



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 2, 2022 – 08:05 AM EDT

PDB ID : 8CSO  
Title : Crystal Structure of Orotidine 5'-phosphate decarboxylase from *Klebsiella pneumoniae* in complex with Uridine-5'-monophosphate  
Authors : Seattle Structural Genomics Center for Infectious Disease; Seattle Structural Genomics Center for Infectious Disease (SSGCID)  
Deposited on : 2022-05-13  
Resolution : 2.60 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.28.1  
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) : 2.28.1

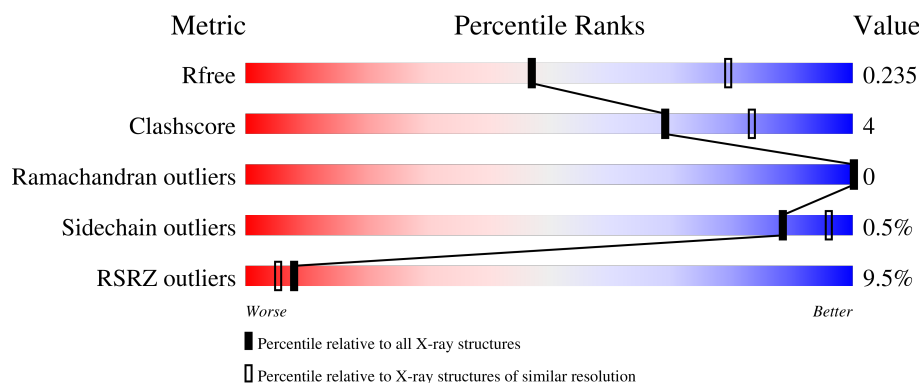
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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  $\geq 3$ , 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	236	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">%</span> </div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>84%</span> <span>12%</span> <span>.</span> </div> </div>
1	B	236	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">%</span> </div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>88%</span> <span>9%</span> <span>.</span> </div> </div>
1	C	236	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">%</span> </div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>90%</span> <span>8%</span> <span>.</span> </div> </div>
1	D	236	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">%</span> </div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>86%</span> <span>8%</span> <span>6%</span> </div> </div>
1	E	236	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">%</span> </div> <div style="display: flex; justify-content: space-between; align-items: center;"> <span>87%</span> <span>11%</span> <span>.</span> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	236	<div><div></div><div>89%</div><div>8%</div><div></div></div>
1	G	236	<div><div>22%</div><div>82%</div><div>6%</div><div>13%</div></div>
1	H	236	<div><div>15%</div><div>84%</div><div>6%</div><div>10%</div></div>
1	I	236	<div><div>26%</div><div>80%</div><div>12%</div><div>8%</div></div>
1	J	236	<div><div>22%</div><div>70%</div><div>8%</div><div>22%</div></div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16033 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 Orotidine 5'-phosphate decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	229	Total	C	N	O	S	0	0	0
			1712	1075	305	321	11			
1	B	229	Total	C	N	O	S	0	0	0
			1703	1069	306	317	11			
1	C	231	Total	C	N	O	S	0	0	0
			1730	1085	309	325	11			
1	D	221	Total	C	N	O	S	0	0	0
			1653	1043	293	306	11			
1	E	231	Total	C	N	O	S	0	0	0
			1709	1074	304	320	11			
1	F	231	Total	C	N	O	S	0	0	0
			1713	1077	306	319	11			
1	G	206	Total	C	N	O	S	0	0	0
			1423	907	248	258	10			
1	H	212	Total	C	N	O	S	0	0	0
			1445	921	247	267	10			
1	I	217	Total	C	N	O	S	0	0	0
			1449	915	255	269	10			
1	J	184	Total	C	N	O	S	0	0	0
			1152	718	204	221	9			

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP A0A0H3GNE4
A	-6	ALA	-	expression tag	UNP A0A0H3GNE4
A	-5	HIS	-	expression tag	UNP A0A0H3GNE4
A	-4	HIS	-	expression tag	UNP A0A0H3GNE4
A	-3	HIS	-	expression tag	UNP A0A0H3GNE4
A	-2	HIS	-	expression tag	UNP A0A0H3GNE4
A	-1	HIS	-	expression tag	UNP A0A0H3GNE4
A	0	MET	-	expression tag	UNP A0A0H3GNE4
A	1	VAL	-	expression tag	UNP A0A0H3GNE4

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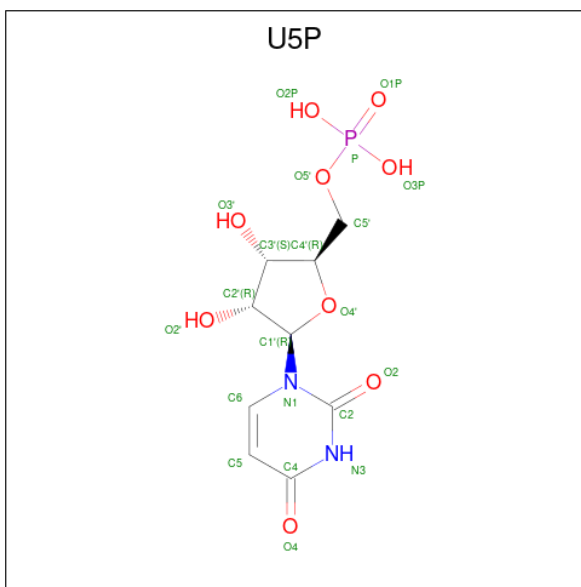
Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	MET	-	initiating methionine	UNP A0A0H3GNE4
B	-6	ALA	-	expression tag	UNP A0A0H3GNE4
B	-5	HIS	-	expression tag	UNP A0A0H3GNE4
B	-4	HIS	-	expression tag	UNP A0A0H3GNE4
B	-3	HIS	-	expression tag	UNP A0A0H3GNE4
B	-2	HIS	-	expression tag	UNP A0A0H3GNE4
B	-1	HIS	-	expression tag	UNP A0A0H3GNE4
B	0	MET	-	expression tag	UNP A0A0H3GNE4
B	1	VAL	-	expression tag	UNP A0A0H3GNE4
C	-7	MET	-	initiating methionine	UNP A0A0H3GNE4
C	-6	ALA	-	expression tag	UNP A0A0H3GNE4
C	-5	HIS	-	expression tag	UNP A0A0H3GNE4
C	-4	HIS	-	expression tag	UNP A0A0H3GNE4
C	-3	HIS	-	expression tag	UNP A0A0H3GNE4
C	-2	HIS	-	expression tag	UNP A0A0H3GNE4
C	-1	HIS	-	expression tag	UNP A0A0H3GNE4
C	0	MET	-	expression tag	UNP A0A0H3GNE4
C	1	VAL	-	expression tag	UNP A0A0H3GNE4
D	-7	MET	-	initiating methionine	UNP A0A0H3GNE4
D	-6	ALA	-	expression tag	UNP A0A0H3GNE4
D	-5	HIS	-	expression tag	UNP A0A0H3GNE4
D	-4	HIS	-	expression tag	UNP A0A0H3GNE4
D	-3	HIS	-	expression tag	UNP A0A0H3GNE4
D	-2	HIS	-	expression tag	UNP A0A0H3GNE4
D	-1	HIS	-	expression tag	UNP A0A0H3GNE4
D	0	MET	-	expression tag	UNP A0A0H3GNE4
D	1	VAL	-	expression tag	UNP A0A0H3GNE4
E	-7	MET	-	initiating methionine	UNP A0A0H3GNE4
E	-6	ALA	-	expression tag	UNP A0A0H3GNE4
E	-5	HIS	-	expression tag	UNP A0A0H3GNE4
E	-4	HIS	-	expression tag	UNP A0A0H3GNE4
E	-3	HIS	-	expression tag	UNP A0A0H3GNE4
E	-2	HIS	-	expression tag	UNP A0A0H3GNE4
E	-1	HIS	-	expression tag	UNP A0A0H3GNE4
E	0	MET	-	expression tag	UNP A0A0H3GNE4
E	1	VAL	-	expression tag	UNP A0A0H3GNE4
F	-7	MET	-	initiating methionine	UNP A0A0H3GNE4
F	-6	ALA	-	expression tag	UNP A0A0H3GNE4
F	-5	HIS	-	expression tag	UNP A0A0H3GNE4
F	-4	HIS	-	expression tag	UNP A0A0H3GNE4
F	-3	HIS	-	expression tag	UNP A0A0H3GNE4
F	-2	HIS	-	expression tag	UNP A0A0H3GNE4

