



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 21, 2026 – 10:47 am BST

PDB ID : 1QK5 / pdb_00001qk5
Title : TOXOPLASMA GONDII HYPOXANTHINE-GUANINE PHOSPHORIBOSYLTRANSFERASE WITH XMP, PYROPHOSPHATE AND TWO MG2+ IONS
Authors : Heroux, A.; White, E.L.; Ross, L.J.; Davis, R.L.; Borhani, D.W.
Deposited on : 1999-07-09
Resolution : 1.60 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

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	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	FAILED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

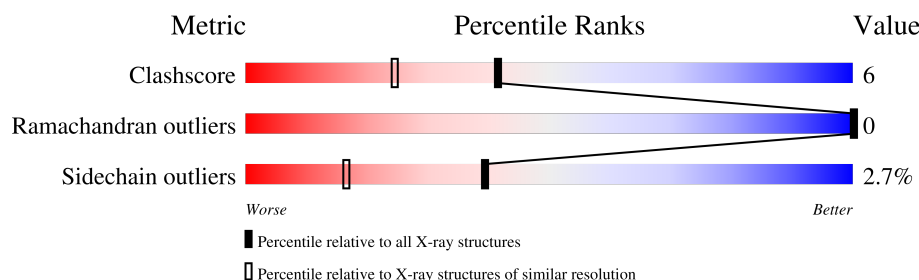
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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))
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)

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 ≥ 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 $\leq 5\%$

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	233	
1	B	233	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3724 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.

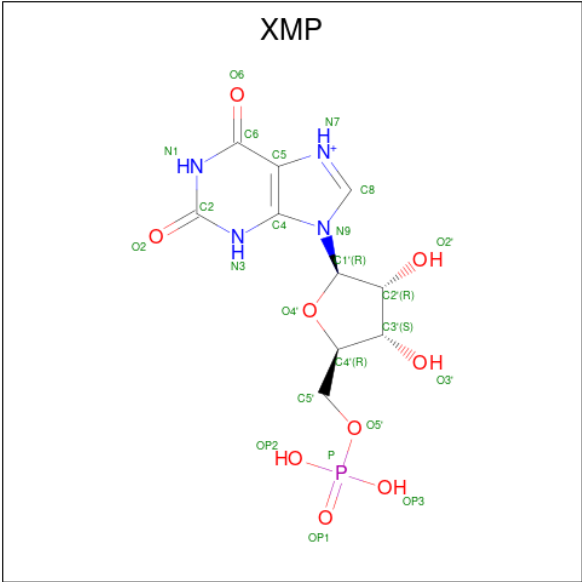
- Molecule 1 is a protein called HYPOXANTHINE-GUANINE PHOSPHORIBOSYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	0	0
			1730	1124	285	313	8			
1	B	217	Total	C	N	O	S	0	0	0
			1742	1133	288	314	7			

There are 8 discrepancies between the modelled and reference sequences:

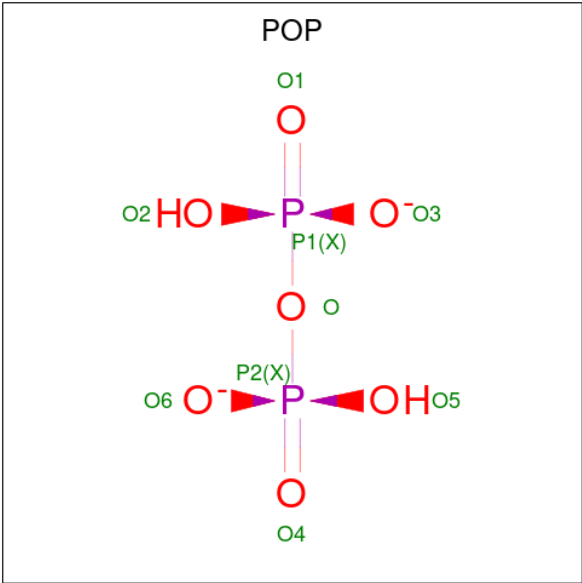
Chain	Residue	Modelled	Actual	Comment	Reference
A	0A	GLY	-	cloning artifact	UNP Q26997
A	0B	SER	-	cloning artifact	UNP Q26997
A	0C	HIS	-	cloning artifact	UNP Q26997
B	0A	GLY	-	cloning artifact	UNP Q26997
B	0B	SER	-	cloning artifact	UNP Q26997
B	0C	HIS	-	cloning artifact	UNP Q26997
A	150	ALA	ASP	engineered mutation	UNP Q26997
B	150	ALA	ASP	engineered mutation	UNP Q26997

- Molecule 2 is XANTHOSINE-5'-MONOPHOSPHATE (CCD ID: XMP) (formula: $C_{10}H_{14}N_4O_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			24	10	4	9	1		
2	B	1	Total	C	N	O	P	0	0
			24	10	4	9	1		

- Molecule 3 is PYROPHOSPHATE 2- (CCD ID: POP) (formula: $\text{H}_2\text{O}_7\text{P}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			9	7	2		
3	B	1	Total	O	P	0	0
			9	7	2		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total 2	Mg 2	0	0
4	B	2	Total 2	Mg 2	0	0

- Molecule 5 is water.

