



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:14 PM UTC

PDB ID : 6MXD / pdb_00006mxd
Title : Crystal structure of Trypanosoma brucei hypoxanthine-guanine-xanthine phosphoribosyltransferase in complex with IMP
Authors : Teran, D.; Guddat, L.
Deposited on : 2018-10-30
Resolution : 2.96 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

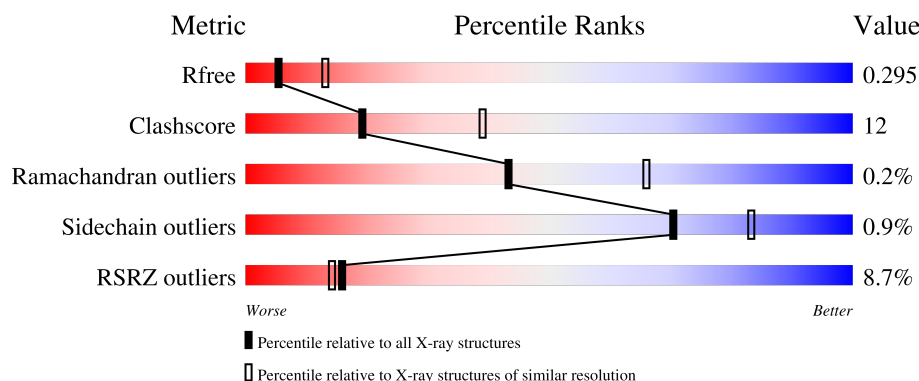
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.96 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	272	7% 58% 21% • 20%
1	B	272	10% 54% 25% 21%
1	C	272	6% 61% 19% 21%
1	D	272	5% 59% 21% 20%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6987 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 Hypoxanthine-guanine phosphoribosyltransferase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1722	1094	305	314	9			
1	B	216	Total	C	N	O	S	0	0	0
			1713	1089	303	312	9			
1	C	216	Total	C	N	O	S	0	0	0
			1715	1090	304	312	9			
1	D	217	Total	C	N	O	S	0	0	0
			1720	1093	305	313	9			

There are 156 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-37	MET	-	expression tag	UNP Q38CA1
A	-36	GLY	-	expression tag	UNP Q38CA1
A	-35	SER	-	expression tag	UNP Q38CA1
A	-34	SER	-	expression tag	UNP Q38CA1
A	-33	HIS	-	expression tag	UNP Q38CA1
A	-32	HIS	-	expression tag	UNP Q38CA1
A	-31	HIS	-	expression tag	UNP Q38CA1
A	-30	HIS	-	expression tag	UNP Q38CA1
A	-29	HIS	-	expression tag	UNP Q38CA1
A	-28	HIS	-	expression tag	UNP Q38CA1
A	-27	ASP	-	expression tag	UNP Q38CA1
A	-26	TYR	-	expression tag	UNP Q38CA1
A	-25	ASP	-	expression tag	UNP Q38CA1
A	-24	ILE	-	expression tag	UNP Q38CA1
A	-23	PRO	-	expression tag	UNP Q38CA1
A	-22	THR	-	expression tag	UNP Q38CA1
A	-21	THR	-	expression tag	UNP Q38CA1
A	-20	GLU	-	expression tag	UNP Q38CA1
A	-19	ASN	-	expression tag	UNP Q38CA1
A	-18	LEU	-	expression tag	UNP Q38CA1
A	-17	TYR	-	expression tag	UNP Q38CA1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	PHE	-	expression tag	UNP Q38CA1
A	-15	GLN	-	expression tag	UNP Q38CA1
A	-14	GLY	-	expression tag	UNP Q38CA1
A	-13	HIS	-	expression tag	UNP Q38CA1
A	-12	MET	-	expression tag	UNP Q38CA1
A	-11	ALA	-	expression tag	UNP Q38CA1
A	-10	SER	-	expression tag	UNP Q38CA1
A	-9	MET	-	expression tag	UNP Q38CA1
A	-8	THR	-	expression tag	UNP Q38CA1
A	-7	GLY	-	expression tag	UNP Q38CA1
A	-6	GLY	-	expression tag	UNP Q38CA1
A	-5	GLN	-	expression tag	UNP Q38CA1
A	-4	GLN	-	expression tag	UNP Q38CA1
A	-3	MET	-	expression tag	UNP Q38CA1
A	-2	GLY	-	expression tag	UNP Q38CA1
A	-1	ARG	-	expression tag	UNP Q38CA1
A	0	GLY	-	expression tag	UNP Q38CA1
A	1	SER	-	expression tag	UNP Q38CA1
B	-37	MET	-	expression tag	UNP Q38CA1
B	-36	GLY	-	expression tag	UNP Q38CA1
B	-35	SER	-	expression tag	UNP Q38CA1
B	-34	SER	-	expression tag	UNP Q38CA1
B	-33	HIS	-	expression tag	UNP Q38CA1
B	-32	HIS	-	expression tag	UNP Q38CA1
B	-31	HIS	-	expression tag	UNP Q38CA1
B	-30	HIS	-	expression tag	UNP Q38CA1
B	-29	HIS	-	expression tag	UNP Q38CA1
B	-28	HIS	-	expression tag	UNP Q38CA1
B	-27	ASP	-	expression tag	UNP Q38CA1
B	-26	TYR	-	expression tag	UNP Q38CA1
B	-25	ASP	-	expression tag	UNP Q38CA1
B	-24	ILE	-	expression tag	UNP Q38CA1
B	-23	PRO	-	expression tag	UNP Q38CA1
B	-22	THR	-	expression tag	UNP Q38CA1
B	-21	THR	-	expression tag	UNP Q38CA1
B	-20	GLU	-	expression tag	UNP Q38CA1
B	-19	ASN	-	expression tag	UNP Q38CA1
B	-18	LEU	-	expression tag	UNP Q38CA1
B	-17	TYR	-	expression tag	UNP Q38CA1
B	-16	PHE	-	expression tag	UNP Q38CA1
B	-15	GLN	-	expression tag	UNP Q38CA1
B	-14	GLY	-	expression tag	UNP Q38CA1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	HIS	-	expression tag	UNP Q38CA1
B	-12	MET	-	expression tag	UNP Q38CA1
B	-11	ALA	-	expression tag	UNP Q38CA1
B	-10	SER	-	expression tag	UNP Q38CA1
B	-9	MET	-	expression tag	UNP Q38CA1
B	-8	THR	-	expression tag	UNP Q38CA1
B	-7	GLY	-	expression tag	UNP Q38CA1
B	-6	GLY	-	expression tag	UNP Q38CA1
B	-5	GLN	-	expression tag	UNP Q38CA1
B	-4	GLN	-	expression tag	UNP Q38CA1
B	-3	MET	-	expression tag	UNP Q38CA1
B	-2	GLY	-	expression tag	UNP Q38CA1
B	-1	ARG	-	expression tag	UNP Q38CA1
B	0	GLY	-	expression tag	UNP Q38CA1
B	1	SER	-	expression tag	UNP Q38CA1
C	-37	MET	-	expression tag	UNP Q38CA1
C	-36	GLY	-	expression tag	UNP Q38CA1
C	-35	SER	-	expression tag	UNP Q38CA1
C	-34	SER	-	expression tag	UNP Q38CA1
C	-33	HIS	-	expression tag	UNP Q38CA1
C	-32	HIS	-	expression tag	UNP Q38CA1
C	-31	HIS	-	expression tag	UNP Q38CA1
C	-30	HIS	-	expression tag	UNP Q38CA1
C	-29	HIS	-	expression tag	UNP Q38CA1
C	-28	HIS	-	expression tag	UNP Q38CA1
C	-27	ASP	-	expression tag	UNP Q38CA1
C	-26	TYR	-	expression tag	UNP Q38CA1
C	-25	ASP	-	expression tag	UNP Q38CA1
C	-24	ILE	-	expression tag	UNP Q38CA1
C	-23	PRO	-	expression tag	UNP Q38CA1
C	-22	THR	-	expression tag	UNP Q38CA1
C	-21	THR	-	expression tag	UNP Q38CA1
C	-20	GLU	-	expression tag	UNP Q38CA1
C	-19	ASN	-	expression tag	UNP Q38CA1
C	-18	LEU	-	expression tag	UNP Q38CA1
C	-17	TYR	-	expression tag	UNP Q38CA1
C	-16	PHE	-	expression tag	UNP Q38CA1
C	-15	GLN	-	expression tag	UNP Q38CA1
C	-14	GLY	-	expression tag	UNP Q38CA1
C	-13	HIS	-	expression tag	UNP Q38CA1
C	-12	MET	-	expression tag	UNP Q38CA1
C	-11	ALA	-	expression tag	UNP Q38CA1

