	1.02
1:A:28:PHE:CE2	1:B:28:PHE:HD1	1.72	1.02
1:C:28:PHE:CE2	1:D:28:PHE:CB	2.42	1.02
1:C:28:PHE:CD2	1:D:28:PHE:CE1	2.48	1.01
1:A:28:PHE:CD2	1:B:28:PHE:CG	2.49	1.01
1:H:28:PHE:CE1	1:H:32:LEU:HD11	1.97	0.99
1:C:28:PHE:CZ	1:D:28:PHE:CB	2.49	0.96
1:A:28:PHE:CG	1:B:28:PHE:CD1	2.53	0.95
1:F:28:PHE:HD2	1:G:28:PHE:CD1	1.75	0.95
1:F:28:PHE:CE2	1:G:28:PHE:HA	2.01	0.95
1:A:28:PHE:HZ	1:B:28:PHE:HA	1.29	0.90
1:C:28:PHE:HE2	1:D:28:PHE:HA	1.37	0.89
1:A:28:PHE:CE2	1:B:28:PHE:CB	2.59	0.86
1:F:28:PHE:CD2	1:G:28:PHE:CG	2.63	0.85
1:C:28:PHE:HD2	1:D:28:PHE:CD1	1.67	0.85
1:H:28:PHE:CE1	1:H:32:LEU:CD1	2.61	0.84
1:C:28:PHE:CZ	1:D:28:PHE:HB2	2.14	0.82
1:A:28:PHE:CD2	1:B:28:PHE:CE1	2.69	0.80
1:A:28:PHE:CD1	1:B:28:PHE:CB	2.46	0.79
1:A:28:PHE:CD1	1:B:28:PHE:CG	2.71	0.78
1:D:28:PHE:HE1	1:D:32:LEU:HD11	1.46	0.78
1:A:28:PHE:CE2	1:B:28:PHE:CA	2.48	0.78
1:C:28:PHE:CE2	1:D:28:PHE:HD1	2.02	0.77
1:H:28:PHE:HE1	1:H:32:LEU:HD11	1.51	0.75
1:E:53:ARG:NH1	1:E:97:GLU:OE2	2.20	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:PHE:CE2	1:B:28:PHE:CG	2.75	0.73
1:C:28:PHE:HB3	1:D:28:PHE:CE2	2.23	0.73
1:A:28:PHE:HZ	1:B:28:PHE:CA	1.89	0.71
1:C:28:PHE:HZ	1:D:28:PHE:HA	1.50	0.71
1:C:28:PHE:CG	1:D:28:PHE:CG	2.77	0.71
1:B:53:ARG:NH1	1:B:97:GLU:OE2	2.24	0.70
1:F:53:ARG:NH1	1:F:97:GLU:OE2	2.26	0.69
1:C:28:PHE:CD2	1:D:28:PHE:CD2	2.79	0.69
1:H:53:ARG:NH1	1:H:97:GLU:OE2	2.26	0.68
1:C:28:PHE:CG	1:D:28:PHE:CD2	2.83	0.67
1:A:28:PHE:HD2	1:B:28:PHE:CD1	2.00	0.65
1:B:28:PHE:HE1	1:B:32:LEU:HD11	1.61	0.65
1:E:215:LEU:HD22	1:F:7:ILE:HG22	1.78	0.65
1:F:28:PHE:CE2	1:G:28:PHE:CA	2.78	0.64
1:C:28:PHE:CE1	1:D:28:PHE:HB2	2.31	0.64
1:A:28:PHE:CD2	1:B:28:PHE:HD1	1.71	0.63
1:C:53:ARG:NH1	1:C:97:GLU:OE2	2.32	0.62
1:D:53:ARG:NH1	1:D:97:GLU:OE2	2.32	0.62
1:B:28:PHE:CE1	1:B:32:LEU:HD11	2.34	0.62
1:G:53:ARG:NH1	1:G:97:GLU:OE2	2.33	0.61
1:E:240:ARG:NH1	1:G:139:GLU:OE1	2.33	0.60
1:D:28:PHE:CE1	1:D:32:LEU:HD11	2.34	0.59
1:B:240:ARG:NH1	1:F:139:GLU:OE1	2.32	0.58
1:E:28:PHE:CD1	1:H:28:PHE:HB2	2.38	0.58
1:A:53:ARG:NH1	1:A:97:GLU:OE2	2.37	0.58
1:F:28:PHE:HE2	1:G:28:PHE:HA	1.66	0.57
1:C:96:ARG:HH21	1:C:96:ARG:HG2	1.69	0.57
1:H:28:PHE:CZ	1:H:32:LEU:HD11	2.40	0.57
1:A:28:PHE:CZ	1:B:28:PHE:N	2.69	0.56
1:E:98:VAL:HG12	1:E:98:VAL:O	2.05	0.56
1:A:98:VAL:HG12	1:A:98:VAL:O	2.04	0.56
1:E:215:LEU:HD22	1:F:7:ILE:CG2	2.35	0.56
1:E:28:PHE:CD2	1:H:28:PHE:CD1	2.93	0.56
1:F:98:VAL:O	1:F:98:VAL:HG12	2.05	0.56
1:B:215:LEU:HD22	1:G:7:ILE:HG22	1.87	0.56
1:D:98:VAL:HG12	1:D:98:VAL:O	2.06	0.55
1:G:98:VAL:HG12	1:G:98:VAL:O	2.07	0.55
1:H:98:VAL:HG12	1:H:98:VAL:O	2.05	0.55
1:B:98:VAL:O	1:B:98:VAL:HG12	2.05	0.55
1:G:98:VAL:HG13	1:G:124:ALA:HB2	1.88	0.55
1:C:98:VAL:HG12	1:C:98:VAL:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:VAL:HG13	1:B:124:ALA:HB2	1.89	0.54
1:F:96:ARG:HG2	1:F:96:ARG:HH21	1.73	0.53
1:D:98:VAL:HG13	1:D:124:ALA:HB2	1.91	0.53
1:F:98:VAL:HG13	1:F:124:ALA:HB2	1.89	0.53
1:E:98:VAL:HG13	1:E:124:ALA:HB2	1.90	0.53
1:C:28:PHE:HE2	1:D:28:PHE:CD1	2.13	0.52
1:C:98:VAL:HG13	1:C:124:ALA:HB2	1.91	0.52
1:A:98:VAL:HG13	1:A:124:ALA:HB2	1.92	0.52
1:G:248:MET:HG2	4:G:403:HOH:O	2.10	0.52
1:C:28:PHE:CB	1:D:28:PHE:CE2	2.92	0.51
1:D:28:PHE:HE1	1:D:32:LEU:CD1	2.21	0.50
1:H:98:VAL:HG13	1:H:124:ALA:HB2	1.92	0.50
1:D:28:PHE:CE1	1:D:32:LEU:CD1	2.95	0.50
1:G:28:PHE:HE1	1:G:32:LEU:HD11	1.76	0.50
1:D:96:ARG:HH21	1:D:96:ARG:HG2	1.76	0.50
1:A:93:GLU:O	1:A:97:GLU:HB2	2.12	0.49
1:C:154:ASN:HA	1:H:80:MET:HG3	1.94	0.49
1:B:214:LEU:HD21	1:B:248:MET:HE2	1.95	0.49
1:H:93:GLU:O	1:H:97:GLU:HB2	2.14	0.48
1:A:37:ALA:O	1:A:172:SER:HA	2.13	0.48
1:A:88:GLU:OE2	1:A:211:GLU:OE2	2.32	0.48
1:E:214:LEU:HD21	1:E:248:MET:HE2	1.95	0.48
1:C:96:ARG:HG2	1:C:96:ARG:NH2	2.28	0.48
1:C:214:LEU:HD21	1:C:248:MET:HE2	1.96	0.48
1:G:28:PHE:CE1	1:G:32:LEU:HD11	2.48	0.48
1:D:28:PHE:CD1	1:D:28:PHE:C	2.92	0.47
1:F:28:PHE:HD2	1:G:28:PHE:CE1	2.29	0.47
1:B:215:LEU:HD22	1:G:7:ILE:CG2	2.43	0.47
1:A:80:MET:HG3	1:D:154:ASN:HA	1.95	0.47
1:A:139:GLU:OE1	1:G:240:ARG:NH1	2.45	0.47
1:D:214:LEU:HD21	1:D:248:MET:HE2	1.96	0.47
1:H:37:ALA:O	1:H:172:SER:HA	2.13	0.47
1:A:214:LEU:HD21	1:A:248:MET:HE2	1.97	0.47
1:B:28:PHE:CE1	1:B:32:LEU:CD1	2.97	0.47
1:E:92:ASP:O	1:E:96:ARG:HG3	2.15	0.47
1:F:28:PHE:CD2	1:G:28:PHE:CB	2.98	0.47
1:B:92:ASP:O	1:B:96:ARG:HG3	2.15	0.46
1:H:214:LEU:HD21	1:H:248:MET:HE2	1.97	0.46
1:C:28:PHE:HE2	1:D:28:PHE:HD1	1.56	0.45
1:B:88:GLU:OE2	1:B:211:GLU:OE2	2.35	0.45
1:F:37:ALA:O	1:F:172:SER:HA	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:96:ARG:HG2	1:F:96:ARG:NH2	2.32	0.45
1:C:37:ALA:O	1:C:172:SER:HA	2.17	0.45
1:D:88:GLU:OE2	1:D:211:GLU:OE2	2.35	0.45
1:G:214:LEU:HD21	1:G:248:MET:HE2	1.98	0.45
1:F:214:LEU:HD21	1:F:248:MET:HE2	1.97	0.45
1:F:28:PHE:CE2	1:G:28:PHE:CB	3.00	0.45
1:F:240:ARG:NH1	1:H:139:GLU:OE1	2.45	0.45
1:C:88:GLU:OE2	1:C:211:GLU:OE2	2.36	0.44
1:D:220:GLU:OE2	1:D:262:LYS:NZ	2.48	0.44
1:C:93:GLU:O	1:C:97:GLU:HB3	2.17	0.44
1:C:28:PHE:CZ	1:D:28:PHE:N	2.84	0.44
1:H:88:GLU:OE2	1:H:211:GLU:OE2	2.36	0.44
1:H:96:ARG:HH21	1:H:96:ARG:HG2	1.83	0.43
1:A:96:ARG:HG2	1:A:96:ARG:HH21	1.84	0.43
1:G:37:ALA:O	1:G:172:SER:HA	2.18	0.43
1:B:154:ASN:HA	1:C:80:MET:HG3	2.00	0.43
1:E:28:PHE:CG	1:H:28:PHE:CG	3.07	0.43
1:B:259:ILE:O	1:B:263:LEU:HB2	2.19	0.43
1:D:37:ALA:O	1:D:172:SER:HA	2.19	0.43
1:B:68:HIS:O	1:B:134:THR:HB	2.19	0.43
1:E:88:GLU:OE2	1:E:211:GLU:OE2	2.36	0.43
1:F:28:PHE:CG	1:G:28:PHE:CG	3.05	0.43
1:H:28:PHE:CD1	1:H:28:PHE:O	2.71	0.43
1:E:68:HIS:O	1:E:134:THR:HB	2.19	0.43
1:E:215:LEU:HB3	1:F:8:LYS:HA	2.00	0.43
1:A:28:PHE:CG	1:B:28:PHE:CD2	3.03	0.42
1:F:21:LYS:O	1:F:25:LYS:HB2	2.20	0.42
1:H:28:PHE:HE1	1:H:32:LEU:CD1	2.20	0.42
1:B:37:ALA:O	1:B:172:SER:HA	2.20	0.42
1:C:28:PHE:CD2	1:D:28:PHE:CZ	3.05	0.41
1:E:142:ILE:HD11	1:E:169:THR:HB	2.02	0.41
1:C:68:HIS:O	1:C:134:THR:HB	2.21	0.41
1:C:220:GLU:OE2	1:C:262:LYS:NZ	2.47	0.41
1:D:94:LEU:HD23	1:D:94:LEU:HA	1.85	0.41
1:G:88:GLU:OE2	1:G:211:GLU:OE2	2.39	0.41
1:H:28:PHE:CD1	1:H:28:PHE:C	2.98	0.41
1:A:179:THR:HB	2:A:301:48F:C8	2.51	0.40
1:G:182:GLU:CG	4:G:404:HOH:O	2.69	0.40