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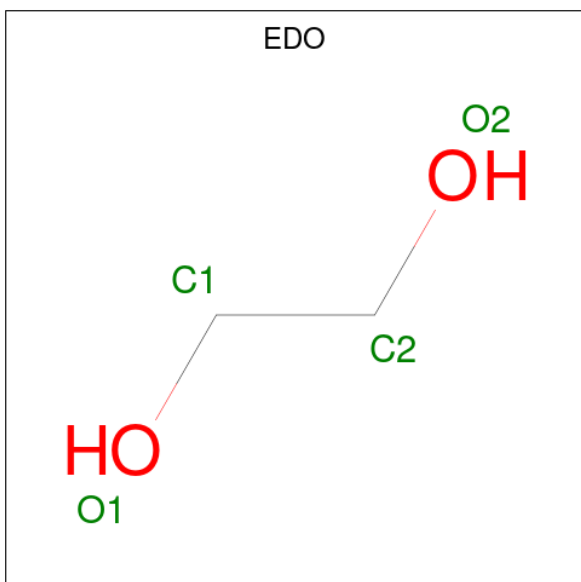
Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	HIS	-	expression tag	UNP A0A0H3GNE4
F	0	MET	-	expression tag	UNP A0A0H3GNE4
F	1	VAL	-	expression tag	UNP A0A0H3GNE4
G	-7	MET	-	initiating methionine	UNP A0A0H3GNE4
G	-6	ALA	-	expression tag	UNP A0A0H3GNE4
G	-5	HIS	-	expression tag	UNP A0A0H3GNE4
G	-4	HIS	-	expression tag	UNP A0A0H3GNE4
G	-3	HIS	-	expression tag	UNP A0A0H3GNE4
G	-2	HIS	-	expression tag	UNP A0A0H3GNE4
G	-1	HIS	-	expression tag	UNP A0A0H3GNE4
G	0	MET	-	expression tag	UNP A0A0H3GNE4
G	1	VAL	-	expression tag	UNP A0A0H3GNE4
H	-7	MET	-	initiating methionine	UNP A0A0H3GNE4
H	-6	ALA	-	expression tag	UNP A0A0H3GNE4
H	-5	HIS	-	expression tag	UNP A0A0H3GNE4
H	-4	HIS	-	expression tag	UNP A0A0H3GNE4
H	-3	HIS	-	expression tag	UNP A0A0H3GNE4
H	-2	HIS	-	expression tag	UNP A0A0H3GNE4
H	-1	HIS	-	expression tag	UNP A0A0H3GNE4
H	0	MET	-	expression tag	UNP A0A0H3GNE4
H	1	VAL	-	expression tag	UNP A0A0H3GNE4
I	-7	MET	-	initiating methionine	UNP A0A0H3GNE4
I	-6	ALA	-	expression tag	UNP A0A0H3GNE4
I	-5	HIS	-	expression tag	UNP A0A0H3GNE4
I	-4	HIS	-	expression tag	UNP A0A0H3GNE4
I	-3	HIS	-	expression tag	UNP A0A0H3GNE4
I	-2	HIS	-	expression tag	UNP A0A0H3GNE4
I	-1	HIS	-	expression tag	UNP A0A0H3GNE4
I	0	MET	-	expression tag	UNP A0A0H3GNE4
I	1	VAL	-	expression tag	UNP A0A0H3GNE4
J	-7	MET	-	initiating methionine	UNP A0A0H3GNE4
J	-6	ALA	-	expression tag	UNP A0A0H3GNE4
J	-5	HIS	-	expression tag	UNP A0A0H3GNE4
J	-4	HIS	-	expression tag	UNP A0A0H3GNE4
J	-3	HIS	-	expression tag	UNP A0A0H3GNE4
J	-2	HIS	-	expression tag	UNP A0A0H3GNE4
J	-1	HIS	-	expression tag	UNP A0A0H3GNE4
J	0	MET	-	expression tag	UNP A0A0H3GNE4
J	1	VAL	-	expression tag	UNP A0A0H3GNE4

- Molecule 2 is URIDINE-5'-MONOPHOSPHATE (three-letter code: U5P) (formula:  $C_9H_{13}N_2O_9P$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			21	9	2	9	1		
2	B	1	Total	C	N	O	P	0	0
			21	9	2	9	1		
2	C	1	Total	C	N	O	P	0	0
			21	9	2	9	1		
2	E	1	Total	C	N	O	P	0	0
			21	9	2	9	1		
2	F	1	Total	C	N	O	P	0	0
			21	9	2	9	1		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0
3	F	1	Total C O 4 2 2	0	0

- Molecule 4 is water.

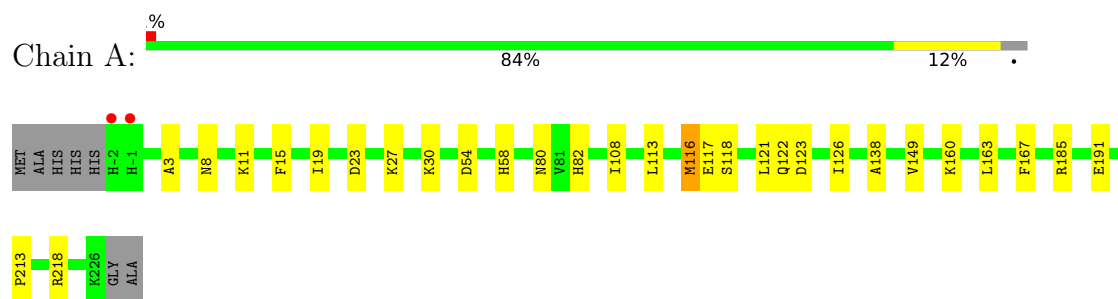
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	33	Total O 33 33	0	0
4	B	21	Total O 21 21	0	0
4	C	55	Total O 55 55	0	0
4	D	46	Total O 46 46	0	0
4	E	41	Total O 41 41	0	0
4	F	30	Total O 30 30	0	0
4	H	1	Total O 1 1	0	0



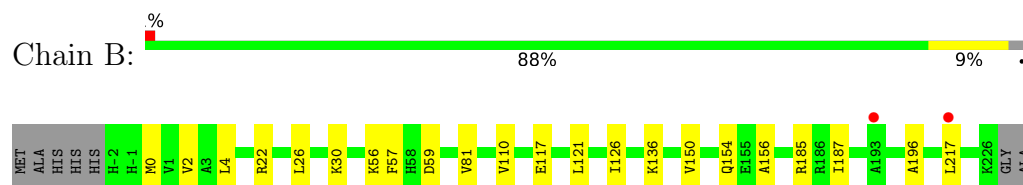
### 3 Residue-property plots [i](#)

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.

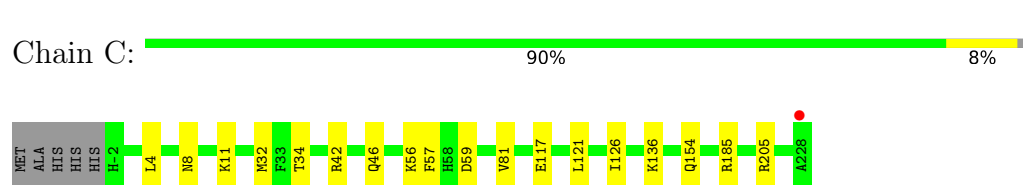
- Molecule 1: Orotidine 5'-phosphate decarboxylase



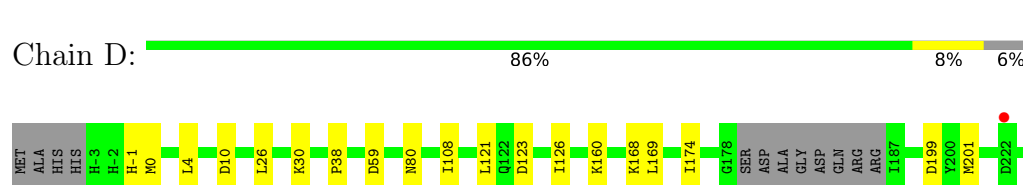
- Molecule 1: Orotidine 5'-phosphate decarboxylase



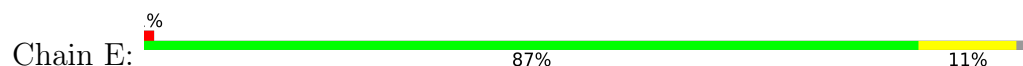
- Molecule 1: Orotidine 5'-phosphate decarboxylase



- Molecule 1: Orotidine 5'-phosphate decarboxylase



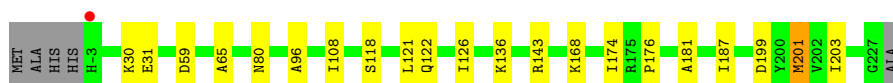
- Molecule 1: Orotidine 5'-phosphate decarboxylase