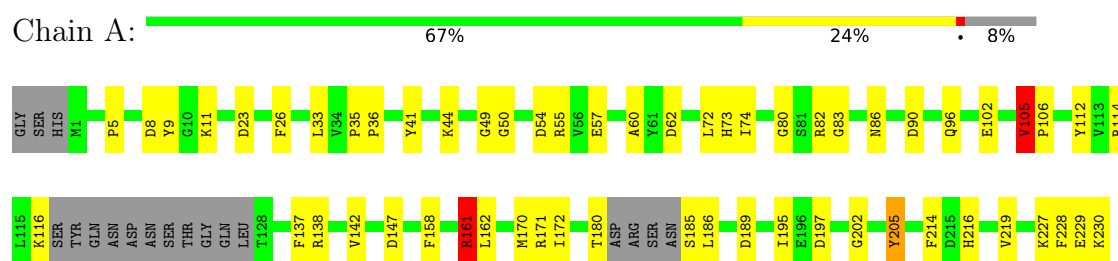
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	102	Total 102	O 102	0	0
5	B	80	Total 80	O 80	0	0

3 Residue-property plots [i](#)

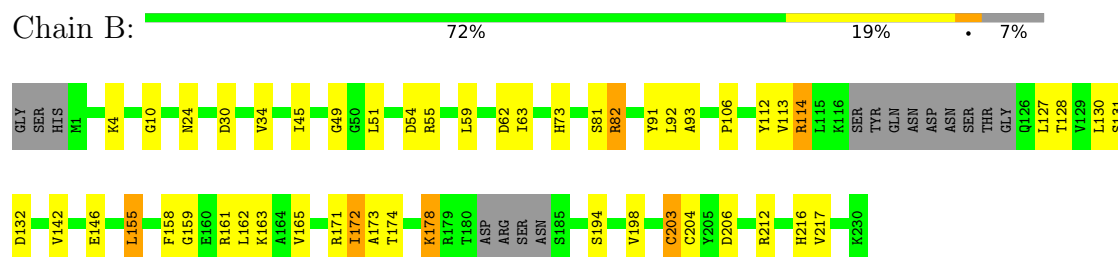
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS failed to run properly.

• Molecule 1: HYPOXANTHINE-GUANINE PHOSPHORIBOSYLTRANSFERASE



• Molecule 1: HYPOXANTHINE-GUANINE PHOSPHORIBOSYLTRANSFERASE



4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	55.21Å 112.25Å 144.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	13.00 – 1.60	Depositor
% Data completeness (in resolution range)	98.0 (13.00-1.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 1.60Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.234 , 0.261	Depositor
Wilson B-factor (Å ²)	21.9	Xtriage
Anisotropy	0.337	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3724	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.20% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: POP, XMP, 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	1.03	1/1771 (0.1%)	1.83	35/2388 (1.5%)
1	B	1.03	1/1783 (0.1%)	1.99	43/2405 (1.8%)
All	All	1.03	2/3554 (0.1%)	1.91	78/4793 (1.6%)

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	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	82	ARG	CD-NE	-5.58	1.38	1.46
1	A	172	ILE	N-CA	5.02	1.51	1.46

All (78) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	82	ARG	CD-NE-CZ	25.64	160.30	124.40
1	B	212	ARG	NE-CZ-NH1	12.21	133.71	121.50
1	B	82	ARG	CG-CD-NE	9.28	132.42	112.00
1	B	82	ARG	NE-CZ-NH2	8.77	127.09	119.20
1	B	55	ARG	CD-NE-CZ	8.64	136.50	124.40
1	B	159	GLY	O-C-N	8.46	130.40	122.19
1	B	54	ASP	CA-CB-CG	8.35	120.95	112.60
1	B	171	ARG	NE-CZ-NH1	8.23	129.73	121.50
1	A	73	HIS	CA-CB-CG	-8.15	105.65	113.80