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	258/290 (89%)	249 (96%)	9 (4%)	0	100	100
1	B	254/290 (88%)	246 (97%)	8 (3%)	0	100	100
1	C	253/290 (87%)	245 (97%)	8 (3%)	0	100	100
1	D	253/290 (87%)	245 (97%)	8 (3%)	0	100	100
1	E	254/290 (88%)	246 (97%)	8 (3%)	0	100	100
1	F	260/290 (90%)	251 (96%)	9 (4%)	0	100	100
1	G	260/290 (90%)	250 (96%)	10 (4%)	0	100	100
1	H	257/290 (89%)	248 (96%)	9 (4%)	0	100	100
All	All	2049/2320 (88%)	1980 (97%)	69 (3%)	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	195/229 (85%)	192 (98%)	3 (2%)	57	79
1	B	190/229 (83%)	186 (98%)	4 (2%)	47	73
1	C	191/229 (83%)	187 (98%)	4 (2%)	47	73
1	D	191/229 (83%)	188 (98%)	3 (2%)	55	78
1	E	192/229 (84%)	187 (97%)	5 (3%)	40	69
1	F	197/229 (86%)	194 (98%)	3 (2%)	57	79

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	197/229 (86%)	194 (98%)	3 (2%)	57	79
1	H	192/229 (84%)	189 (98%)	3 (2%)	55	78
All	All	1545/1832 (84%)	1517 (98%)	28 (2%)	51	76

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	95	LEU
1	A	145	GLU
1	B	95	LEU
1	B	145	GLU
1	B	155	SER
1	B	263	LEU
1	C	95	LEU
1	C	97	GLU
1	C	145	GLU
1	C	155	SER
1	D	97	GLU
1	D	145	GLU
1	D	155	SER
1	E	95	LEU
1	E	145	GLU
1	E	155	SER
1	E	263	LEU
1	E	264	LYS
1	F	95	LEU
1	F	145	GLU
1	F	155	SER
1	G	95	LEU
1	G	145	GLU
1	G	155	SER
1	H	95	LEU
1	H	145	GLU
1	H	155	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	236	HIS

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Mol	Chain	Res	Type
1	B	236	HIS
1	B	265	ASN
1	C	236	HIS
1	D	236	HIS
1	E	236	HIS
1	E	265	ASN
1	F	236	HIS
1	G	236	HIS
1	H	236	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](#)

13 ligands are modelled in this entry.

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$
2	48F	C	301	-	38,38,38	1.55	7 (18%)	49,57,57	1.80	11 (22%)
2	48F	G	301	-	38,38,38	1.65	6 (15%)	49,57,57	1.84	11 (22%)
3	NHE	H	302	-	13,13,13	2.93	2 (15%)	16,17,17	1.47	2 (12%)
2	48F	E	301	-	38,38,38	1.72	6 (15%)	49,57,57	1.67	7 (14%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	48F	D	301	-	38,38,38	1.57	5 (13%)	49,57,57	1.86	10 (20%)
3	NHE	C	302	-	13,13,13	2.32	2 (15%)	16,17,17	1.73	3 (18%)
2	48F	F	301	-	38,38,38	1.60	5 (13%)	49,57,57	1.78	12 (24%)
2	48F	H	301	-	38,38,38	1.70	6 (15%)	49,57,57	1.80	11 (22%)
3	NHE	A	302	-	13,13,13	2.93	2 (15%)	16,17,17	1.81	3 (18%)
2	48F	A	301	-	38,38,38	1.57	6 (15%)	49,57,57	1.88	12 (24%)
3	NHE	G	302	-	13,13,13	3.06	2 (15%)	16,17,17	1.56	1 (6%)
3	NHE	F	302	-	13,13,13	2.91	2 (15%)	16,17,17	1.79	3 (18%)
2	48F	B	301	-	38,38,38	1.67	7 (18%)	49,57,57	1.70	9 (18%)