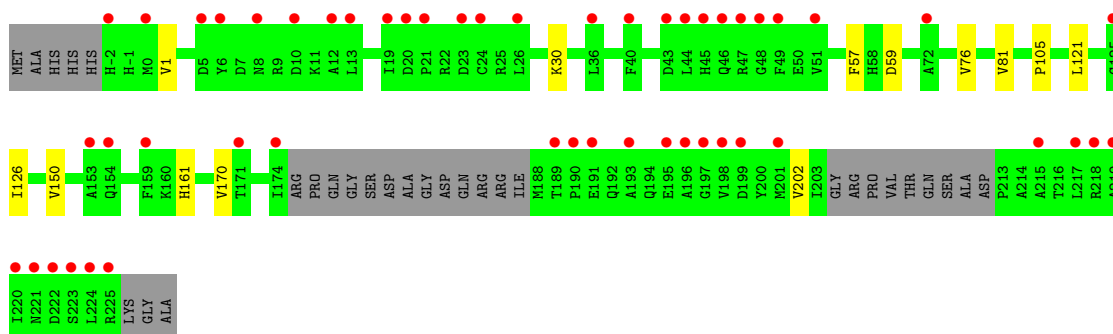
- Molecule 1: Orotidine 5'-phosphate decarboxylase

Chain F: 89% 8%



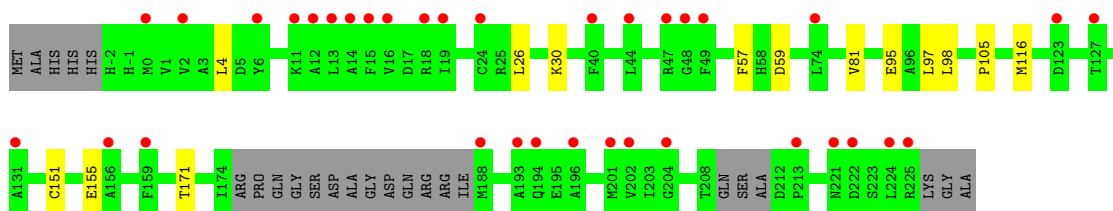
- Molecule 1: Orotidine 5'-phosphate decarboxylase

Chain G: 22% 82% 6% 13%



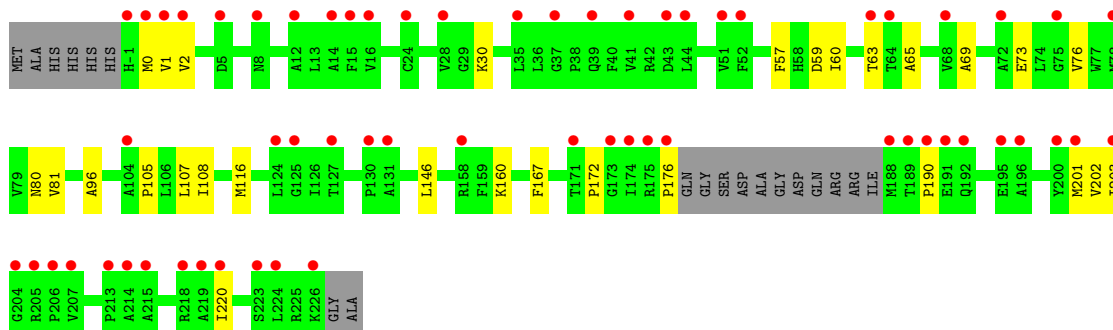
- Molecule 1: Orotidine 5'-phosphate decarboxylase

Chain H: 15% 84% 6% 10%

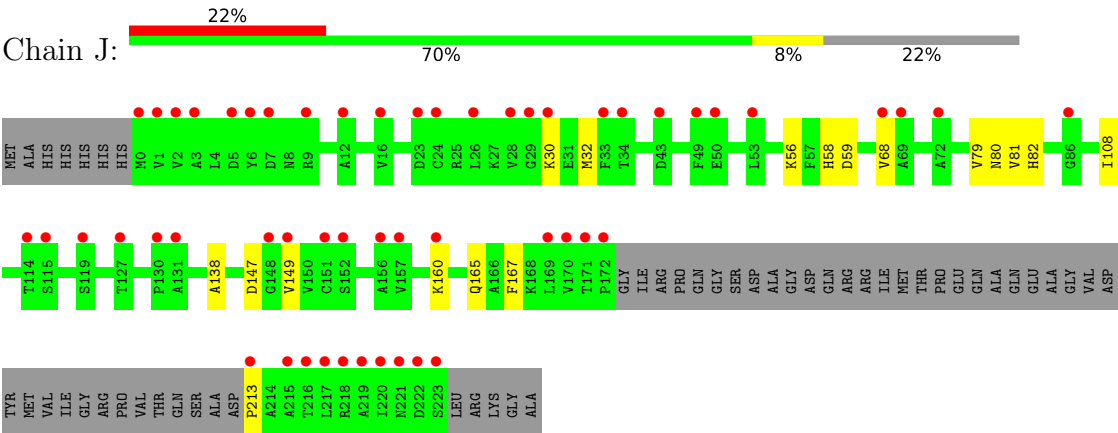


- Molecule 1: Orotidine 5'-phosphate decarboxylase

Chain I: 26% 80% 12% 8%



● Molecule 1: Orotidine 5'-phosphate decarboxylase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.71Å 91.31Å 127.73Å 100.94° 100.36° 91.11°	Depositor
Resolution (Å)	48.34 – 2.60 48.94 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.1 (48.34-2.60) 98.1 (48.94-2.60)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.18 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.20.1	Depositor
R, $R_{free}$	0.197 , 0.235 0.197 , 0.235	Depositor DCC
$R_{free}$ test set	1959 reflections (2.41%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.8	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 51.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	16033	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: U5P, EDO

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	0.33	0/1742	0.56	0/2366
1	B	0.31	0/1733	0.55	0/2356
1	C	0.35	0/1760	0.58	0/2387
1	D	0.34	0/1681	0.58	0/2282
1	E	0.32	0/1739	0.54	0/2364
1	F	0.31	0/1744	0.54	0/2370
1	G	0.26	0/1447	0.47	0/1974
1	H	0.26	0/1471	0.47	0/2014
1	I	0.27	0/1474	0.48	0/2021
1	J	0.25	0/1170	0.44	0/1607
All	All	0.31	0/15961	0.53	0/21741

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

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	1712	0	1696	20	0
1	B	1703	0	1679	15	0
1	C	1730	0	1721	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1653	0	1651	12	0
1	E	1709	0	1681	18	0
1	F	1713	0	1689	14	0
1	G	1423	0	1329	8	0
1	H	1445	0	1320	8	0
1	I	1449	0	1306	15	0
1	J	1152	0	950	11	0
2	A	21	0	11	0	0
2	B	21	0	11	1	0
2	C	21	0	11	1	0
2	E	21	0	11	1	0
2	F	21	0	11	0	0
3	B	4	0	6	0	0
3	C	4	0	6	0	0
3	F	4	0	6	0	0
4	A	33	0	0	0	0
4	B	21	0	0	0	0
4	C	55	0	0	1	0
4	D	46	0	0	0	0
4	E	41	0	0	0	0
4	F	30	0	0	1	0
4	H	1	0	0	0	0
All	All	16033	0	15095	114	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 4.