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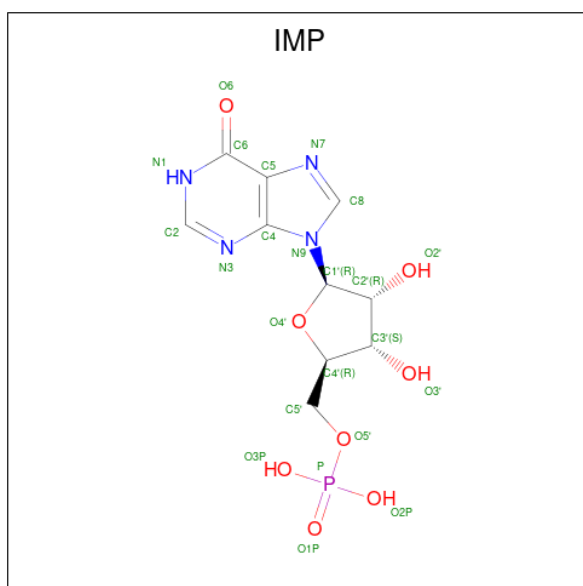
Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	SER	-	expression tag	UNP Q38CA1
C	-9	MET	-	expression tag	UNP Q38CA1
C	-8	THR	-	expression tag	UNP Q38CA1
C	-7	GLY	-	expression tag	UNP Q38CA1
C	-6	GLY	-	expression tag	UNP Q38CA1
C	-5	GLN	-	expression tag	UNP Q38CA1
C	-4	GLN	-	expression tag	UNP Q38CA1
C	-3	MET	-	expression tag	UNP Q38CA1
C	-2	GLY	-	expression tag	UNP Q38CA1
C	-1	ARG	-	expression tag	UNP Q38CA1
C	0	GLY	-	expression tag	UNP Q38CA1
C	1	SER	-	expression tag	UNP Q38CA1
D	-37	MET	-	expression tag	UNP Q38CA1
D	-36	GLY	-	expression tag	UNP Q38CA1
D	-35	SER	-	expression tag	UNP Q38CA1
D	-34	SER	-	expression tag	UNP Q38CA1
D	-33	HIS	-	expression tag	UNP Q38CA1
D	-32	HIS	-	expression tag	UNP Q38CA1
D	-31	HIS	-	expression tag	UNP Q38CA1
D	-30	HIS	-	expression tag	UNP Q38CA1
D	-29	HIS	-	expression tag	UNP Q38CA1
D	-28	HIS	-	expression tag	UNP Q38CA1
D	-27	ASP	-	expression tag	UNP Q38CA1
D	-26	TYR	-	expression tag	UNP Q38CA1
D	-25	ASP	-	expression tag	UNP Q38CA1
D	-24	ILE	-	expression tag	UNP Q38CA1
D	-23	PRO	-	expression tag	UNP Q38CA1
D	-22	THR	-	expression tag	UNP Q38CA1
D	-21	THR	-	expression tag	UNP Q38CA1
D	-20	GLU	-	expression tag	UNP Q38CA1
D	-19	ASN	-	expression tag	UNP Q38CA1
D	-18	LEU	-	expression tag	UNP Q38CA1
D	-17	TYR	-	expression tag	UNP Q38CA1
D	-16	PHE	-	expression tag	UNP Q38CA1
D	-15	GLN	-	expression tag	UNP Q38CA1
D	-14	GLY	-	expression tag	UNP Q38CA1
D	-13	HIS	-	expression tag	UNP Q38CA1
D	-12	MET	-	expression tag	UNP Q38CA1
D	-11	ALA	-	expression tag	UNP Q38CA1
D	-10	SER	-	expression tag	UNP Q38CA1
D	-9	MET	-	expression tag	UNP Q38CA1
D	-8	THR	-	expression tag	UNP Q38CA1

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-7	GLY	-	expression tag	UNP Q38CA1
D	-6	GLY	-	expression tag	UNP Q38CA1
D	-5	GLN	-	expression tag	UNP Q38CA1
D	-4	GLN	-	expression tag	UNP Q38CA1
D	-3	MET	-	expression tag	UNP Q38CA1
D	-2	GLY	-	expression tag	UNP Q38CA1
D	-1	ARG	-	expression tag	UNP Q38CA1
D	0	GLY	-	expression tag	UNP Q38CA1
D	1	SER	-	expression tag	UNP Q38CA1

- Molecule 2 is INOSINIC ACID (CCD ID: IMP) (formula: $C_{10}H_{13}N_4O_8P$).



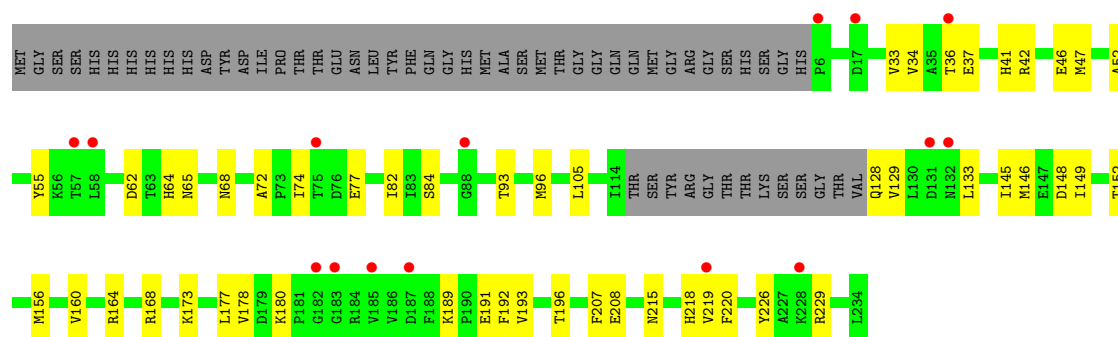
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	C	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	D	1	Total	C	N	O	P	0	0
			23	10	4	8	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

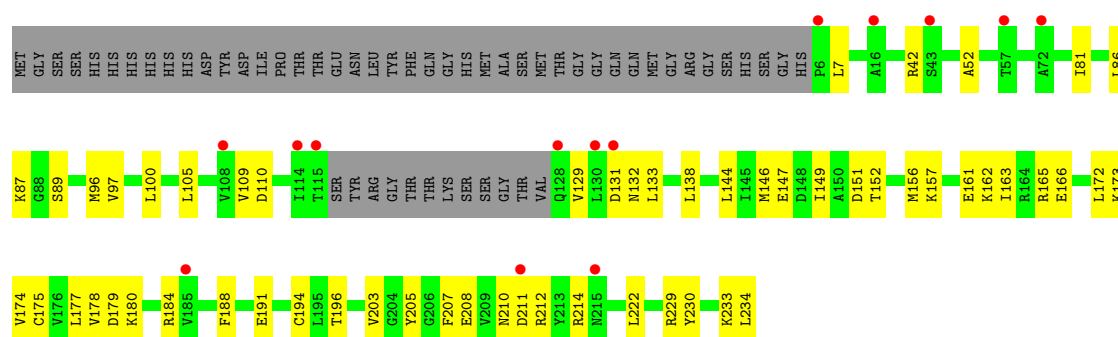
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0
3	C	1	Total Mg 1 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	6	Total O 6 6	0	0
4	B	9	Total O 9 9	0	0
4	C	4	Total O 4 4	0	0
4	D	3	Total O 3 3	0	0