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
2	48F	C	301	-	-	4/32/48/48	0/3/3/3
2	48F	G	301	-	-	5/32/48/48	0/3/3/3
3	NHE	H	302	-	-	1/7/15/15	0/1/1/1
2	48F	E	301	-	-	14/32/48/48	0/3/3/3
2	48F	D	301	-	-	8/32/48/48	0/3/3/3
3	NHE	C	302	-	-	5/7/15/15	0/1/1/1
2	48F	F	301	-	-	5/32/48/48	0/3/3/3
2	48F	H	301	-	-	3/32/48/48	0/3/3/3
3	NHE	A	302	-	-	4/7/15/15	0/1/1/1
2	48F	A	301	-	-	3/32/48/48	0/3/3/3
3	NHE	G	302	-	-	2/7/15/15	0/1/1/1
3	NHE	F	302	-	-	7/7/15/15	0/1/1/1
2	48F	B	301	-	-	8/32/48/48	0/3/3/3

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	302	NHE	C2-S	-8.36	1.65	1.77
3	A	302	NHE	C2-S	-7.94	1.66	1.77
3	F	302	NHE	C2-S	-7.85	1.66	1.77
3	H	302	NHE	C2-S	-7.76	1.66	1.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	302	NHE	O1-S	7.00	1.64	1.45
3	H	302	NHE	O1-S	6.93	1.64	1.45
3	C	302	NHE	C2-S	-6.87	1.68	1.77
3	A	302	NHE	O2-S	6.78	1.64	1.45
3	F	302	NHE	O1-S	6.69	1.64	1.45
2	E	301	48F	C5-C4	5.91	1.49	1.39
2	B	301	48F	C5-C4	5.68	1.49	1.39
2	G	301	48F	C5-C4	5.26	1.48	1.39
2	A	301	48F	C5-C4	5.21	1.48	1.39
2	H	301	48F	C5-C4	5.12	1.48	1.39
2	C	301	48F	C5-C4	5.06	1.48	1.39
2	F	301	48F	C5-C4	5.05	1.48	1.39
2	D	301	48F	C5-C4	4.95	1.47	1.39
2	H	301	48F	PB-O3A	4.69	1.64	1.59
2	H	301	48F	PA-O3A	4.54	1.64	1.59
3	C	302	NHE	O3-S	4.44	1.63	1.47
2	F	301	48F	PA-O3A	4.15	1.64	1.59
2	D	301	48F	PB-O3A	3.82	1.63	1.59
2	B	301	48F	PA-O3A	3.80	1.63	1.59
2	G	301	48F	PA-O3A	3.77	1.63	1.59
2	G	301	48F	PB-O3A	3.73	1.63	1.59
2	A	301	48F	PA-O3A	3.72	1.63	1.59
2	E	301	48F	PA-O3A	3.69	1.63	1.59
2	C	301	48F	PA-O3A	3.40	1.63	1.59
2	D	301	48F	PA-O3A	3.36	1.63	1.59
2	E	301	48F	PB-O3A	3.34	1.63	1.59
2	A	301	48F	PB-O3A	3.28	1.63	1.59
2	F	301	48F	PB-O3A	3.21	1.63	1.59
2	C	301	48F	PB-O3A	3.19	1.62	1.59
2	A	301	48F	C8-N7	3.14	1.37	1.31
2	C	301	48F	C5-C6	3.11	1.49	1.41
2	B	301	48F	PB-O3A	3.10	1.62	1.59
2	B	301	48F	C8-N7	3.06	1.37	1.31
2	F	301	48F	C8-N7	3.02	1.37	1.31
2	D	301	48F	C5-C6	3.01	1.49	1.41
2	E	301	48F	C8-N7	3.01	1.37	1.31
2	F	301	48F	C5-C6	3.00	1.49	1.41
2	G	301	48F	C5-C6	2.99	1.49	1.41
2	E	301	48F	C5-C6	2.98	1.49	1.41
2	B	301	48F	C5-C6	2.82	1.48	1.41
2	H	301	48F	C8-N7	2.77	1.37	1.31
2	G	301	48F	C8-N7	2.75	1.37	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	48F	C8-N7	2.74	1.36	1.31
2	C	301	48F	C8-N7	2.68	1.36	1.31
2	A	301	48F	C5-C6	2.62	1.48	1.41
2	G	301	48F	C5D-C4D	2.57	1.55	1.51
2	H	301	48F	C5-C6	2.50	1.48	1.41
2	E	301	48F	C5D-C4D	2.38	1.55	1.51
2	H	301	48F	C5-N7	-2.29	1.34	1.39
2	B	301	48F	C5-N7	-2.19	1.35	1.39
2	A	301	48F	C5-N7	-2.17	1.35	1.39
2	C	301	48F	C5D-C4D	2.11	1.54	1.51
2	B	301	48F	C5D-C4D	2.10	1.54	1.51
2	C	301	48F	C5-N7	-2.06	1.35	1.39