All (114) 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:E:174:ILE:HG13	1:E:201:MET:HB3	1.62	0.81
1:E:59:ASP:HB2	1:F:30:LYS:HE3	1.63	0.79
1:F:174:ILE:HG13	1:F:201:MET:HB3	1.65	0.78
1:C:154:GLN:NE2	4:C:401:HOH:O	2.22	0.71
1:E:30:LYS:HE3	1:F:59:ASP:HB2	1.75	0.68
1:F:30:LYS:NZ	4:F:401:HOH:O	2.12	0.67
1:E:66:ARG:HB3	1:F:31:GLU:HG3	1.77	0.67
1:G:59:ASP:HB2	1:H:30:LYS:HE3	1.78	0.64
1:I:30:LYS:HE3	1:J:59:ASP:HB2	1.79	0.64
1:B:154:GLN:NE2	1:B:187:ILE:O	2.30	0.63
1:I:57:PHE:HB2	1:I:81:VAL:HG12	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:174:ILE:HG13	1:D:201:MET:HB3	1.82	0.61
1:F:143:ARG:NH1	1:J:147:ASP:OD1	2.29	0.61
1:E:-3:HIS:CG	1:E:-2:HIS:H	2.21	0.59
1:I:0:MET:HE1	1:I:190:PRO:HB3	1.84	0.59
1:H:4:LEU:HD11	1:H:26:LEU:HD13	1.83	0.58
1:B:56:LYS:HZ1	2:B:301:U5P:H6	1.68	0.58
1:I:176:PRO:HG3	1:I:220:ILE:HD11	1.86	0.58
1:D:121:LEU:HB3	1:D:126:ILE:HB	1.85	0.57
1:E:123:ASP:O	1:F:136:LYS:NZ	2.37	0.57
1:H:57:PHE:HB2	1:H:81:VAL:HG12	1.87	0.57
1:F:121:LEU:HB3	1:F:126:ILE:HB	1.86	0.56
1:C:57:PHE:HB2	1:C:81:VAL:HG12	1.88	0.56
1:A:123:ASP:O	1:B:136:LYS:NZ	2.38	0.56
1:E:181:ALA:HB2	1:E:187:ILE:HD13	1.88	0.56
1:E:205:ARG:NH2	2:E:301:U5P:O2P	2.34	0.55
1:B:4:LEU:HD11	1:B:26:LEU:HD13	1.88	0.55
1:C:121:LEU:HB3	1:C:126:ILE:HB	1.88	0.55
1:I:172:PRO:HA	1:I:202:VAL:HB	1.89	0.54
1:C:117:GLU:HG3	1:C:185:ARG:HB3	1.89	0.54
1:G:30:LYS:HE3	1:H:59:ASP:HB2	1.90	0.53
1:B:121:LEU:HB3	1:B:126:ILE:HB	1.92	0.52
1:C:59:ASP:HB2	1:D:30:LYS:HE3	1.92	0.52
1:A:23:ASP:OD2	1:A:218:ARG:NH2	2.42	0.52
1:I:60:ILE:HG13	1:I:63:THR:H	1.75	0.51
1:A:58:HIS:H	1:B:30:LYS:NZ	2.08	0.51
1:A:121:LEU:HB3	1:A:126:ILE:HB	1.92	0.51
1:I:160:LYS:NZ	1:I:167:PHE:O	2.43	0.49
1:E:165:GLN:HG2	1:G:161:HIS:O	2.13	0.49
1:F:176:PRO:HD3	1:F:203:ILE:HD11	1.95	0.49
1:C:136:LYS:HE2	1:D:123:ASP:O	2.14	0.48
1:G:121:LEU:HB3	1:G:126:ILE:HB	1.95	0.48
1:C:56:LYS:NZ	1:D:59:ASP:OD2	2.40	0.48
1:H:95:GLU:HA	1:H:98:LEU:HG	1.95	0.48
1:J:68:VAL:HG13	1:J:79:VAL:HG11	1.95	0.48
1:B:156:ALA:HB3	1:B:196:ALA:HB1	1.95	0.48
1:F:118:SER:O	1:F:122:GLN:HG3	2.14	0.48
1:B:117:GLU:HG3	1:B:185:ARG:HB3	1.95	0.48
1:J:160:LYS:NZ	1:J:165:GLN:O	2.46	0.48
1:H:155:GLU:N	1:H:155:GLU:OE1	2.47	0.47
1:C:8:ASN:HB3	1:C:11:LYS:HD2	1.96	0.47
1:E:25:ARG:NE	1:E:50:GLU:OE1	2.36	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:57:PHE:HB2	1:G:81:VAL:HG12	1.97	0.47
1:A:191:GLU:OE1	1:A:191:GLU:N	2.46	0.47
1:D:4:LEU:HD11	1:D:26:LEU:HD13	1.97	0.46
1:E:-3:HIS:CG	1:E:-2:HIS:N	2.83	0.46
1:E:4:LEU:HD11	1:E:26:LEU:HD13	1.97	0.46
1:D:80:ASN:HA	1:D:108:ILE:O	2.15	0.46
1:F:65:ALA:HB1	1:F:96:ALA:HB2	1.98	0.46
1:A:27:LYS:NZ	1:A:54:ASP:OD2	2.33	0.46
1:E:76:VAL:O	1:E:105:PRO:HB3	2.16	0.45
1:E:121:LEU:HB3	1:E:126:ILE:HB	1.98	0.45
1:I:59:ASP:HB2	1:J:30:LYS:HE3	1.99	0.45
1:I:107:LEU:HG	1:I:146:LEU:HG	1.99	0.45
1:I:2:VAL:HA	1:I:203:ILE:O	2.17	0.45
1:A:118:SER:O	1:A:122:GLN:HG3	2.16	0.44
1:J:58:HIS:HB2	1:J:82:HIS:HB2	1.99	0.44
1:C:4:LEU:O	1:C:32:MET:HG3	2.17	0.44
1:A:160:LYS:HD3	1:A:160:LYS:HA	1.85	0.44
1:H:151:CYS:O	1:H:171:THR:HA	2.16	0.44
1:A:15:PHE:CD1	1:A:213:PRO:HB2	2.52	0.44
1:A:58:HIS:HB2	1:A:82:HIS:HB2	2.00	0.44
1:F:181:ALA:HB2	1:F:187:ILE:HD13	1.99	0.43
1:B:22:ARG:HE	1:B:22:ARG:HB2	1.43	0.43
1:J:149:VAL:HG13	1:J:167:PHE:CZ	2.53	0.43
1:E:80:ASN:HA	1:E:108:ILE:O	2.18	0.43
1:C:205:ARG:NH2	2:C:301:U5P:O1P	2.49	0.43
1:E:0:MET:HE1	1:E:201:MET:SD	2.59	0.43
1:G:76:VAL:O	1:G:105:PRO:HB3	2.17	0.43
1:B:0:MET:HG2	1:B:217:LEU:HD11	2.00	0.43
1:I:1:VAL:O	1:I:202:VAL:HA	2.18	0.43
1:I:69:ALA:O	1:I:73:GLU:HG2	2.18	0.43
1:B:57:PHE:HB2	1:B:81:VAL:HG12	1.99	0.43
1:A:117:GLU:HG3	1:A:185:ARG:HB3	2.01	0.43
1:A:3:ALA:HA	1:A:27:LYS:HB3	2.01	0.42
1:B:110:VAL:HA	1:B:150:VAL:HB	2.01	0.42
1:A:30:LYS:HE3	1:B:59:ASP:HB2	2.02	0.42
1:A:138:ALA:HB1	1:A:163:LEU:HD11	2.00	0.42
1:D:-1:HIS:HB2	1:D:0:MET:HE3	2.02	0.42
1:C:34:THR:O	1:D:38:PRO:HG3	2.18	0.42
1:G:1:VAL:O	1:G:202:VAL:HA	2.19	0.42
1:E:118:SER:O	1:E:122:GLN:HG3	2.20	0.42
1:I:76:VAL:O	1:I:105:PRO:HB3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:80:ASN:HA	1:I:108:ILE:O	2.19	0.42
1:F:80:ASN:HA	1:F:108:ILE:O	2.20	0.42
1:D:10:ASP:OD2	1:E:186:ARG:HA	2.20	0.41
1:J:81:VAL:N	1:J:108:ILE:O	2.45	0.41
1:A:113:LEU:HB2	1:A:116:MET:HG3	2.03	0.41
1:G:150:VAL:HA	1:G:170:VAL:O	2.20	0.41
1:H:97:LEU:HD22	1:H:105:PRO:HD2	2.03	0.41
1:J:56:LYS:HB3	1:J:82:HIS:CD2	2.55	0.41
1:J:80:ASN:HA	1:J:108:ILE:O	2.21	0.41
1:A:58:HIS:H	1:B:30:LYS:HZ2	1.69	0.41
1:B:2:VAL:HG21	1:B:217:LEU:HD22	2.03	0.41
1:A:149:VAL:HG13	1:A:167:PHE:CE2	2.56	0.40
1:D:160:LYS:HE2	1:D:169:LEU:HG	2.03	0.40
1:J:138:ALA:HA	1:J:149:VAL:HG11	2.02	0.40
1:A:80:ASN:HA	1:A:108:ILE:O	2.21	0.40
1:D:168:LYS:HA	1:D:199:ASP:OD2	2.21	0.40
1:I:65:ALA:HB1	1:I:96:ALA:HB2	2.03	0.40
1:A:8:ASN:HB3	1:A:11:LYS:HD2	2.02	0.40
1:A:15:PHE:CZ	1:A:19:ILE:HD13	2.56	0.40
1:C:42:ARG:O	1:C:46:GLN:HG3	2.21	0.40
1:F:168:LYS:HA	1:F:199:ASP:OD2	2.21	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	227/236 (96%)	222 (98%)	5 (2%)	0	100	100
1	B	227/236 (96%)	220 (97%)	7 (3%)	0	100	100
1	C	229/236 (97%)	221 (96%)	8 (4%)	0	100	100
1	D	217/236 (92%)	212 (98%)	5 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	229/236 (97%)	220 (96%)	9 (4%)	0	100	100
1	F	229/236 (97%)	223 (97%)	6 (3%)	0	100	100
1	G	200/236 (85%)	196 (98%)	4 (2%)	0	100	100
1	H	206/236 (87%)	201 (98%)	5 (2%)	0	100	100
1	I	213/236 (90%)	208 (98%)	5 (2%)	0	100	100
1	J	180/236 (76%)	173 (96%)	7 (4%)	0	100	100
All	All	2157/2360 (91%)	2096 (97%)	61 (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	175/186 (94%)	174 (99%)	1 (1%)	86	95
1	B	172/186 (92%)	172 (100%)	0	100	100
1	C	177/186 (95%)	177 (100%)	0	100	100
1	D	169/186 (91%)	169 (100%)	0	100	100
1	E	172/186 (92%)	172 (100%)	0	100	100
1	F	173/186 (93%)	172 (99%)	1 (1%)	86	95
1	G	125/186 (67%)	125 (100%)	0	100	100
1	H	126/186 (68%)	125 (99%)	1 (1%)	81	92
1	I	121/186 (65%)	119 (98%)	2 (2%)	60	81
1	J	82/186 (44%)	80 (98%)	2 (2%)	49	74
All	All	1492/1860 (80%)	1485 (100%)	7 (0%)	88	96