- Molecule 1: Hypoxanthine-guanine phosphoribosyltransferase, putative



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	59.28Å 64.98Å 79.08Å 85.74° 74.58° 62.87°	Depositor
Resolution (Å)	47.27 – 2.96 47.27 – 2.96	Depositor EDS
% Data completeness (in resolution range)	98.5 (47.27-2.96) 98.5 (47.27-2.96)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.96Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.252 , 0.302 0.249 , 0.295	Depositor DCC
R_{free} test set	2000 reflections (9.53%)	wwPDB-VP
Wilson B-factor (Å ²)	25.2	Xtriage
Anisotropy	0.434	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 62.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	6987	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.23 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.0863e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.13	0/1755	0.35	0/2374
1	B	0.15	0/1746	0.39	1/2362 (0.0%)
1	C	0.13	0/1748	0.36	0/2364
1	D	0.15	0/1753	0.33	0/2371
All	All	0.14	0/7002	0.36	1/9471 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	185	VAL	N-CA-C	-5.21	107.33	111.81

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	208	GLU	Peptide

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	1722	0	1756	44	1
1	B	1713	0	1748	52	1
1	C	1715	0	1749	32	0
1	D	1720	0	1751	44	0
2	A	23	0	11	1	0
2	B	23	0	11	3	0
2	C	23	0	11	2	0
2	D	23	0	11	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	6	0	0	0	0
4	B	9	0	0	1	0
4	C	4	0	0	0	0
4	D	3	0	0	1	0
All	All	6987	0	7048	164	1

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 12.