All (95) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	48F	C5-C4-N3	-5.84	118.67	126.72
2	D	301	48F	C5-C4-N3	-5.80	118.73	126.72
2	B	301	48F	C5-C4-N3	-5.77	118.78	126.72
2	A	301	48F	C5-C4-N3	-5.65	118.94	126.72
2	F	301	48F	C5-C4-N3	-5.53	119.10	126.72
2	H	301	48F	C5-C4-N3	-5.40	119.28	126.72
2	G	301	48F	C5-C4-N3	-5.33	119.38	126.72
3	G	302	NHE	C1-N-C1'	5.28	124.34	114.18
2	E	301	48F	C5-C4-N3	-5.24	119.50	126.72
3	C	302	NHE	C1-N-C1'	4.95	123.72	114.18
3	A	302	NHE	C1-N-C1'	4.95	123.71	114.18
2	B	301	48F	N3-C4-N9	4.89	135.49	127.17
2	C	301	48F	N3-C4-N9	4.63	135.04	127.17
2	H	301	48F	N3-C4-N9	4.56	134.92	127.17
2	G	301	48F	N3-C2-N1	-4.51	121.76	128.58
2	D	301	48F	N3-C4-N9	4.50	134.82	127.17
2	A	301	48F	N3-C2-N1	-4.40	121.93	128.58
2	A	301	48F	N3-C4-N9	4.38	134.61	127.17
2	F	301	48F	N3-C2-N1	-4.37	121.96	128.58
3	H	302	NHE	C1-N-C1'	4.37	122.60	114.18
2	G	301	48F	N3-C4-N9	4.36	134.58	127.17
2	E	301	48F	N3-C4-N9	4.31	134.49	127.17
3	F	302	NHE	C1-N-C1'	4.29	122.44	114.18
3	F	302	NHE	O1-S-C2	4.21	113.10	106.73
2	F	301	48F	C2-N3-C4	4.12	121.88	111.83
2	A	301	48F	C2-N3-C4	4.11	121.87	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	301	48F	C2-N3-C4	4.10	121.85	111.83
2	H	301	48F	N3-C2-N1	-4.06	122.43	128.58
2	D	301	48F	C2-N3-C4	4.01	121.63	111.83
2	F	301	48F	N3-C4-N9	3.99	133.95	127.17
2	C	301	48F	C2-N3-C4	3.97	121.52	111.83
2	B	301	48F	N3-C2-N1	-3.96	122.58	128.58
2	B	301	48F	C2-N3-C4	3.84	121.21	111.83
2	D	301	48F	C4-C5-N7	-3.83	106.21	110.58
2	E	301	48F	N3-C2-N1	-3.81	122.82	128.58
2	H	301	48F	C2-N3-C4	3.80	121.11	111.83
2	D	301	48F	N3-C2-N1	-3.77	122.88	128.58
2	C	301	48F	N3-C2-N1	-3.63	123.08	128.58
2	C	301	48F	C4-C5-N7	-3.61	106.46	110.58
2	E	301	48F	C2-N3-C4	3.52	120.42	111.83
2	A	301	48F	C4-C5-N7	-3.43	106.66	110.58
3	A	302	NHE	O1-S-C2	3.32	111.74	106.73
2	F	301	48F	C4-C5-N7	-3.29	106.82	110.58
2	G	301	48F	O4D-C4D-C3D	-3.29	104.90	109.67
2	H	301	48F	C4-N9-C8	3.28	109.18	105.74
2	H	301	48F	C4-C5-N7	-3.15	106.98	110.58
2	H	301	48F	O4D-C4D-C3D	-3.05	105.24	109.67
2	G	301	48F	C4-C5-N7	-3.03	107.12	110.58
2	D	301	48F	C5-N7-C8	2.99	108.16	103.45
2	D	301	48F	C4-N9-C8	2.93	108.82	105.74
2	C	301	48F	C5-N7-C8	2.89	108.00	103.45
2	E	301	48F	C4-C5-N7	-2.88	107.29	110.58
2	A	301	48F	C4-N9-C8	2.88	108.76	105.74
2	G	301	48F	C4-N9-C8	2.83	108.71	105.74
2	C	301	48F	C4-N9-C8	2.77	108.64	105.74
3	C	302	NHE	O1-S-C2	2.75	110.88	106.73
2	A	301	48F	C4'-O4'-C1'	-2.73	103.43	109.47
2	D	301	48F	C6-C5-N7	2.73	137.35	132.09
2	G	301	48F	C2-N1-C6	2.61	123.02	118.73
2	E	301	48F	C4-N9-C8	2.59	108.46	105.74
2	G	301	48F	C6-C5-N7	2.58	137.07	132.09
3	C	302	NHE	O3-S-C2	2.58	111.05	106.00
3	F	302	NHE	C5'-C6'-C1'	2.55	115.66	111.09
2	H	301	48F	C5-N7-C8	2.54	107.44	103.45
2	E	301	48F	C2-N1-C6	2.52	122.88	118.73
2	D	301	48F	N9-C8-N7	-2.51	110.37	113.94
2	F	301	48F	C6-C5-N7	2.51	136.93	132.09
2	F	301	48F	O4D-C4D-C3D	-2.51	106.03	109.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	48F	C5-N7-C8	2.51	107.39	103.45
2	C	301	48F	C6-C5-N7	2.49	136.90	132.09
2	F	301	48F	C2-N1-C6	2.48	122.81	118.73
2	B	301	48F	C4-N9-C8	2.47	108.33	105.74
2	A	301	48F	C6-C5-N7	2.41	136.73	132.09
2	H	301	48F	C2-N1-C6	2.40	122.67	118.73
2	G	301	48F	C2'-C3'-C4'	2.37	107.20	102.61
2	B	301	48F	C4-C5-N7	-2.36	107.88	110.58
2	F	301	48F	O1A-PA-O3A	2.36	113.66	107.27
3	H	302	NHE	O3-S-C2	2.36	110.62	106.00
2	C	301	48F	N9-C8-N7	-2.35	110.61	113.94
2	H	301	48F	N9-C8-N7	-2.35	110.61	113.94
2	A	301	48F	C2-N1-C6	2.26	122.45	118.73
2	B	301	48F	N6-C6-N1	2.26	123.40	118.38
2	A	301	48F	N9-C8-N7	-2.25	110.74	113.94
3	A	302	NHE	O2-S-C2	2.24	110.12	106.73
2	C	301	48F	O3A-PB-O1B	-2.22	104.02	110.70
2	F	301	48F	C5-N7-C8	2.22	106.93	103.45
2	G	301	48F	C5-N7-C8	2.17	106.86	103.45
2	B	301	48F	C2-N1-C6	2.11	122.19	118.73
2	D	301	48F	O3A-PB-O1B	-2.10	104.38	110.70
2	A	301	48F	O4'-C1'-C2'	-2.09	102.13	106.62
2	F	301	48F	C2'-C3'-C4'	2.08	106.63	102.61
2	H	301	48F	C6-C5-N7	2.07	136.08	132.09
2	B	301	48F	C5-C6-N6	-2.03	118.26	123.29
2	C	301	48F	O4D-C4D-C3D	-2.02	106.75	109.67
2	F	301	48F	C4-N9-C8	2.00	107.84	105.74

There are no chirality outliers.

All (69) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	48F	O1D-C1D-C2D-O2D
2	B	301	48F	O1D-C1D-C2D-O2D
2	C	301	48F	O1D-C1D-C2D-O2D
2	D	301	48F	C5'-O5'-PA-O2A
2	E	301	48F	C1D-C2D-C3D-C4D
2	E	301	48F	O2D-C2D-C3D-C4D
2	E	301	48F	O2D-C2D-C3D-O3D
2	E	301	48F	O1D-C1D-C2D-O2D
2	F	301	48F	O1D-C1D-C2D-C3D
2	F	301	48F	O1D-C1D-C2D-O2D

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Mol	Chain	Res	Type	Atoms
2	G	301	48F	O1D-C1D-C2D-C3D
2	G	301	48F	O1D-C1D-C2D-O2D
2	H	301	48F	PA-O3A-PB-O5D
2	H	301	48F	O1D-C1D-C2D-O2D
3	A	302	NHE	C6'-C1'-N-C1
3	A	302	NHE	C1-C2-S-O2
3	A	302	NHE	C1-C2-S-O3
3	C	302	NHE	C6'-C1'-N-C1
3	C	302	NHE	C1-C2-S-O2
3	F	302	NHE	C2-C1-N-C1'
3	F	302	NHE	N-C1-C2-S
3	G	302	NHE	C2'-C1'-N-C1
3	H	302	NHE	C2-C1-N-C1'
2	D	301	48F	O4'-C4'-C5'-O5'
2	D	301	48F	C3'-C4'-C5'-O5'
3	C	302	NHE	C1-C2-S-O3
3	F	302	NHE	C1-C2-S-O3
3	G	302	NHE	C2-C1-N-C1'
3	F	302	NHE	C6'-C1'-N-C1
2	B	301	48F	PA-O3A-PB-O5D
2	C	301	48F	PA-O3A-PB-O5D
2	D	301	48F	PA-O3A-PB-O5D
2	E	301	48F	PA-O3A-PB-O5D
2	F	301	48F	PA-O3A-PB-O5D
2	G	301	48F	PA-O3A-PB-O5D
2	E	301	48F	O3D-C3D-C4D-C5D
3	A	302	NHE	C1-C2-S-O1
3	C	302	NHE	C1-C2-S-O1
3	F	302	NHE	C1-C2-S-O1
3	F	302	NHE	C1-C2-S-O2
2	F	301	48F	PB-O3A-PA-O1A
2	E	301	48F	C2D-C3D-C4D-C5D
2	B	301	48F	C2D-C3D-C4D-O4D
2	D	301	48F	C5'-O5'-PA-O1A
2	D	301	48F	C5'-O5'-PA-O3A
2	E	301	48F	C2D-C3D-C4D-O4D
2	D	301	48F	O1D-C1D-C2D-O2D
2	E	301	48F	PB-O3A-PA-O2A
2	A	301	48F	O1D-C1D-C2D-C3D
2	B	301	48F	O1D-C1D-C2D-C3D
2	C	301	48F	O1D-C1D-C2D-C3D
2	E	301	48F	O1D-C1D-C2D-C3D

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Mol	Chain	Res	Type	Atoms
2	H	301	48F	O1D-C1D-C2D-C3D
2	B	301	48F	O4'-C4'-C5'-O5'
3	C	302	NHE	C2'-C1'-N-C1
3	F	302	NHE	C2'-C1'-N-C1
2	E	301	48F	O3D-C3D-C4D-O4D
2	G	301	48F	PB-O3A-PA-O2A
2	A	301	48F	PA-O3A-PB-O5D
2	B	301	48F	O3D-C3D-C4D-C5D
2	E	301	48F	C1D-C2D-C3D-O3D
2	D	301	48F	PA-O3A-PB-O2B
2	F	301	48F	PB-O3A-PA-O2A
2	E	301	48F	O4'-C4'-C5'-O5'
2	B	301	48F	PB-O3A-PA-O1A
2	C	301	48F	PB-O3A-PA-O1A
2	E	301	48F	PB-O3A-PA-O1A
2	G	301	48F	PB-O3A-PA-O1A
2	B	301	48F	C3'-C4'-C5'-O5'