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	116	MET

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Mol	Chain	Res	Type
1	F	201	MET
1	H	116	MET
1	I	116	MET
1	I	201	MET
1	J	32	MET
1	J	213	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	122	GLN
1	E	-3	HIS
1	E	184	GLN
1	F	-3	HIS
1	F	209	GLN
1	J	45	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

8 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	U5P	C	301	-	18,22,22	1.11	1 (5%)	21,33,33	1.19	1 (4%)
2	U5P	B	301	-	18,22,22	0.98	1 (5%)	21,33,33	1.08	1 (4%)
2	U5P	A	301	-	18,22,22	1.11	2 (11%)	21,33,33	1.22	2 (9%)
2	U5P	F	302	-	18,22,22	1.14	2 (11%)	21,33,33	1.07	1 (4%)
3	EDO	C	302	-	3,3,3	0.62	0	2,2,2	0.42	0
3	EDO	B	302	-	3,3,3	0.56	0	2,2,2	0.27	0
3	EDO	F	301	-	3,3,3	0.55	0	2,2,2	0.29	0
2	U5P	E	301	-	18,22,22	1.08	2 (11%)	21,33,33	1.16	2 (9%)

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	U5P	C	301	-	-	2/8/26/26	0/2/2/2
2	U5P	B	301	-	-	4/8/26/26	0/2/2/2
2	U5P	A	301	-	-	5/8/26/26	0/2/2/2
2	U5P	F	302	-	-	4/8/26/26	0/2/2/2
3	EDO	C	302	-	-	0/1/1/1	-
3	EDO	B	302	-	-	0/1/1/1	-
3	EDO	F	301	-	-	0/1/1/1	-
2	U5P	E	301	-	-	5/8/26/26	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	302	U5P	C4-N3	3.57	1.39	1.33
2	E	301	U5P	C4-N3	3.25	1.38	1.33
2	B	301	U5P	C4-N3	3.23	1.38	1.33
2	C	301	U5P	C4-N3	3.21	1.38	1.33
2	A	301	U5P	C4-N3	3.19	1.38	1.33
2	F	302	U5P	C6-N1	2.67	1.39	1.35
2	A	301	U5P	C6-N1	2.32	1.38	1.35
2	E	301	U5P	C6-N1	2.24	1.38	1.35

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	U5P	C5-C4-N3	-3.96	114.60	123.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	301	U5P	C5-C4-N3	-3.92	114.69	123.31
2	B	301	U5P	C5-C4-N3	-3.88	114.78	123.31
2	A	301	U5P	C5-C4-N3	-3.86	114.81	123.31
2	F	302	U5P	C5-C4-N3	-3.76	115.03	123.31
2	A	301	U5P	O3P-P-O2P	2.96	118.96	107.64
2	E	301	U5P	O3P-P-O2P	2.25	116.22	107.64

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	U5P	C2'-C1'-N1-C6
2	A	301	U5P	C5'-O5'-P-O3P
2	B	301	U5P	C2'-C1'-N1-C6
2	B	301	U5P	C5'-O5'-P-O1P
2	C	301	U5P	C2'-C1'-N1-C6
2	E	301	U5P	C2'-C1'-N1-C6
2	E	301	U5P	C5'-O5'-P-O3P
2	F	302	U5P	C2'-C1'-N1-C6
2	F	302	U5P	O4'-C1'-N1-C6
2	A	301	U5P	C5'-O5'-P-O1P
2	E	301	U5P	C5'-O5'-P-O1P
2	F	302	U5P	C5'-O5'-P-O1P
2	B	301	U5P	C5'-O5'-P-O3P
2	E	301	U5P	C5'-O5'-P-O2P
2	E	301	U5P	O4'-C4'-C5'-O5'
2	A	301	U5P	O4'-C4'-C5'-O5'
2	F	302	U5P	O4'-C4'-C5'-O5'
2	A	301	U5P	C5'-O5'-P-O2P
2	B	301	U5P	O4'-C4'-C5'-O5'
2	C	301	U5P	O4'-C4'-C5'-O5'

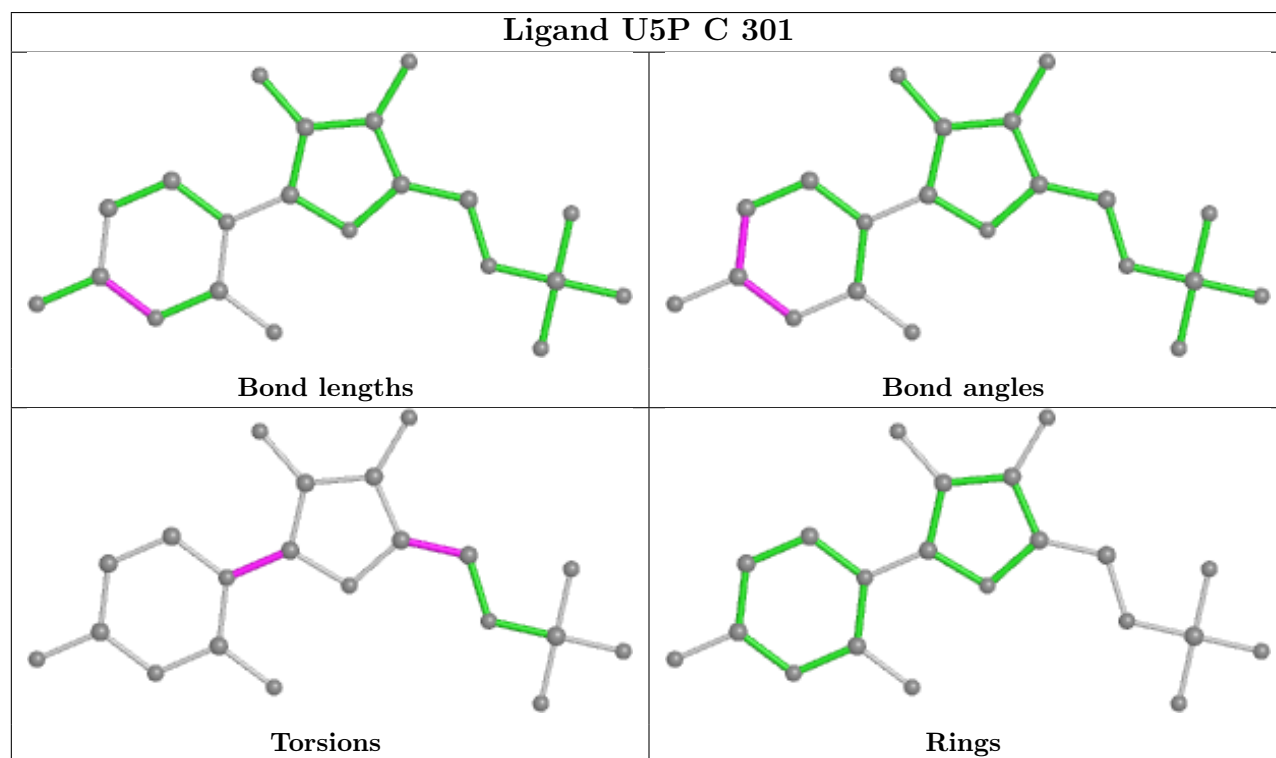
There are no ring outliers.

3 monomers are involved in 3 short contacts:

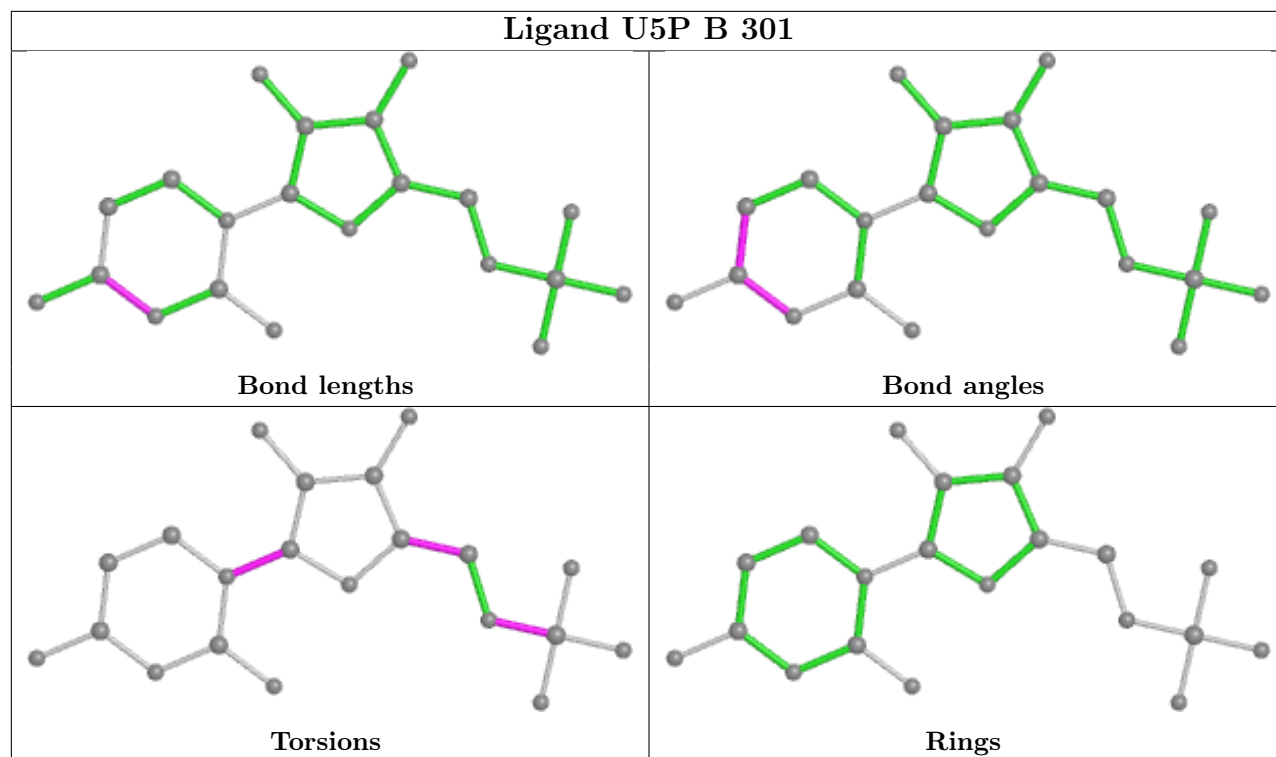
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	301	U5P	1	0
2	B	301	U5P	1	0
2	E	301	U5P	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

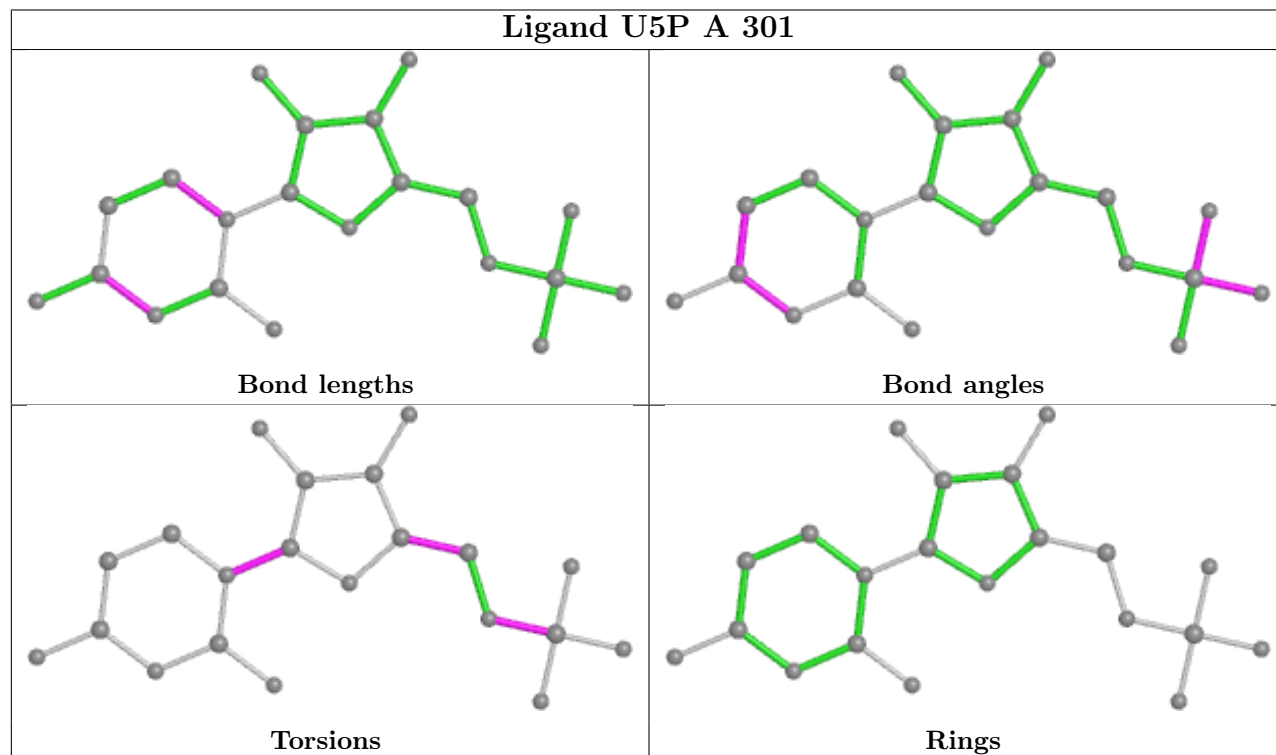
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 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.

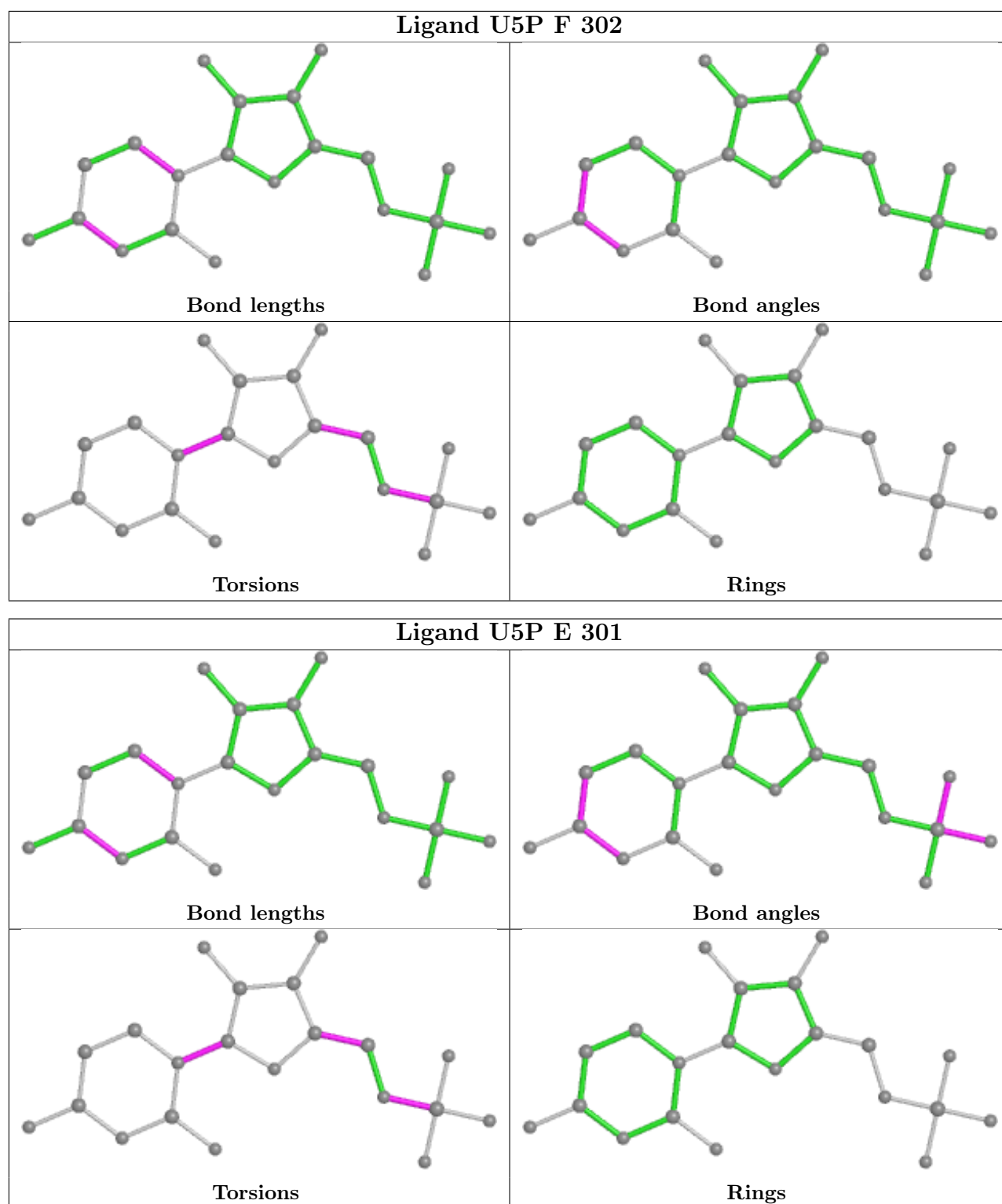


## Ligand U5P B 301



## Ligand U5P A 301





## 5.7 Other polymers ⓘ

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

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	229/236 (97%)	0.02	2 (0%) 84 82	37, 52, 81, 127	0
1	B	229/236 (97%)	0.12	2 (0%) 84 82	42, 59, 84, 125	0
1	C	231/236 (97%)	-0.03	1 (0%) 92 91	31, 40, 57, 99	0
1	D	221/236 (93%)	0.09	1 (0%) 91 89	30, 40, 75, 131	0
1	E	231/236 (97%)	0.05	2 (0%) 84 82	37, 48, 74, 130	0
1	F	231/236 (97%)	0.02	1 (0%) 92 91	38, 54, 80, 133	0
1	G	206/236 (87%)	1.14	51 (24%) 0 0	60, 92, 132, 164	0
1	H	212/236 (89%)	0.89	35 (16%) 1 1	69, 100, 130, 149	0
1	I	217/236 (91%)	1.45	61 (28%) 0 0	65, 104, 128, 155	0
1	J	184/236 (77%)	1.50	53 (28%) 0 0	81, 115, 152, 180	0
All	All	2191/2360 (92%)	0.49	209 (9%) 8 5	30, 62, 128, 180	0