All (164) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:208:GLU:OE1	1:D:211:ASP:N	1.92	1.01
1:B:30:ALA:O	1:B:223:LYS:HE2	1.78	0.84
1:B:37:GLU:H	1:B:218:HIS:HD2	1.26	0.83
1:A:179:ASP:OD1	1:A:184:ARG:NH2	2.16	0.76
1:B:28:PRO:O	1:B:223:LYS:HE3	1.83	0.76
1:C:149:ILE:HD11	1:C:180:LYS:HE2	1.69	0.74
1:A:230:TYR:HE1	1:B:77:GLU:HG2	1.52	0.73
1:A:98:ARG:HB3	1:B:217:ARG:HH21	1.55	0.72
2:D:301:IMP:O2P	4:D:401:HOH:O	2.09	0.71
1:A:217:ARG:HH12	1:B:102:ASP:CG	1.99	0.70
1:D:7:LEU:HD13	1:D:42:ARG:HD2	1.75	0.66
1:C:180:LYS:NZ	2:C:301:IMP:O6	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:151:ASP:O	1:D:184:ARG:HG2	1.95	0.66
1:D:165:ARG:HG3	1:D:166:GLU:HG3	1.78	0.65
1:C:64:HIS:CE1	1:C:77:GLU:HG2	2.32	0.65
1:B:31:GLU:OE1	1:B:200:ARG:NH2	2.31	0.64
1:C:52:ALA:HA	1:C:105:LEU:HD13	1.81	0.62
1:A:36:THR:HA	1:A:218:HIS:HB3	1.80	0.62
1:B:29:MET:HE1	1:B:216:TYR:HE1	1.64	0.62
1:B:37:GLU:H	1:B:218:HIS:CD2	2.14	0.62
1:A:184:ARG:NH1	1:A:188:PHE:O	2.33	0.62
1:C:96:MET:HB3	1:C:146:MET:HE1	1.82	0.61
1:C:33:VAL:HG12	1:C:220:PHE:CD2	2.35	0.60
1:A:15:ASP:OD1	1:A:19:ASN:N	2.33	0.60
1:B:212:ARG:N	1:B:212:ARG:HD2	2.17	0.60
1:A:97:VAL:HG11	1:A:109:VAL:HG21	1.84	0.59
1:A:180:LYS:NZ	2:A:301:IMP:O6	2.34	0.59
1:D:129:VAL:HG11	1:D:132:ASN:ND2	2.18	0.59
1:A:87:LYS:HD2	1:B:111:PHE:HD1	1.66	0.58
1:B:7:LEU:HD11	1:B:38:SER:OG	2.02	0.58
1:A:15:ASP:OD1	1:A:18:GLY:N	2.37	0.58
1:A:87:LYS:HD2	1:B:111:PHE:CD1	2.39	0.58
1:B:203:VAL:HA	1:B:207:PHE:HB2	1.86	0.57
1:A:7:LEU:HD13	1:A:42:ARG:HD2	1.86	0.57
1:A:10:ASN:ND2	1:A:38:SER:H	2.03	0.57
1:D:162:LYS:HA	1:D:165:ARG:HG2	1.86	0.57
1:A:33:VAL:HG22	1:A:220:PHE:HD1	1.70	0.56
1:A:209:VAL:HG21	1:A:222:LEU:HD22	1.87	0.56
1:C:47:MET:HB3	1:C:96:MET:HE1	1.87	0.56
1:D:174:VAL:HG21	1:D:188:PHE:HZ	1.69	0.56
1:C:34:VAL:HG12	1:C:219:VAL:HG23	1.88	0.56
1:D:152:THR:HA	1:D:184:ARG:HA	1.87	0.56
1:B:97:VAL:HB	1:B:107:ASN:HD21	1.70	0.56
1:D:149:ILE:HD11	1:D:180:LYS:HD3	1.87	0.56
1:A:147:GLU:HB3	1:A:176:VAL:HG12	1.88	0.55
1:A:201:TYR:OH	1:A:208:GLU:O	2.22	0.55
2:B:3301:IMP:O1P	4:B:3401:HOH:O	2.18	0.55
1:B:81:ILE:HB	1:B:143:VAL:HG12	1.89	0.55
1:D:149:ILE:HB	1:D:207:PHE:CE1	2.42	0.54
1:B:42:ARG:HH12	1:B:45:LYS:HZ2	1.54	0.54
1:B:147:GLU:HB2	1:B:176:VAL:HG12	1.90	0.54
1:A:37:GLU:CD	1:A:217:ARG:HD3	2.33	0.54
1:C:145:ILE:HG21	1:C:156:MET:HE1	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:LYS:HE3	1:A:231:PRO:HA	1.90	0.53
1:A:64:HIS:HB2	1:A:77:GLU:HB3	1.91	0.53
1:D:100:LEU:HD22	1:D:105:LEU:HD23	1.90	0.53
1:C:226:TYR:CD1	1:C:229:ARG:HD3	2.44	0.52
1:B:44:MET:HE1	1:B:205:TYR:OH	2.10	0.52
1:B:145:ILE:HD11	1:B:172:LEU:HD11	1.91	0.52
1:A:51:LEU:HD22	1:A:144:LEU:HD21	1.92	0.52
1:B:9:PRO:HA	1:B:38:SER:CB	2.40	0.52
1:D:144:LEU:HD21	1:D:175:CYS:HB3	1.92	0.52
1:D:173:LYS:HG2	1:D:191:GLU:HG2	1.92	0.51
1:D:208:GLU:HB3	1:D:214:ARG:HG3	1.91	0.51
1:D:177:LEU:O	1:D:194:CYS:HB2	2.10	0.51
1:B:30:ALA:C	1:B:223:LYS:HE2	2.36	0.51
1:A:80:LEU:HD11	1:A:144:LEU:HB2	1.92	0.50
1:B:15:ASP:HB2	1:B:19:ASN:H	1.77	0.50
1:B:89:SER:O	1:B:93:THR:HG23	2.12	0.50
1:C:33:VAL:HG12	1:C:220:PHE:HD2	1.75	0.50
1:A:37:GLU:OE2	1:A:217:ARG:HD3	2.12	0.50
1:B:149:ILE:HD12	1:B:178:VAL:HB	1.93	0.49
1:C:82:ILE:HG23	1:C:146:MET:HG3	1.94	0.49
1:A:149:ILE:HD11	1:A:180:LYS:HD2	1.95	0.48
1:A:9:PRO:HD2	1:C:196:THR:O	2.13	0.48
1:B:100:LEU:HD13	1:B:105:LEU:HD23	1.94	0.48
1:C:36:THR:HA	1:C:218:HIS:HB3	1.96	0.48
1:D:52:ALA:HA	1:D:105:LEU:HD22	1.96	0.47
1:D:97:VAL:HG11	1:D:109:VAL:HG21	1.95	0.47
1:A:148:ASP:OD1	1:A:149:ILE:N	2.47	0.47
1:A:6:PRO:HG3	1:C:193:VAL:HB	1.96	0.47
1:C:173:LYS:HD3	1:C:191:GLU:HG2	1.96	0.47
1:B:45:LYS:HE2	1:B:99:TYR:CE1	2.50	0.46
1:A:149:ILE:HB	1:A:207:PHE:CE1	2.50	0.46
1:C:47:MET:SD	1:C:192:PHE:HB3	2.55	0.46
1:D:96:MET:HB3	1:D:146:MET:HE1	1.97	0.46
1:C:152:THR:OG1	2:C:301:IMP:O1P	2.33	0.46
1:A:29:MET:HE3	1:A:29:MET:H	1.79	0.46
1:A:156:MET:O	1:A:160:VAL:HG13	2.15	0.45
1:C:55:TYR:O	1:C:74:ILE:HG13	2.16	0.45
1:D:203:VAL:HG21	1:D:222:LEU:HD22	1.97	0.45
1:D:208:GLU:OE1	1:D:210:ASN:N	2.50	0.45
1:B:179:ASP:OD1	1:B:184:ARG:NH2	2.40	0.45
1:D:149:ILE:CG2	2:D:301:IMP:H2'	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:164:ARG:NH2	1:B:169:PRO:O	2.49	0.45
1:B:152:THR:OG1	2:B:3301:IMP:O3P	2.34	0.45
1:B:33:VAL:HG11	1:B:36:THR:HG22	1.97	0.45
1:D:149:ILE:HB	1:D:207:PHE:HE1	1.80	0.45
1:B:60:HIS:HB2	1:B:140:GLY:O	2.17	0.44
1:B:145:ILE:HG21	1:B:156:MET:HE1	2.00	0.44
1:B:7:LEU:HD13	1:B:42:ARG:HG3	1.99	0.44
1:D:110:ASP:HB3	1:D:133:LEU:HD22	1.99	0.44
1:D:233:LYS:NZ	1:D:234:LEU:H	2.15	0.44
1:A:29:MET:O	1:A:222:LEU:HD12	2.17	0.44
1:A:132:ASN:OD1	1:A:133:LEU:N	2.51	0.44
1:D:149:ILE:HD13	1:D:178:VAL:HG13	1.99	0.44
1:B:211:ASP:OD1	1:B:214:ARG:NH1	2.51	0.44
1:B:29:MET:CE	1:B:216:TYR:HE1	2.30	0.44
1:D:86:LEU:HA	1:D:87:LYS:HA	1.76	0.44
1:A:59:LYS:HZ2	1:A:73:PRO:HB3	1.82	0.44
1:C:215:ASN:OD1	1:C:215:ASN:N	2.50	0.44
1:A:145:ILE:HG13	1:A:172:LEU:HD11	2.00	0.43
1:B:92:PHE:CD2	1:B:177:LEU:HB2	2.53	0.43
1:A:225:GLU:H	1:A:225:GLU:HG3	1.57	0.43
1:A:9:PRO:HA	1:A:38:SER:CB	2.48	0.43
1:A:234:LEU:C	1:B:136:THR:H	2.27	0.43
1:C:177:LEU:HG	1:C:178:VAL:HG23	2.00	0.43
1:C:62:ASP:HB2	1:C:72:ALA:HB2	2.01	0.43
1:C:148:ASP:OD1	1:C:148:ASP:N	2.51	0.43
1:D:149:ILE:HD12	1:D:178:VAL:O	2.18	0.43
1:C:149:ILE:HD12	1:C:178:VAL:O	2.18	0.43
1:B:40:ILE:HG22	1:B:44:MET:HE2	2.00	0.42
1:D:131:ASP:OD1	1:D:131:ASP:N	2.53	0.42
1:A:145:ILE:HD12	1:A:160:VAL:HG12	2.02	0.42
1:B:29:MET:HE2	1:B:29:MET:HB3	1.92	0.42
1:B:86:LEU:N	1:B:112:ILE:O	2.52	0.42
1:C:65:ASN:OD1	1:C:68:ASN:ND2	2.53	0.42
1:C:84:SER:OG	1:C:93:THR:OG1	2.31	0.42
1:D:157:LYS:O	1:D:161:GLU:HG2	2.19	0.42
1:C:160:VAL:O	1:C:164:ARG:HB2	2.18	0.42
1:A:44:MET:HE1	1:A:205:TYR:CE2	2.53	0.42
1:B:212:ARG:NH2	1:B:234:LEU:HD11	2.35	0.42
1:D:156:MET:HB3	1:D:188:PHE:CE1	2.54	0.42
1:B:34:VAL:HB	1:B:219:VAL:HG12	2.00	0.42
1:B:154:ARG:HE	1:B:154:ARG:HB2	1.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:229:ARG:HG2	1:D:230:TYR:CE1	2.55	0.42
1:A:60:HIS:HB3	1:A:142:HIS:CD2	2.54	0.42
1:A:229:ARG:HD3	1:A:230:TYR:CE2	2.54	0.42
1:B:233:LYS:HE3	1:B:233:LYS:HB2	1.84	0.42
1:B:9:PRO:HD2	1:D:196:THR:O	2.20	0.42
1:B:180:LYS:NZ	2:B:3301:IMP:O6	2.40	0.41
1:B:210:ASN:O	1:B:211:ASP:HB2	2.20	0.41
1:D:81:ILE:HG21	1:D:138:LEU:HD11	2.02	0.41
1:A:10:ASN:HD21	1:A:38:SER:H	1.68	0.41
1:D:212:ARG:NH2	1:D:230:TYR:HB3	2.35	0.41
1:D:212:ARG:HH21	1:D:230:TYR:HB3	1.84	0.41
1:B:29:MET:HE1	1:B:216:TYR:CE1	2.51	0.41
1:C:208:GLU:N	1:C:208:GLU:OE1	2.53	0.41
1:B:9:PRO:HA	1:B:38:SER:HB2	2.03	0.41
1:D:96:MET:HG2	1:D:146:MET:HE1	2.01	0.41
1:D:177:LEU:HD21	1:D:205:TYR:HD2	1.86	0.41
1:D:179:ASP:OD1	1:D:184:ARG:NH2	2.48	0.41
1:B:208:GLU:HB3	1:B:214:ARG:HG3	2.03	0.41
1:C:149:ILE:HB	1:C:207:PHE:CE2	2.56	0.41
1:D:147:GLU:HG2	1:D:156:MET:HE3	2.02	0.41
1:D:152:THR:OG1	2:D:301:IMP:O3P	2.37	0.41
1:C:42:ARG:O	1:C:46:GLU:HG3	2.20	0.41
1:D:163:ILE:HB	1:D:172:LEU:HD21	2.03	0.41
1:A:130:LEU:HD23	1:A:130:LEU:HA	1.91	0.40
1:B:29:MET:SD	1:B:216:TYR:HE1	2.44	0.40
1:D:208:GLU:HB3	1:D:214:ARG:CG	2.51	0.40
1:C:37:GLU:HG3	1:C:41:HIS:ND1	2.36	0.40
1:C:189:LYS:NZ	1:D:184:ARG:O	2.44	0.40
1:D:129:VAL:HG11	1:D:132:ASN:HD22	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:GLU:OE1	1:B:14:ARG:NH2[1_565]	2.07	0.13