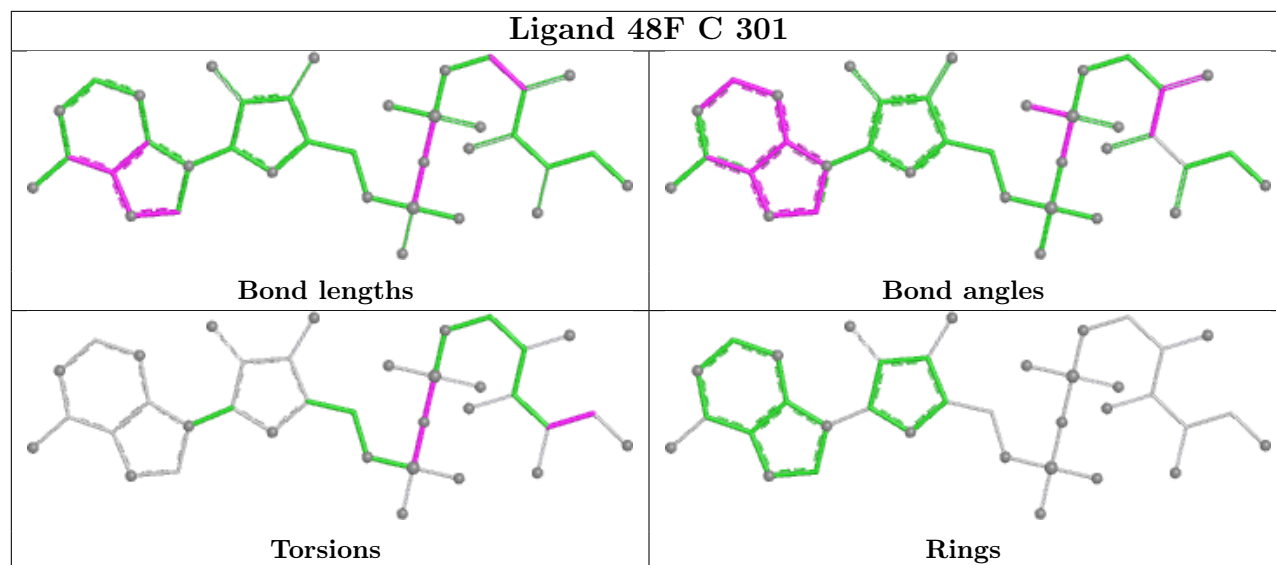
There are no ring outliers.

1 monomer is involved in 1 short contact:

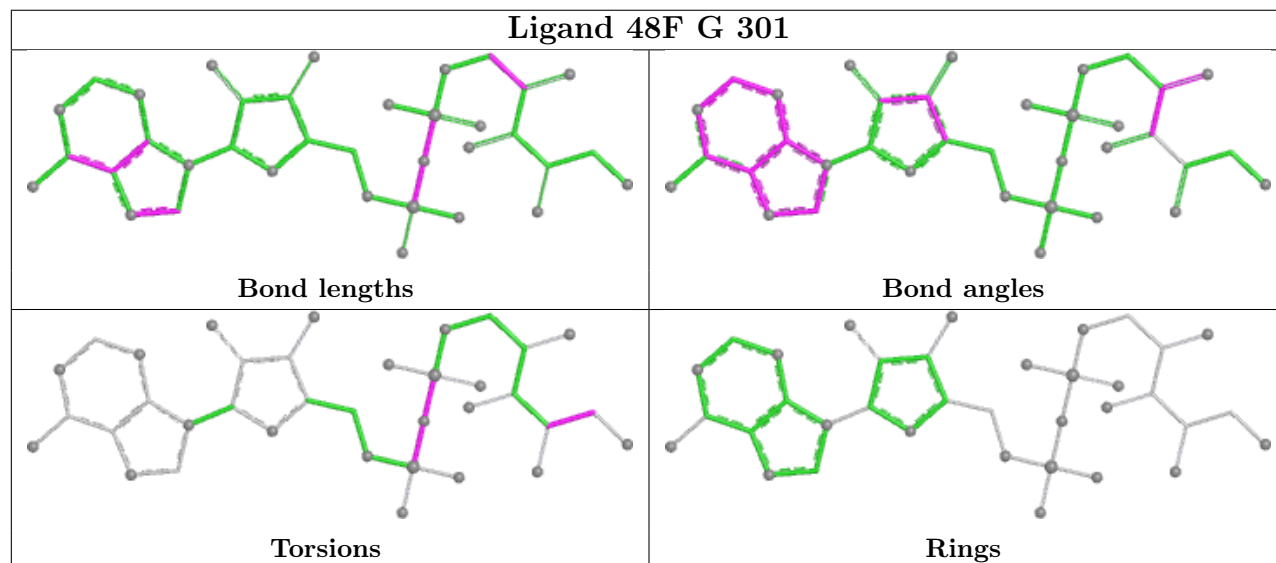
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	48F	1	0

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.