All (209) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	224	LEU	7.8
1	I	196	ALA	7.2
1	I	43	ASP	6.5
1	H	14	ALA	6.5
1	G	218	ARG	6.3
1	G	219	ALA	5.9
1	G	20	ASP	5.5
1	J	220	ILE	5.3
1	J	217	LEU	5.3
1	I	16	VAL	5.2
1	I	68	VAL	5.2
1	I	220	ILE	5.1
1	J	24	CYS	5.1

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Mol	Chain	Res	Type	RSRZ
1	J	28	VAL	5.1
1	J	119	SER	5.1
1	J	223	SER	5.0
1	I	173	GLY	5.0
1	G	19	ILE	4.9
1	J	222	ASP	4.9
1	J	6	TYR	4.7
1	G	49	PHE	4.7
1	G	221	ASN	4.7
1	J	34	THR	4.7
1	I	37	GLY	4.7
1	I	64	THR	4.6
1	J	156	ALA	4.6
1	G	174	ILE	4.5
1	I	51	VAL	4.5
1	G	190	PRO	4.5
1	J	216	THR	4.5
1	J	215	ALA	4.5
1	J	12	ALA	4.4
1	J	1	VAL	4.2
1	H	156	ALA	4.2
1	J	115	SER	4.2
1	J	50	GLU	4.2
1	I	24	CYS	4.2
1	J	169	LEU	4.1
1	I	2	VAL	4.1
1	H	47	ARG	4.1
1	G	220	ILE	4.0
1	J	148	GLY	4.0
1	J	171	THR	4.0
1	J	16	VAL	4.0
1	I	191	GLU	4.0
1	G	153	ALA	4.0
1	G	215	ALA	3.9
1	H	6	TYR	3.9
1	G	222	ASP	3.9
1	I	28	VAL	3.8
1	H	225	ARG	3.8
1	J	7	ASP	3.8
1	J	170	VAL	3.8
1	G	21	PRO	3.7
1	I	192	GLN	3.7

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Mol	Chain	Res	Type	RSRZ
1	I	189	THR	3.6
1	J	69	ALA	3.5
1	I	175	ARG	3.5
1	J	3	ALA	3.5
1	I	12	ALA	3.5
1	J	151	CYS	3.4
1	J	29	GLY	3.4
1	H	213	PRO	3.4
1	J	130	PRO	3.4
1	G	23	ASP	3.4
1	I	174	ILE	3.4
1	H	2	VAL	3.4
1	J	160	LYS	3.4
1	G	191	GLU	3.4
1	I	176	PRO	3.3
1	I	200	TYR	3.3
1	H	19	ILE	3.3
1	J	49	PHE	3.3
1	I	219	ALA	3.3
1	I	130	PRO	3.3
1	I	-1	HIS	3.3
1	H	193	ALA	3.3
1	G	159	PHE	3.3
1	J	172	PRO	3.3
1	G	47	ARG	3.3
1	I	171	THR	3.2
1	H	48	GLY	3.2
1	I	8	ASN	3.2
1	I	224	LEU	3.2
1	I	215	ALA	3.2
1	I	0	MET	3.2
1	I	214	ALA	3.2
1	G	8	ASN	3.2
1	J	68	VAL	3.2
1	G	36	LEU	3.2
1	G	201	MET	3.1
1	H	16	VAL	3.1
1	G	154	GLN	3.1
1	I	44	LEU	3.1
1	G	217	LEU	3.1
1	I	52	PHE	3.1
1	G	48	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	I	75	GLY	3.1
1	G	72	ALA	3.1
1	G	196	ALA	3.1
1	J	30	LYS	3.1
1	J	152	SER	3.0
1	I	63	THR	3.0
1	B	193	ALA	3.0
1	F	-3	HIS	3.0
1	J	33	PHE	3.0
1	I	223	SER	3.0
1	I	207	VAL	3.0
1	G	46	GLN	3.0
1	H	24	CYS	3.0
1	G	223	SER	2.9
1	G	125	GLY	2.9
1	J	2	VAL	2.9
1	I	39	GLN	2.9
1	J	221	ASN	2.9
1	B	217	LEU	2.9
1	G	189	THR	2.9
1	G	10	ASP	2.9
1	I	124	LEU	2.8
1	G	6	TYR	2.8
1	H	194	GLN	2.8
1	I	78	MET	2.8
1	G	198	VAL	2.8
1	J	127	THR	2.8
1	H	18	ARG	2.8
1	G	0	MET	2.7
1	H	15	PHE	2.7
1	C	228	ALA	2.7
1	J	218	ARG	2.7
1	A	-2	HIS	2.7
1	I	205	ARG	2.7
1	J	219	ALA	2.7
1	G	193	ALA	2.6
1	J	131	ALA	2.6
1	I	226	LYS	2.6
1	I	158	ARG	2.6
1	G	24	CYS	2.6
1	J	149	VAL	2.6
1	I	206	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	G	40	PHE	2.6
1	I	5	ASP	2.6
1	I	15	PHE	2.6
1	I	14	ALA	2.6
1	E	-3	HIS	2.6
1	I	190	PRO	2.6
1	J	114	THR	2.5
1	G	12	ALA	2.5
1	G	195	GLU	2.5
1	G	5	ASP	2.5
1	I	213	PRO	2.5
1	A	-1	HIS	2.5
1	H	127	THR	2.4
1	I	104	ALA	2.4
1	G	199	ASP	2.4
1	J	9	ARG	2.4
1	I	35	LEU	2.3
1	G	45	HIS	2.3
1	H	49	PHE	2.3
1	I	1	VAL	2.3
1	I	201	MET	2.3
1	J	157	VAL	2.3
1	H	204	GLY	2.3
1	I	188	MET	2.3
1	J	213	PRO	2.3
1	H	12	ALA	2.3
1	H	131	ALA	2.3
1	I	131	ALA	2.3
1	H	123	ASP	2.3
1	G	171	THR	2.3
1	I	127	THR	2.3
1	G	43	ASP	2.3
1	G	44	LEU	2.2
1	D	222	ASP	2.2
1	J	86	GLY	2.2
1	H	221	ASN	2.2
1	E	21	PRO	2.2
1	G	225	ARG	2.2
1	G	26	LEU	2.2
1	H	11	LYS	2.2
1	I	218	ARG	2.2
1	H	44	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	I	204	GLY	2.2
1	H	201	MET	2.2
1	J	23	ASP	2.2
1	J	43	ASP	2.2
1	J	0	MET	2.2
1	J	72	ALA	2.2
1	I	203	ILE	2.1
1	I	72	ALA	2.1
1	H	13	LEU	2.1
1	J	26	LEU	2.1
1	G	197	GLY	2.1
1	H	0	MET	2.1
1	H	40	PHE	2.1
1	J	53	LEU	2.1
1	I	195	GLU	2.1
1	H	188	MET	2.1
1	H	222	ASP	2.1
1	G	51	VAL	2.1
1	H	224	LEU	2.1
1	H	159	PHE	2.1
1	I	125	GLY	2.1
1	G	13	LEU	2.1
1	H	74	LEU	2.1
1	I	41	VAL	2.1
1	J	5	ASP	2.0
1	H	202	VAL	2.0
1	H	196	ALA	2.0
1	G	-2	HIS	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no monosaccharides 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,