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	213/272 (78%)	203 (95%)	10 (5%)	0	100	100
1	B	212/272 (78%)	199 (94%)	13 (6%)	0	100	100
1	C	212/272 (78%)	202 (95%)	8 (4%)	2 (1%)	14	34
1	D	213/272 (78%)	206 (97%)	7 (3%)	0	100	100
All	All	850/1088 (78%)	810 (95%)	38 (4%)	2 (0%)	43	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	168	ARG
1	C	129	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	190/236 (80%)	186 (98%)	4 (2%)	47	69
1	B	189/236 (80%)	189 (100%)	0	100	100
1	C	189/236 (80%)	187 (99%)	2 (1%)	65	80
1	D	189/236 (80%)	188 (100%)	1 (0%)	81	90
All	All	757/944 (80%)	750 (99%)	7 (1%)	70	83

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

Mol	Chain	Res	Type
1	A	80	LEU
1	A	89	SER
1	A	146	MET
1	A	180	LYS
1	C	128	GLN
1	C	133	LEU
1	D	89	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	B	19	ASN
1	B	78	ASN
1	B	218	HIS
1	C	64	HIS
1	C	65	ASN
1	C	68	ASN
1	D	107	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	IMP	B	3301	-	25,25,25	1.13	1 (4%)	37,38,38	2.05	9 (24%)
2	IMP	A	301	-	25,25,25	1.12	1 (4%)	37,38,38	2.06	9 (24%)
2	IMP	C	301	-	25,25,25	1.12	1 (4%)	37,38,38	2.07	10 (27%)
2	IMP	D	301	-	25,25,25	1.16	1 (4%)	37,38,38	2.13	10 (27%)

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	IMP	B	3301	-	-	3/10/26/26	0/3/3/3
2	IMP	A	301	-	-	2/10/26/26	0/3/3/3
2	IMP	C	301	-	-	5/10/26/26	0/3/3/3
2	IMP	D	301	-	-	1/10/26/26	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	IMP	C2-N3	4.69	1.38	1.30
2	B	3301	IMP	C2-N3	4.55	1.38	1.30
2	C	301	IMP	C2-N3	4.48	1.38	1.30
2	A	301	IMP	C2-N3	4.48	1.38	1.30

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	3301	IMP	C5-C6-N1	6.08	119.53	110.78
2	D	301	IMP	C5-C6-N1	6.07	119.51	110.78
2	A	301	IMP	C5-C6-N1	5.98	119.37	110.78
2	C	301	IMP	C5-C6-N1	5.74	119.03	110.78
2	D	301	IMP	C5-C4-N3	-4.97	119.96	127.27
2	D	301	IMP	C2-N3-C4	4.96	120.26	111.53
2	B	3301	IMP	C2-N3-C4	4.95	120.24	111.53
2	A	301	IMP	C2-N3-C4	4.92	120.19	111.53
2	C	301	IMP	N1-C2-N3	-4.70	120.08	125.50
2	C	301	IMP	C2-N3-C4	4.67	119.75	111.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	IMP	N1-C2-N3	-4.64	120.16	125.50
2	B	3301	IMP	N1-C2-N3	-4.63	120.17	125.50
2	B	3301	IMP	C5-C4-N3	-4.60	120.49	127.27
2	A	301	IMP	C5-C4-N3	-4.56	120.56	127.27
2	D	301	IMP	N1-C2-N3	-4.39	120.45	125.50
2	C	301	IMP	C5-C4-N3	-4.32	120.91	127.27
2	D	301	IMP	C2-N1-C6	-3.94	119.68	125.09
2	B	3301	IMP	C2-N1-C6	-3.74	119.96	125.09
2	A	301	IMP	C2-N1-C6	-3.69	120.04	125.09
2	C	301	IMP	C2-N1-C6	-3.51	120.28	125.09
2	C	301	IMP	O6-C6-C5	-3.21	118.05	126.53
2	C	301	IMP	N9-C8-N7	-2.99	107.86	113.40
2	C	301	IMP	O3P-P-O2P	2.86	118.54	107.80
2	A	301	IMP	N9-C8-N7	-2.69	108.41	113.40
2	D	301	IMP	O3P-P-O2P	2.68	117.86	107.80
2	D	301	IMP	N9-C4-N3	2.66	132.96	126.90
2	C	301	IMP	C8-N7-C5	2.58	108.85	104.26
2	A	301	IMP	O6-C6-C5	-2.57	119.75	126.53
2	A	301	IMP	C8-N7-C5	2.55	108.81	104.26
2	B	3301	IMP	N9-C8-N7	-2.55	108.68	113.40
2	D	301	IMP	O6-C6-C5	-2.48	120.00	126.53
2	B	3301	IMP	O6-C6-C5	-2.47	120.02	126.53
2	B	3301	IMP	C8-N7-C5	2.46	108.65	104.26
2	D	301	IMP	N9-C8-N7	-2.40	108.94	113.40
2	D	301	IMP	C8-N7-C5	2.39	108.52	104.26
2	C	301	IMP	N9-C4-N3	2.32	132.18	126.90
2	B	3301	IMP	N9-C4-N3	2.27	132.08	126.90
2	A	301	IMP	N9-C4-N3	2.18	131.88	126.90

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	301	IMP	C5'-O5'-P-O2P
2	C	301	IMP	C5'-O5'-P-O3P
2	D	301	IMP	C5'-O5'-P-O2P
2	A	301	IMP	O4'-C4'-C5'-O5'
2	A	301	IMP	C3'-C4'-C5'-O5'
2	B	3301	IMP	O4'-C4'-C5'-O5'
2	C	301	IMP	O4'-C4'-C5'-O5'
2	C	301	IMP	C5'-O5'-P-O1P
2	B	3301	IMP	C3'-C4'-C5'-O5'