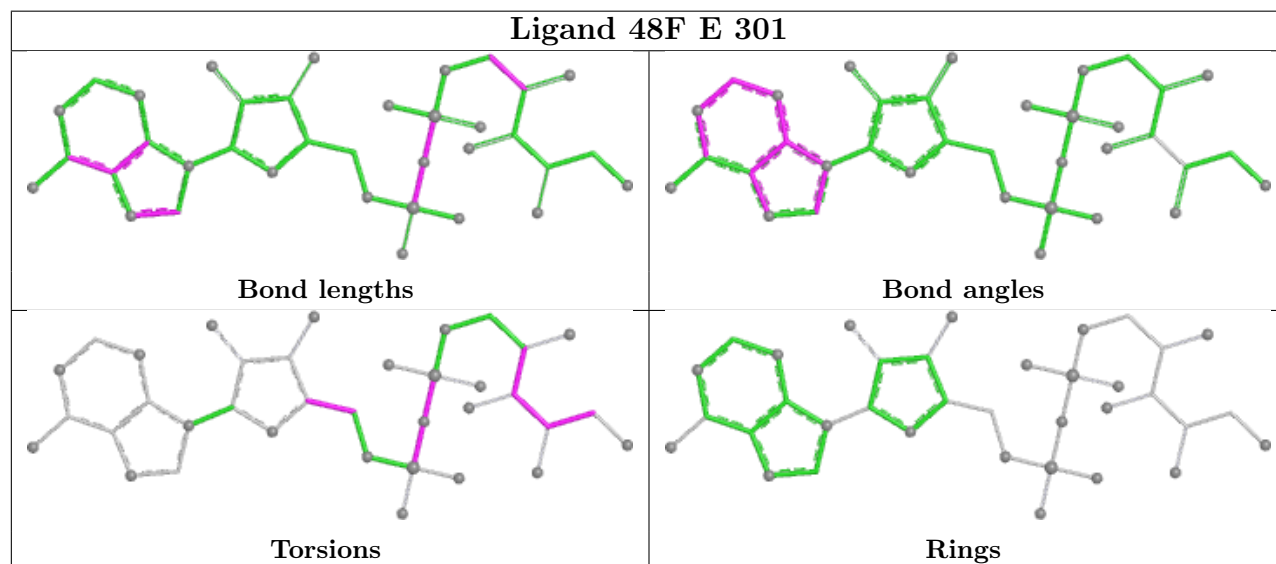
Ligand 48F C 301



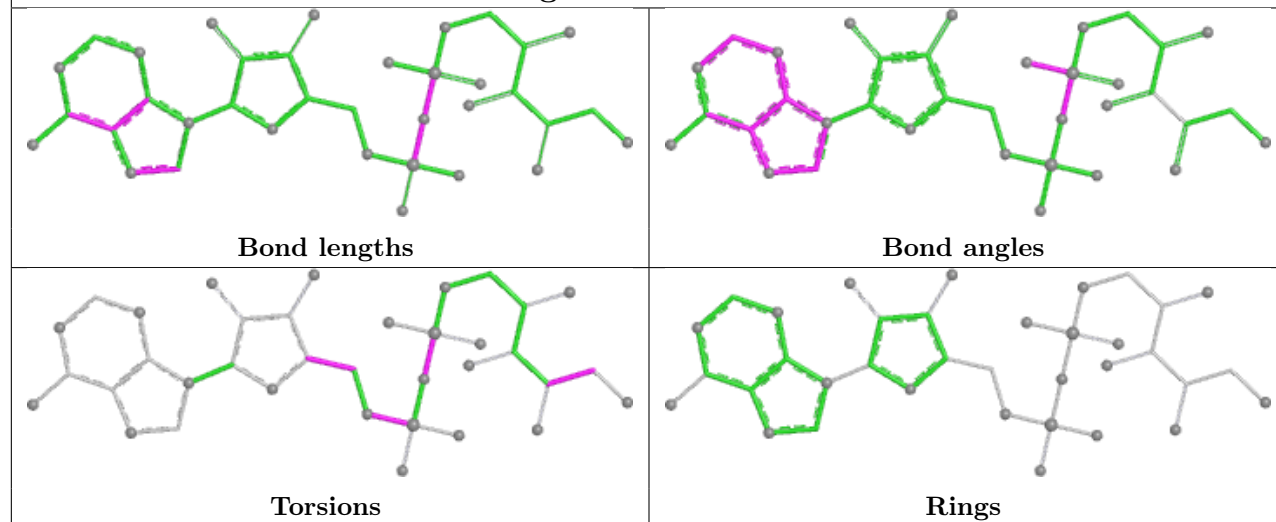
Ligand 48F G 301



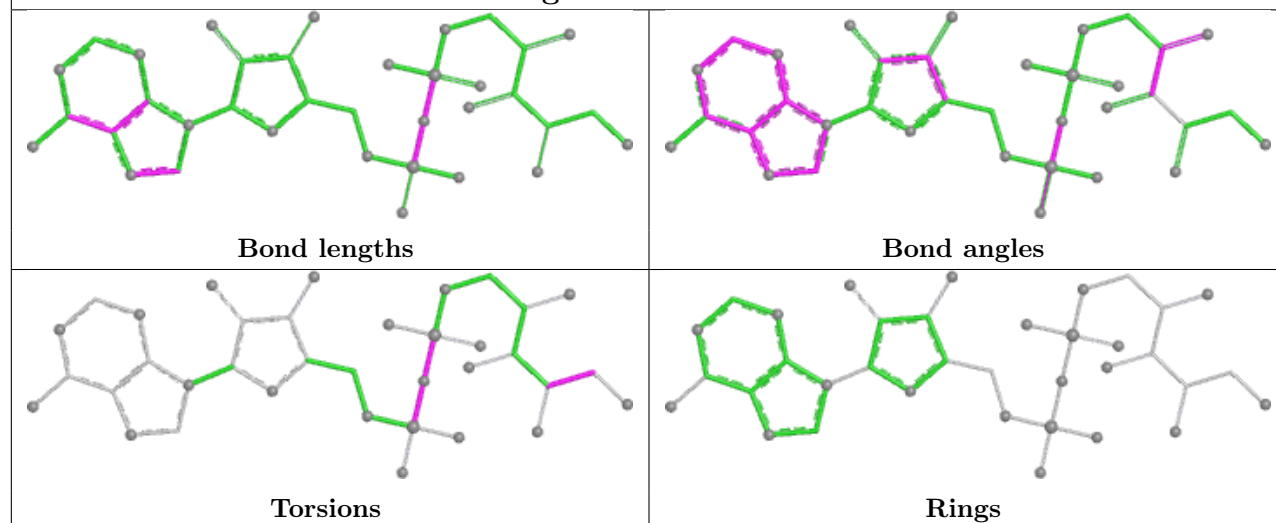
Ligand 48F E 301



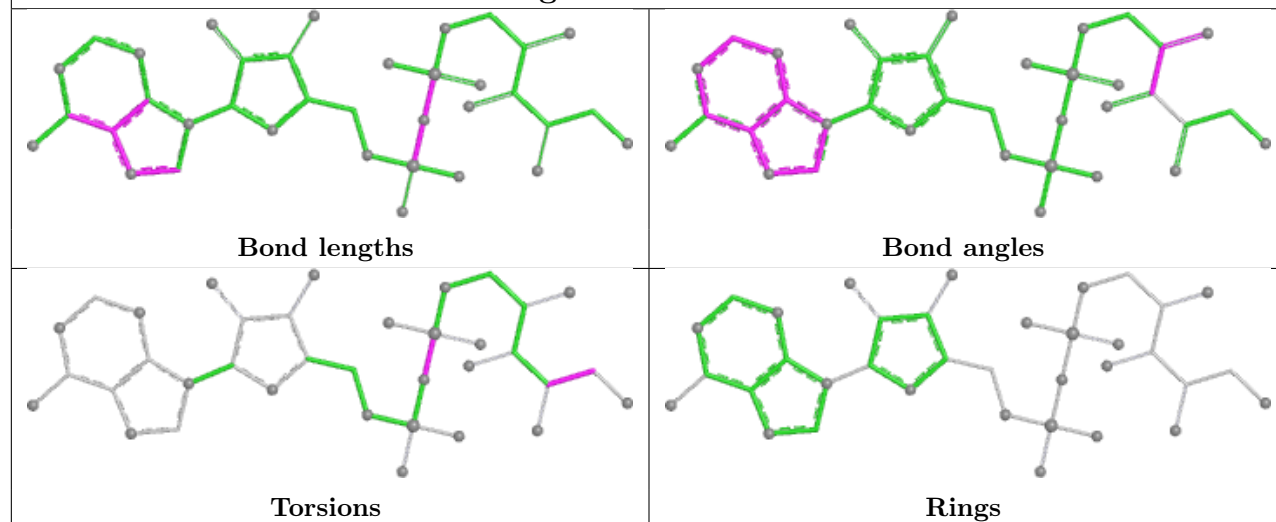
Ligand 48F D 301

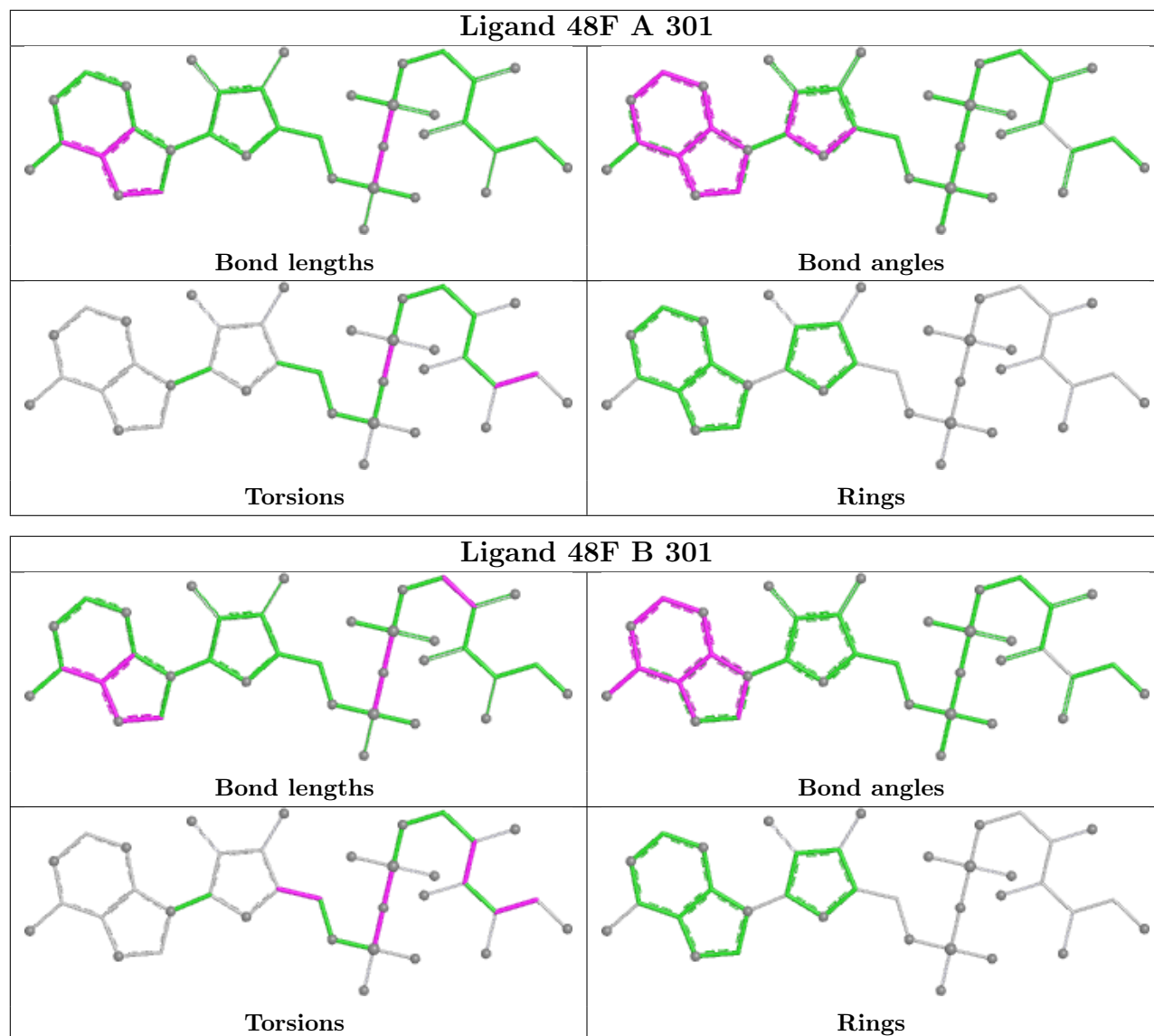


Ligand 48F F 301



Ligand 48F H 301





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	260/290 (89%)	-1.60	0 100 100	39, 55, 80, 121	0
1	B	256/290 (88%)	-1.67	0 100 100	37, 47, 70, 116	0
1	C	255/290 (87%)	-1.60	0 100 100	40, 54, 80, 143	0
1	D	255/290 (87%)	-1.59	0 100 100	37, 55, 81, 139	0
1	E	256/290 (88%)	-1.65	0 100 100	37, 47, 71, 114	0
1	F	262/290 (90%)	-1.63	0 100 100	37, 51, 72, 115	0
1	G	262/290 (90%)	-1.65	0 100 100	37, 50, 73, 115	0
1	H	259/290 (89%)	-1.60	0 100 100	39, 55, 79, 120	0
All	All	2065/2320 (89%)	-1.62	0 100 100	37, 52, 78, 143	0

There are no RSRZ outliers to report.

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](#)

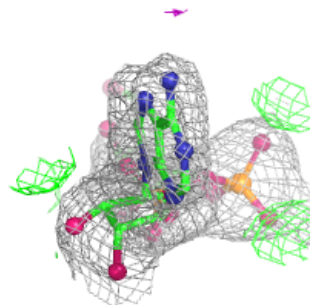
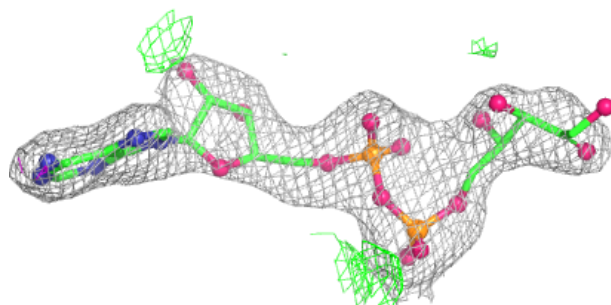