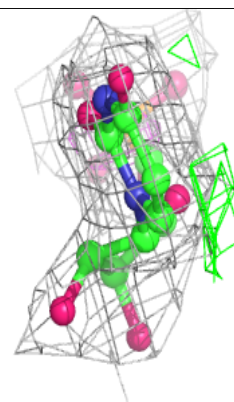
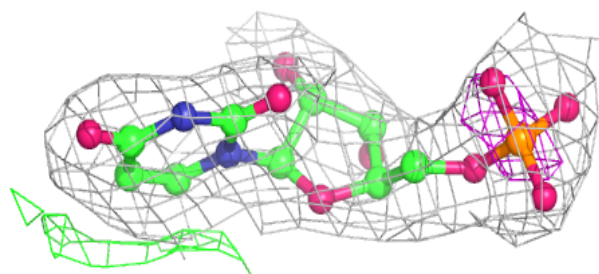
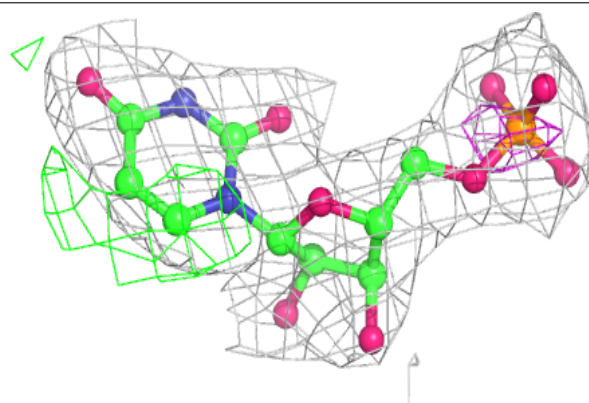
median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	EDO	F	301	4/4	0.88	0.41	40,47,47,49	0
3	EDO	C	302	4/4	0.89	0.30	38,39,40,41	0
3	EDO	B	302	4/4	0.91	0.25	43,43,44,47	0
2	U5P	E	301	21/21	0.92	0.18	50,54,58,62	0
2	U5P	F	302	21/21	0.93	0.15	56,62,66,67	0
2	U5P	B	301	21/21	0.97	0.14	47,50,53,56	0
2	U5P	A	301	21/21	0.97	0.14	43,50,55,58	0
2	U5P	C	301	21/21	0.98	0.18	30,34,35,38	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 U5P E 301:

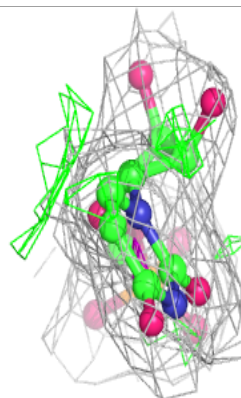
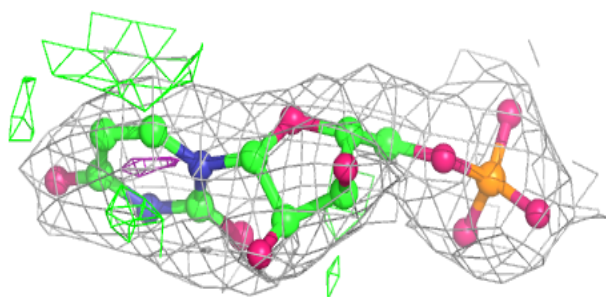
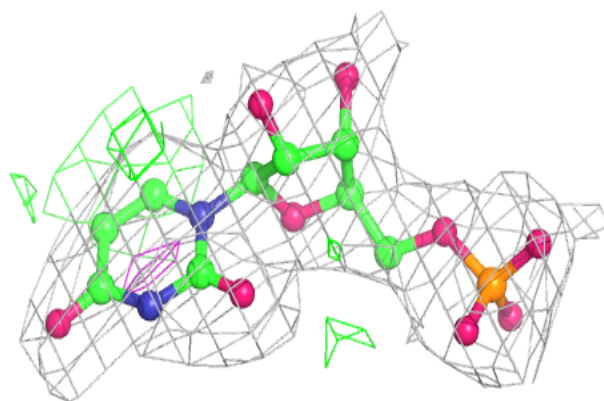
2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



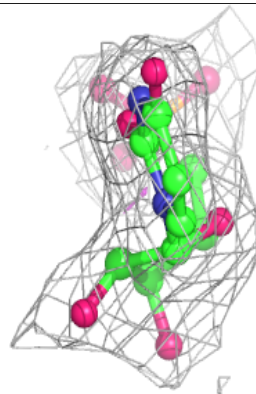
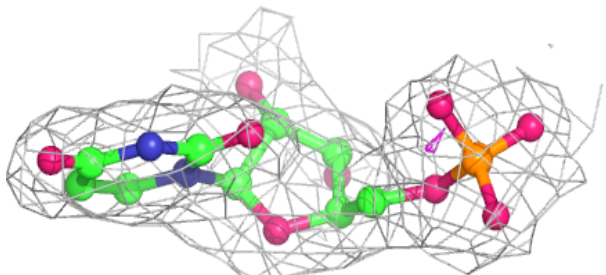
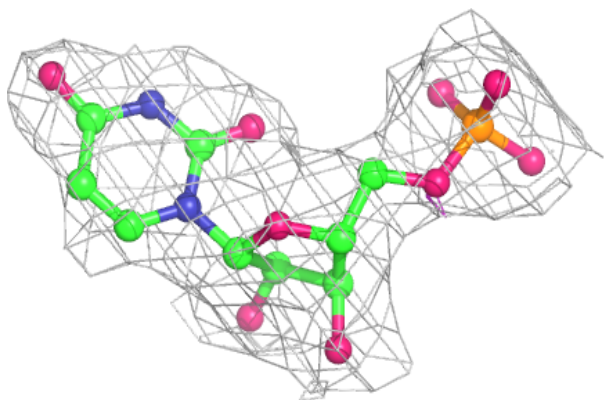


**Electron density around U5P F 302:**

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

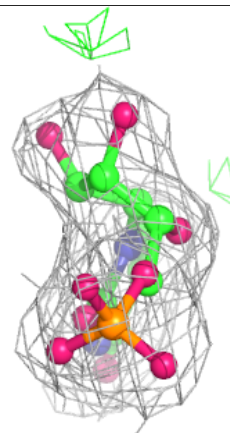
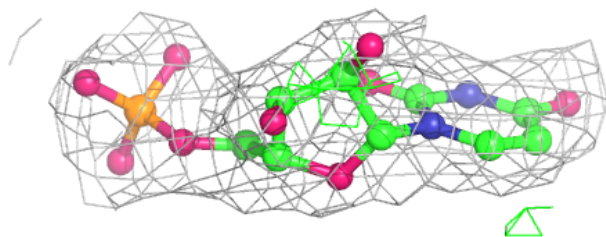
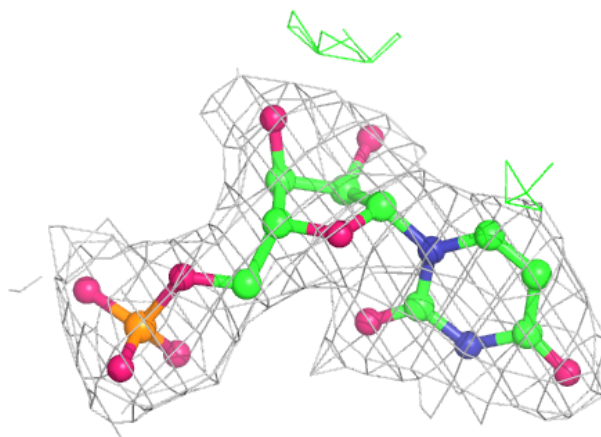
**Electron density around U5P B 301:**

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

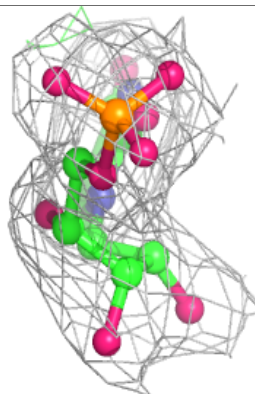
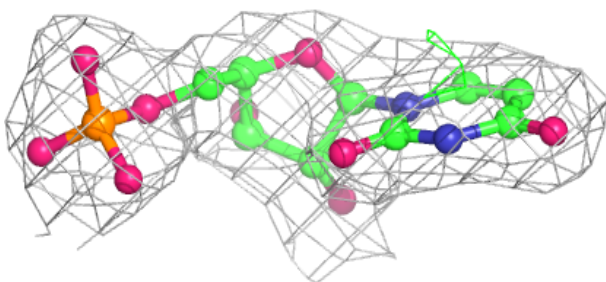
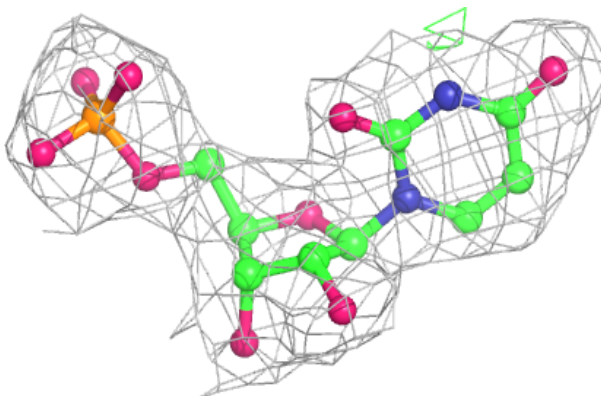


**Electron density around U5P A 301:**

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

**Electron density around U5P 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)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.