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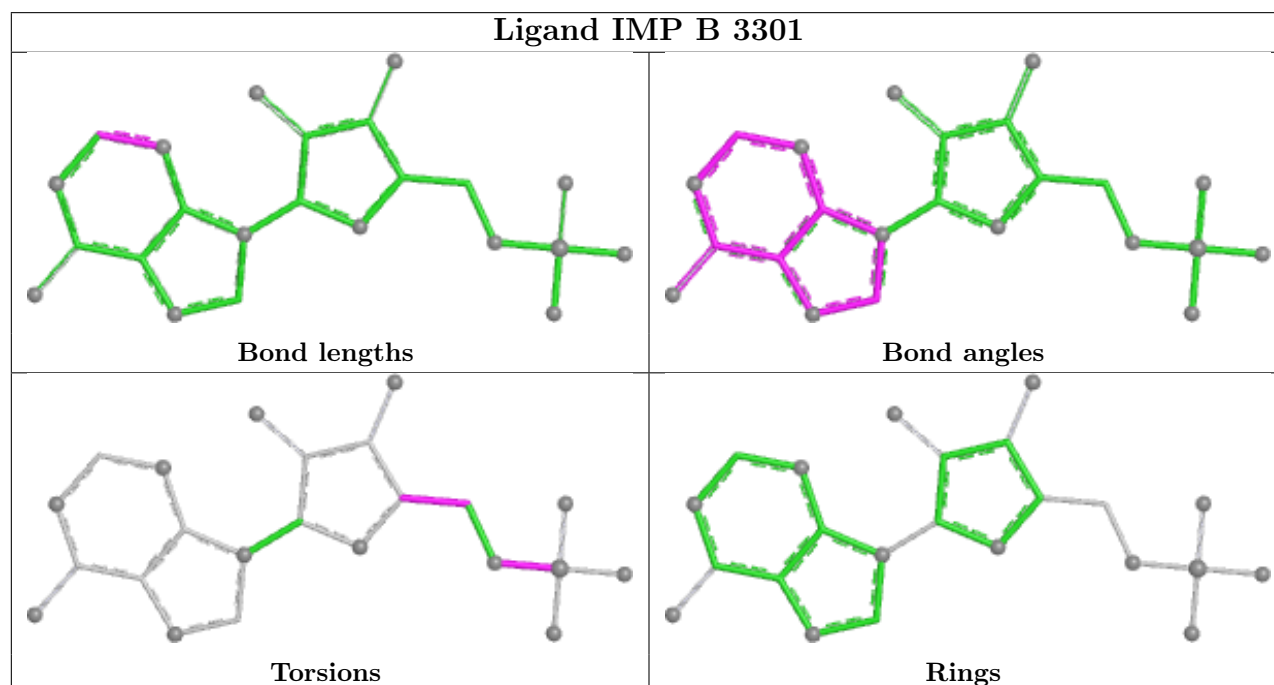
Mol	Chain	Res	Type	Atoms
2	C	301	IMP	C3'-C4'-C5'-O5'
2	B	3301	IMP	C5'-O5'-P-O2P

There are no ring outliers.

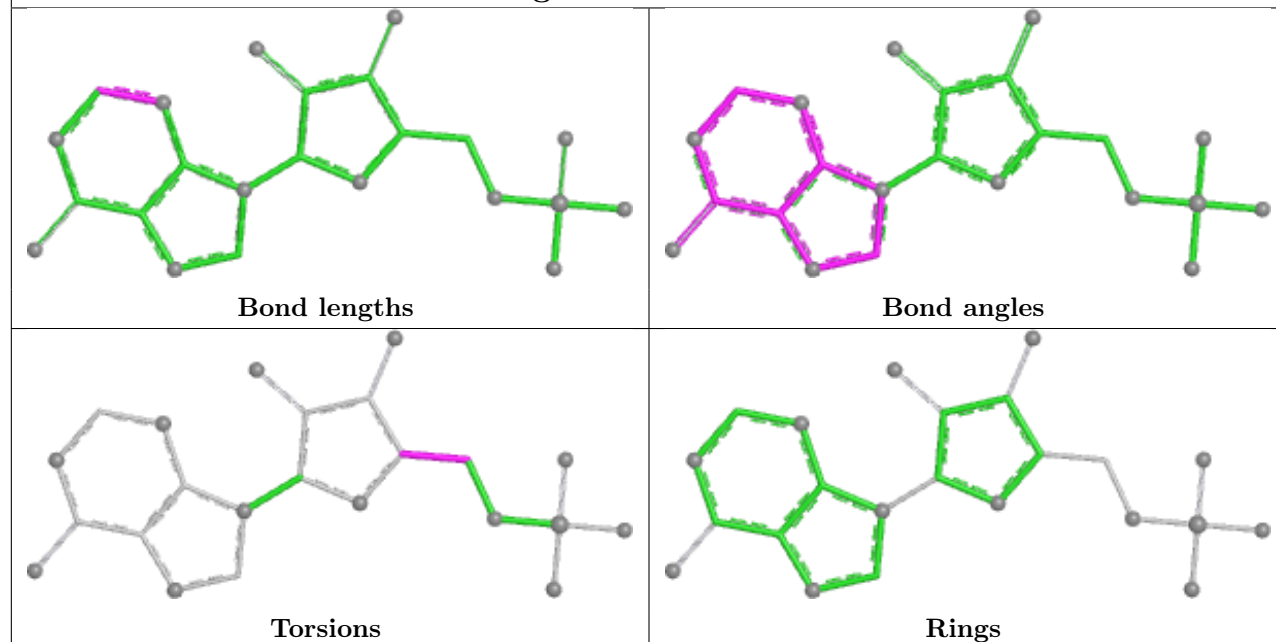
4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3301	IMP	3	0
2	A	301	IMP	1	0
2	C	301	IMP	2	0
2	D	301	IMP	3	0

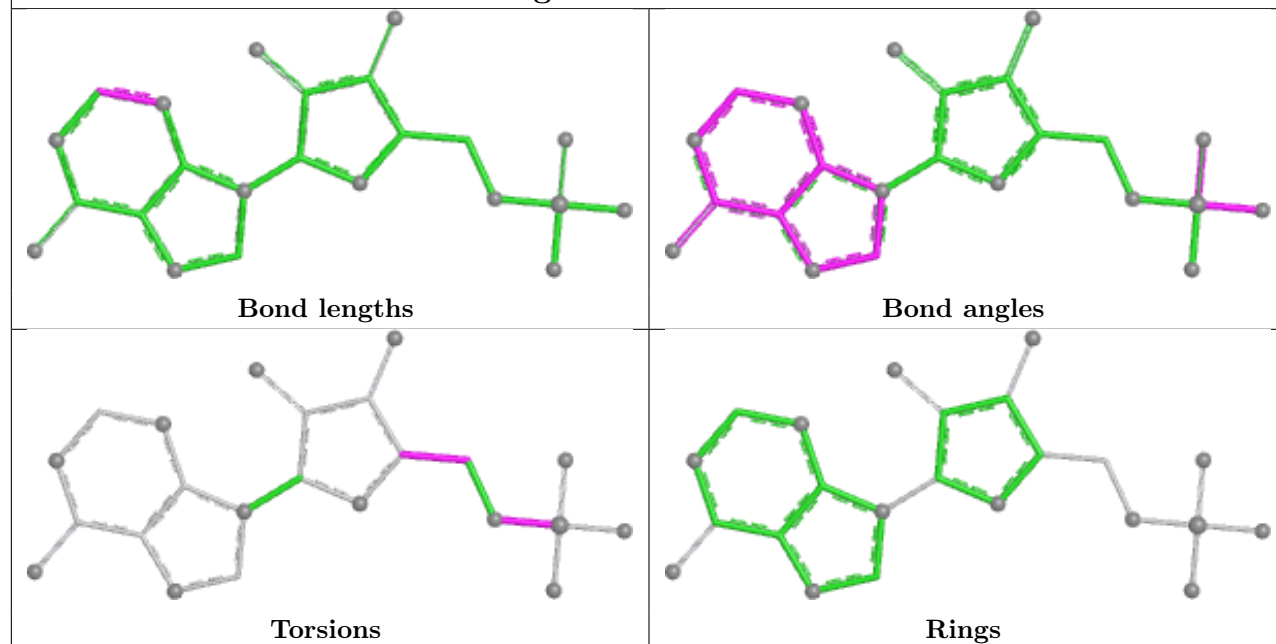
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