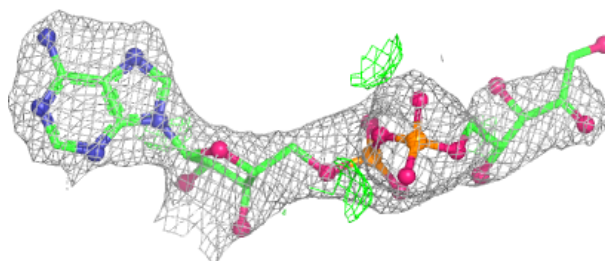
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NHE	C	302	13/13	0.98	0.10	108,115,122,131	0
2	48F	B	301	36/36	0.99	0.04	50,63,88,101	0
2	48F	C	301	36/36	0.99	0.04	56,68,79,84	0
2	48F	D	301	36/36	0.99	0.04	56,68,79,87	0
2	48F	E	301	36/36	0.99	0.04	49,64,89,106	0
2	48F	F	301	36/36	0.99	0.03	45,58,78,96	0
2	48F	G	301	36/36	0.99	0.03	44,57,75,91	0
2	48F	H	301	36/36	0.99	0.04	52,61,80,81	0
3	NHE	A	302	13/13	0.99	0.05	94,99,116,119	0
2	48F	A	301	36/36	0.99	0.04	47,61,78,81	0
3	NHE	F	302	13/13	0.99	0.04	69,72,81,83	0
3	NHE	H	302	13/13	0.99	0.05	89,94,111,112	0
3	NHE	G	302	13/13	1.00	0.03	64,75,83,87	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.

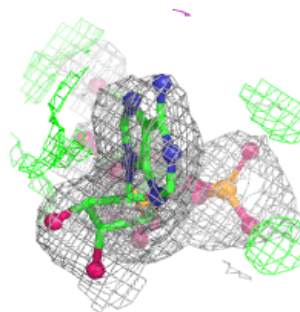
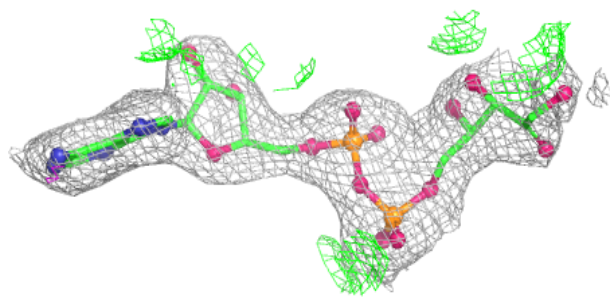
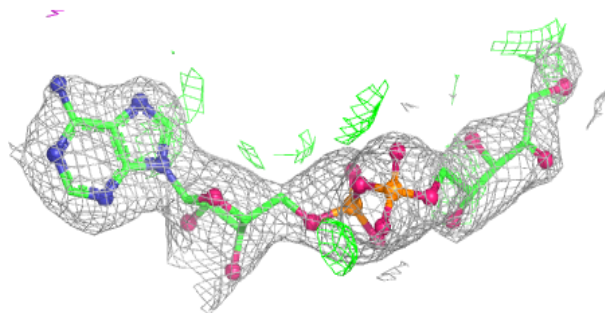
Electron density around 48F B 301:

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

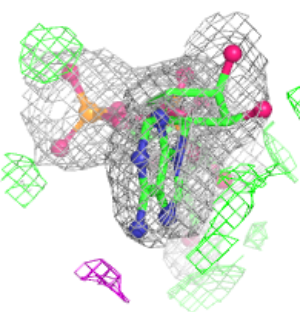
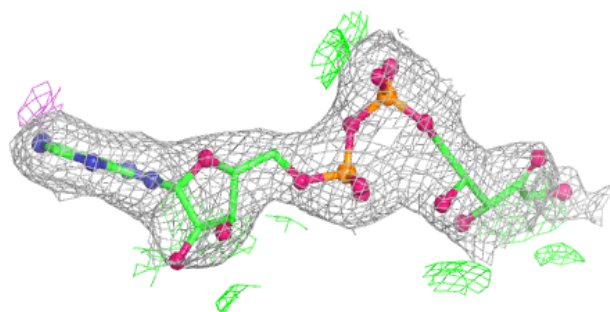
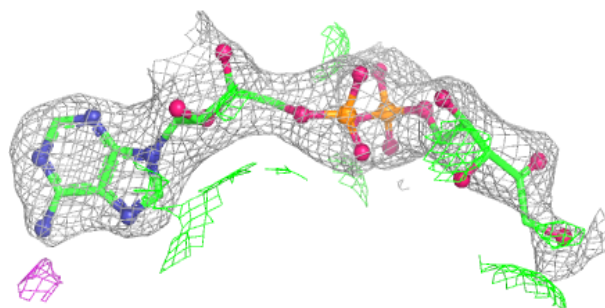