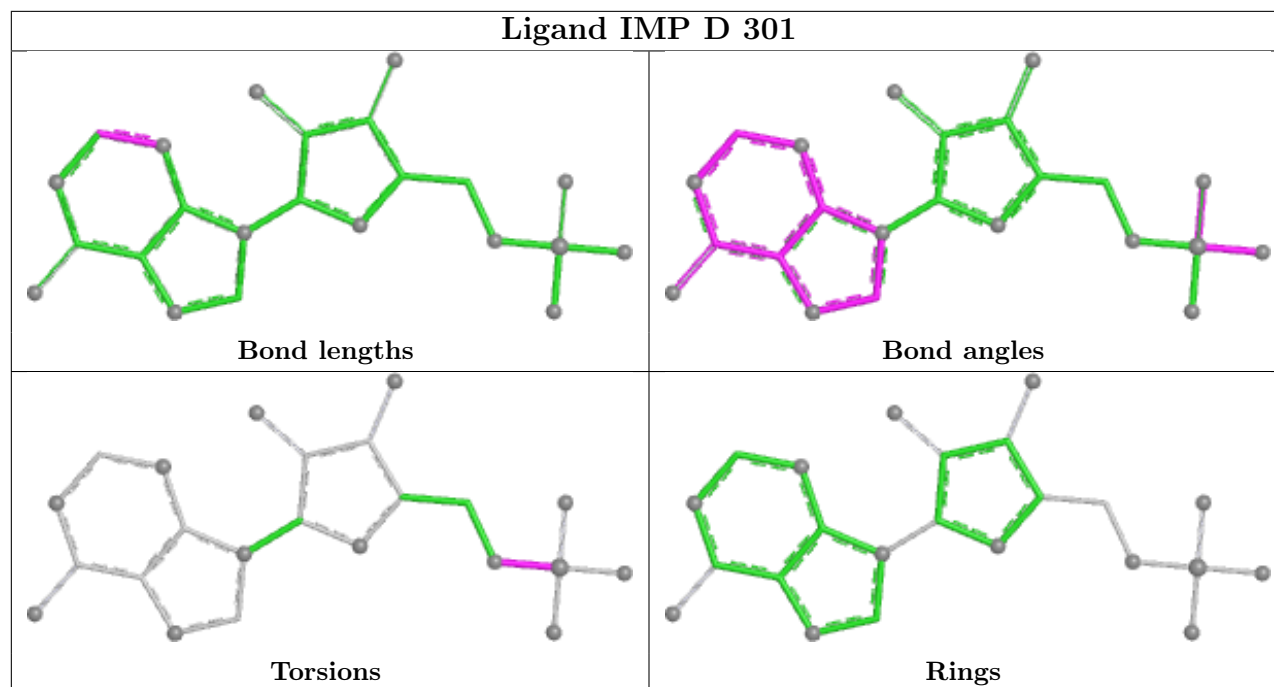


Ligand IMP A 301



Ligand IMP C 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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	217/272 (79%)	0.89	20 (9%) 14 13	23, 40, 62, 80	0
1	B	216/272 (79%)	1.10	26 (12%) 9 8	24, 43, 64, 98	0
1	C	216/272 (79%)	0.91	15 (6%) 23 19	26, 41, 62, 80	0
1	D	217/272 (79%)	0.92	14 (6%) 25 21	22, 39, 62, 87	0
All	All	866/1088 (79%)	0.95	75 (8%) 16 14	22, 41, 63, 98	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	182	GLY	5.0
1	B	199	THR	3.7
1	B	130	LEU	3.6
1	B	115	THR	3.5
1	A	43	SER	3.4
1	A	185	VAL	3.2
1	B	129	VAL	3.2
1	D	115	THR	3.2
1	B	148	ASP	3.1
1	A	131	ASP	3.1
1	D	185	VAL	3.1
1	B	6	PRO	3.1
1	D	130	LEU	3.1
1	A	128	GLN	3.1
1	B	45	LYS	3.0
1	B	38	SER	3.0
1	D	43	SER	3.0
1	B	132	ASN	2.9
1	D	72	ALA	2.9
1	D	131	ASP	2.9
1	D	108	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	15	ASP	2.8
1	C	57	THR	2.7
1	C	6	PRO	2.7
1	D	57	THR	2.7
1	B	43	SER	2.7
1	B	218	HIS	2.6
1	A	152	THR	2.6
1	D	114	ILE	2.6
1	C	185	VAL	2.6
1	A	215	ASN	2.6
1	D	215	ASN	2.6
1	B	131	ASP	2.6
1	B	187	ASP	2.6
1	D	128	GLN	2.5
1	C	88	GLY	2.5
1	B	24	GLY	2.4
1	A	115	THR	2.4
1	D	16	ALA	2.4
1	A	23	ASP	2.3
1	A	196	THR	2.3
1	B	114	ILE	2.3
1	A	183	GLY	2.3
1	C	187	ASP	2.2
1	B	39	THR	2.2
1	B	92	PHE	2.2
1	A	132	ASN	2.2
1	A	158	LEU	2.2
1	B	225	GLU	2.2
1	B	107	ASN	2.2
1	A	155	THR	2.2
1	C	219	VAL	2.2
1	C	58	LEU	2.2
1	C	131	ASP	2.2
1	A	67	GLY	2.2
1	C	183	GLY	2.2
1	B	147	GLU	2.2
1	D	6	PRO	2.2
1	A	130	LEU	2.2
1	B	182	GLY	2.1
1	B	94	ALA	2.1
1	C	17	ASP	2.1
1	C	36	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	225	GLU	2.1
1	A	29	MET	2.1
1	C	132	ASN	2.1
1	A	11	PHE	2.1
1	A	65	ASN	2.1
1	B	165	ARG	2.1
1	C	228	LYS	2.0
1	B	185	VAL	2.0
1	B	194	CYS	2.0
1	C	75	THR	2.0
1	D	211	ASP	2.0
1	B	213	TYR	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

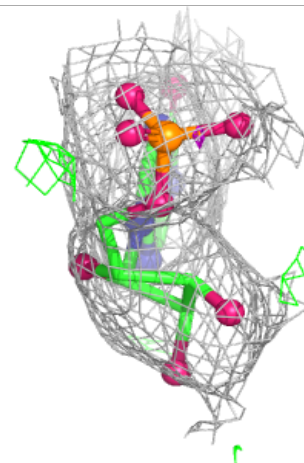
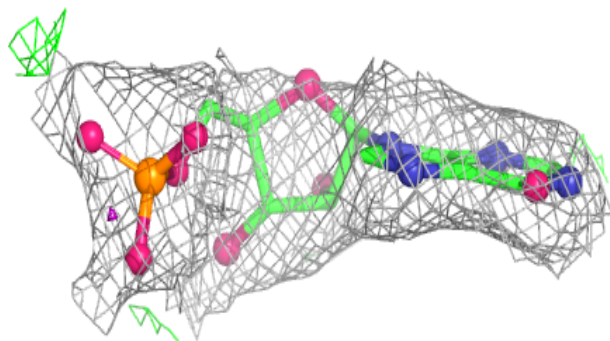
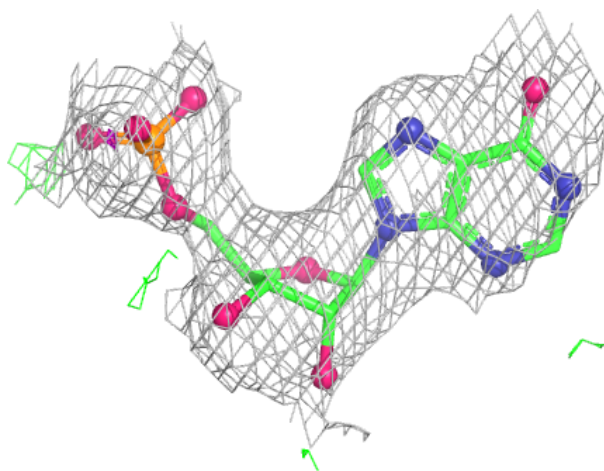
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MG	C	302	1/1	0.80	0.18	48,48,48,48	0
2	IMP	B	3301	23/23	0.83	0.16	48,51,53,54	0
2	IMP	D	301	23/23	0.87	0.14	39,42,46,48	0
3	MG	B	3302	1/1	0.87	0.09	33,33,33,33	0
2	IMP	C	301	23/23	0.87	0.14	45,47,52,54	0
2	IMP	A	301	23/23	0.89	0.13	46,48,52,53	0
3	MG	A	302	1/1	0.90	0.10	26,26,26,26	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.

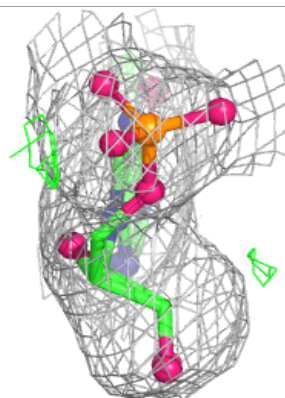
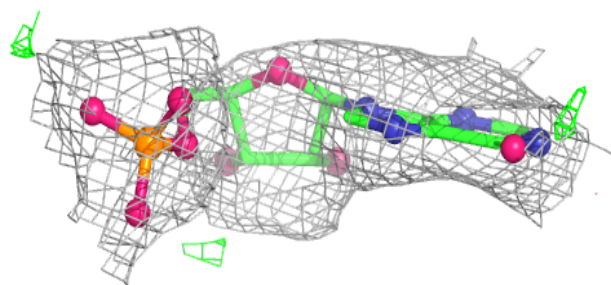
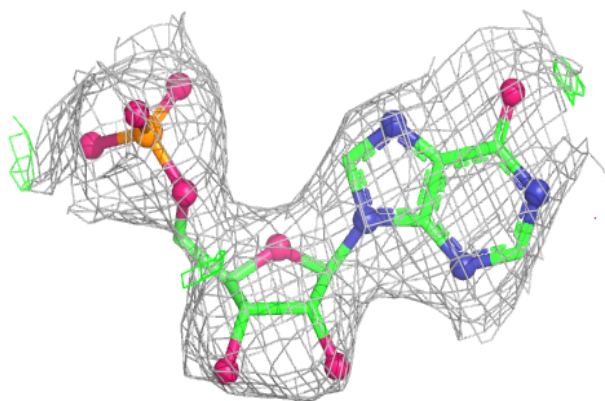
Electron density around IMP B 3301:

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

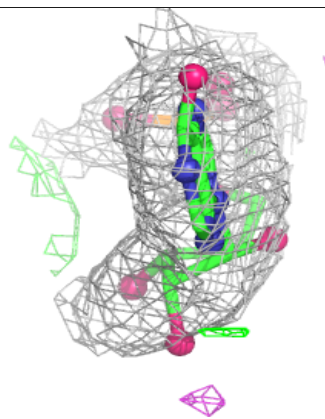
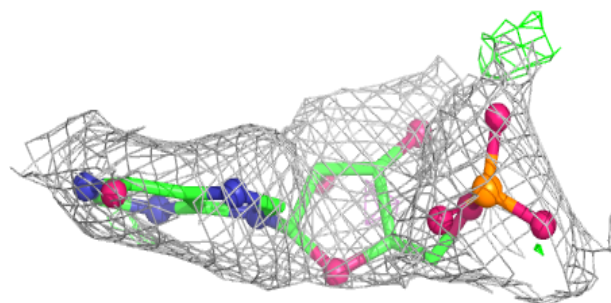
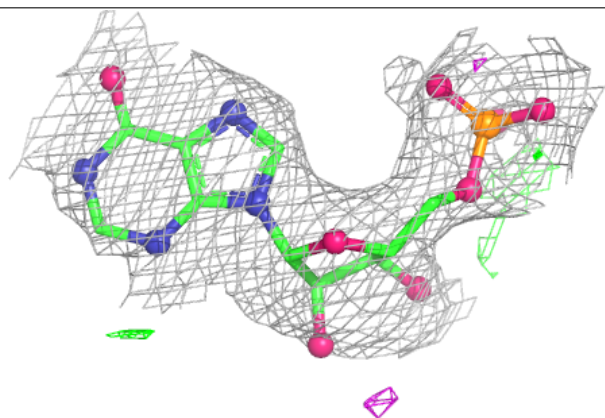


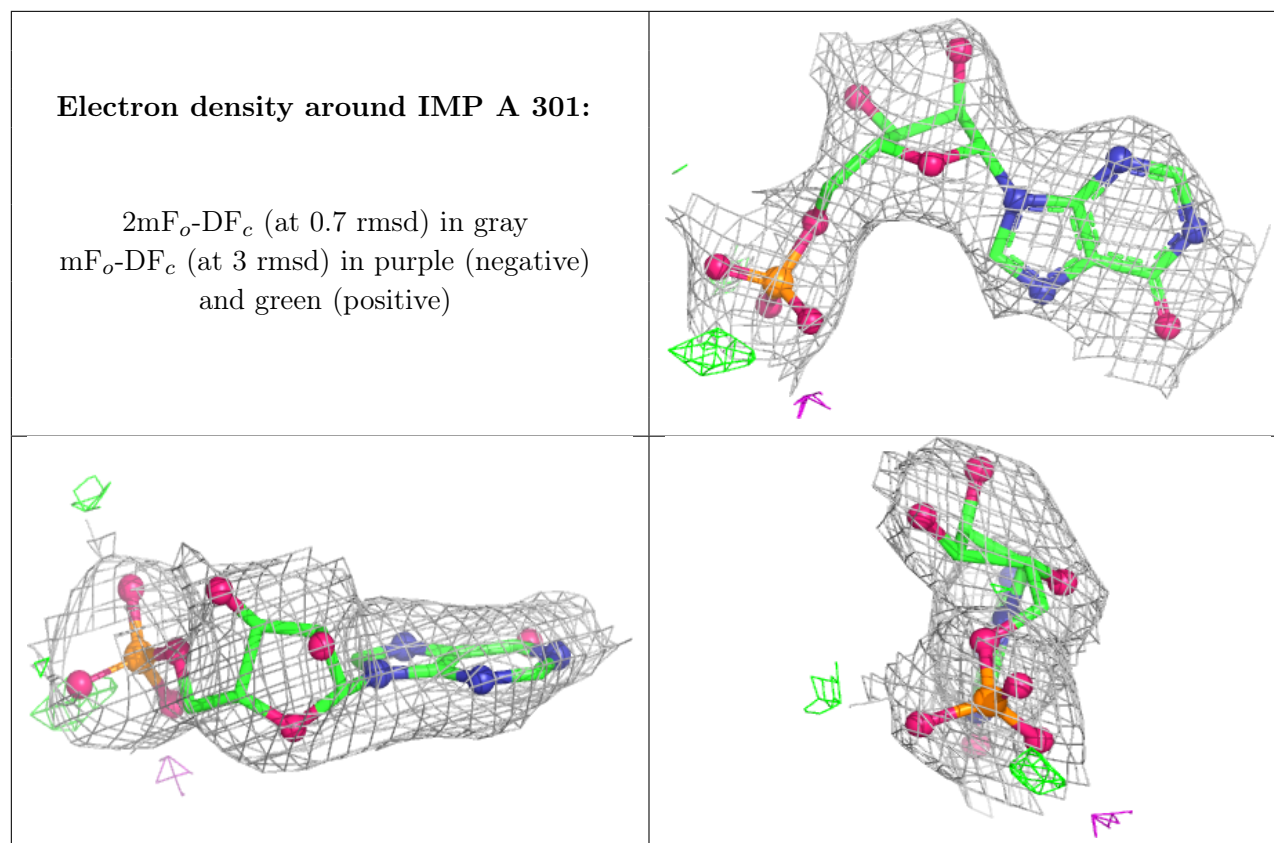
Electron density around IMP D 301:

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

**Electron density around IMP 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.