Electron density around 48F C 301:

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

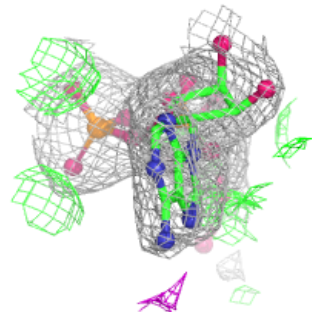
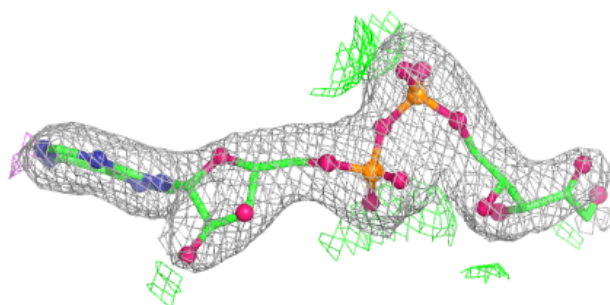
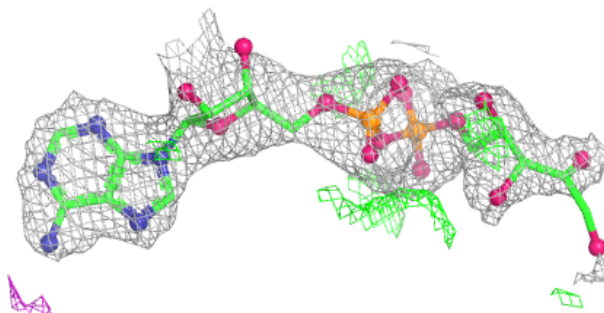
**Electron density around 48F D 301:**

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

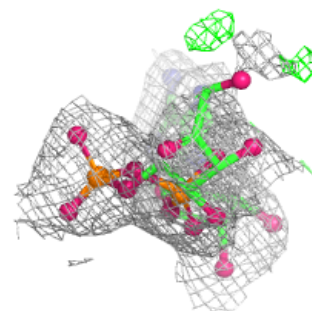
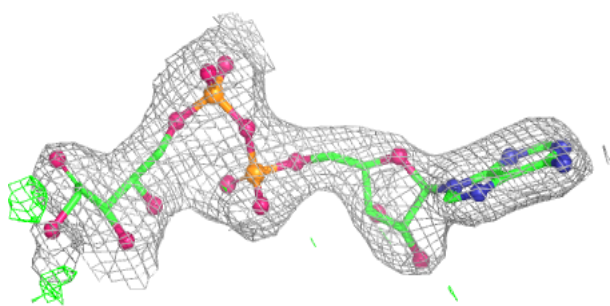
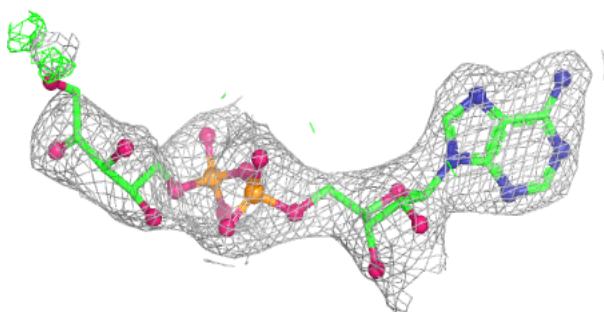


Electron density around 48F E 301:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

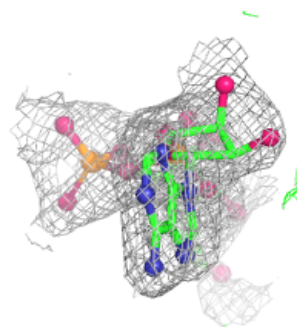
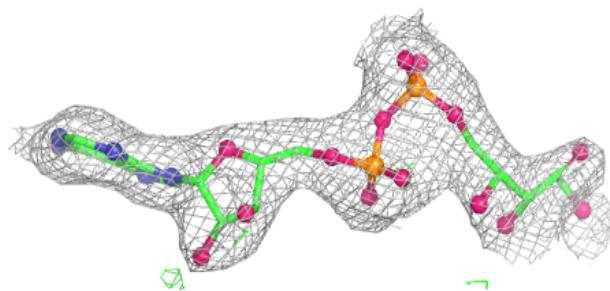
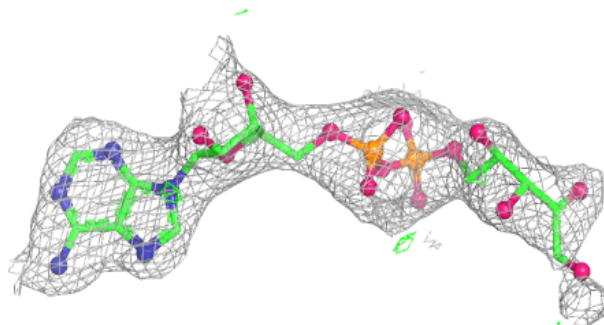
**Electron density around 48F F 301:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

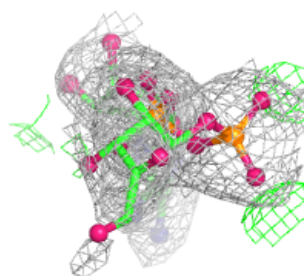
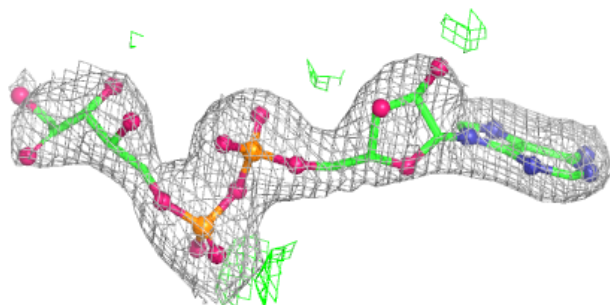
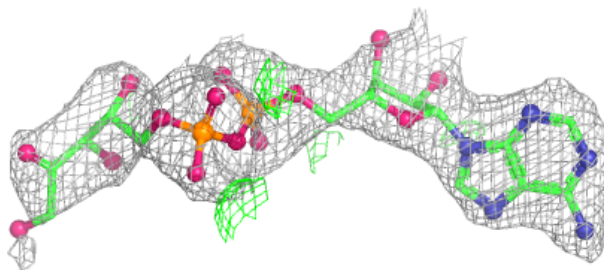


Electron density around 48F G 301:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

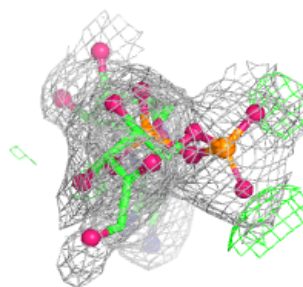
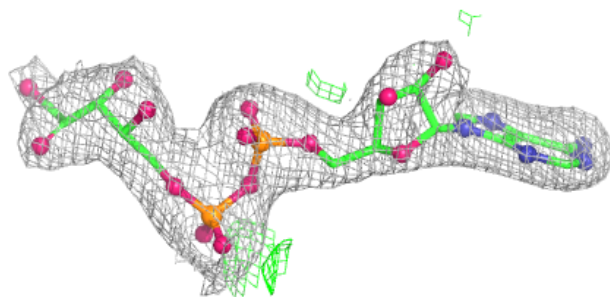
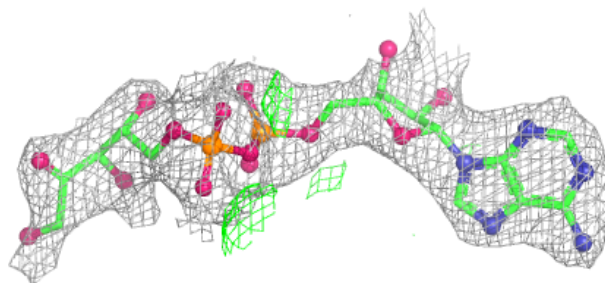
**Electron density around 48F H 301:**

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



Electron density around 48F A 301:

$2mF_o - DF_c$ (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.