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	106	PRO	CA-C-N	8.06	128.15	119.28
1	B	106	PRO	C-N-CA	8.06	128.15	119.28
1	B	171	ARG	CD-NE-CZ	7.94	135.51	124.40
1	B	73	HIS	CA-CB-CG	-7.86	105.94	113.80
1	A	62	ASP	CA-CB-CG	7.78	120.38	112.60
1	B	63	ILE	O-C-N	-7.72	114.38	121.87
1	A	55	ARG	CG-CD-NE	-7.58	95.33	112.00
1	B	62	ASP	CA-CB-CG	7.56	120.16	112.60
1	B	91	TYR	CA-C-O	7.51	128.51	120.55
1	A	202	GLY	CA-C-O	-7.45	117.09	122.23
1	B	128	THR	O-C-N	7.34	131.77	123.19
1	B	63	ILE	CA-C-O	7.26	128.50	120.95
1	A	171	ARG	O-C-N	7.25	132.58	123.15
1	B	24	ASN	CA-CB-CG	-7.23	105.37	112.60
1	A	185	SER	CA-C-O	7.20	133.05	120.80
1	B	45	ILE	CA-C-O	7.18	129.07	120.74
1	A	60	ALA	O-C-N	7.06	129.61	122.12
1	B	45	ILE	O-C-N	-6.78	112.79	122.76
1	A	86	ASN	OD1-CG-ND2	-6.70	115.90	122.60
1	B	206	ASP	O-C-N	6.65	132.45	122.94
1	A	205	TYR	CA-C-O	-6.55	113.58	121.05
1	A	23	ASP	CA-CB-CG	6.53	119.13	112.60
1	A	161	ARG	CD-NE-CZ	6.42	133.38	124.40
1	A	50	GLY	CA-C-O	-6.27	114.01	120.66
1	B	163	LYS	CA-C-O	6.14	127.06	120.55
1	B	92	LEU	CA-C-O	6.10	127.02	120.55
1	B	81	SER	CA-CB-OG	-6.09	98.92	111.10
1	A	186	LEU	CA-C-O	-6.07	114.52	121.19
1	B	82	ARG	O-C-N	6.04	128.37	122.09
1	B	198	VAL	CA-C-O	-5.98	115.39	121.67
1	A	54	ASP	CA-CB-CG	5.98	118.58	112.60
1	A	55	ARG	CD-NE-CZ	5.94	132.72	124.40
1	B	173	ALA	O-C-N	-5.93	116.27	123.27
1	A	228	PHE	CA-C-N	5.91	129.23	120.90
1	A	228	PHE	C-N-CA	5.91	129.23	120.90
1	A	185	SER	O-C-N	-5.83	113.67	123.00
1	A	60	ALA	CA-C-O	-5.80	114.40	120.55
1	A	57	GLU	O-C-N	5.79	128.26	122.12
1	B	217	VAL	O-C-N	5.79	129.95	122.77
1	B	34	VAL	CA-C-O	5.72	123.27	119.38
1	B	194	SER	O-C-N	-5.69	116.56	123.27
1	A	161	ARG	O-C-N	-5.67	116.23	122.07

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	189	ASP	CA-CB-CG	-5.67	106.93	112.60
1	A	90	ASP	CA-CB-CG	-5.59	107.01	112.60
1	A	161	ARG	CA-C-O	5.58	126.68	120.82
1	A	105	VAL	O-C-N	5.57	127.45	121.10
1	B	212	ARG	NH1-CZ-NH2	-5.55	112.08	119.30
1	A	83	GLY	O-C-N	5.55	127.52	122.19
1	B	212	ARG	NE-CZ-NH2	-5.54	114.21	119.20
1	B	114	ARG	NH1-CZ-NH2	5.50	126.45	119.30
1	B	34	VAL	O-C-N	-5.42	116.53	121.41
1	B	59	LEU	CA-C-O	5.42	126.51	120.82
1	A	214	PHE	CA-C-O	-5.39	115.26	121.19
1	B	206	ASP	CA-C-O	-5.38	115.33	121.68
1	B	114	ARG	NE-CZ-NH1	-5.30	116.20	121.50
1	B	172	ILE	CB-CA-C	-5.27	102.83	110.63
1	B	10	GLY	O-C-N	5.22	129.06	122.43
1	B	55	ARG	CA-C-O	5.16	125.89	120.42
1	A	80	GLY	N-CA-C	5.15	119.99	113.24
1	B	132	ASP	CA-CB-CG	5.15	117.75	112.60
1	B	178	LYS	N-CA-CB	5.14	118.79	110.41
1	A	138	ARG	CA-C-N	5.12	129.83	122.40
1	A	138	ARG	C-N-CA	5.12	129.83	122.40
1	A	80	GLY	CA-C-N	-5.09	113.80	122.56
1	A	80	GLY	C-N-CA	-5.09	113.80	122.56
1	A	137	PHE	CA-CB-CG	-5.04	108.76	113.80
1	A	158	PHE	CA-C-N	5.03	125.67	119.99
1	A	158	PHE	C-N-CA	5.03	125.67	119.99
1	B	93	ALA	O-C-N	5.00	127.42	122.12

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	195	ILE	Mainchain

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	1730	0	1716	24	0
1	B	1742	0	1734	16	0
2	A	24	0	12	5	0
2	B	24	0	12	3	0
3	A	9	0	0	0	0
3	B	9	0	0	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	102	0	0	4	0
5	B	80	0	0	0	0
All	All	3724	0	3474	45	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 6.

All (45) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:300:XMP:H8	2:A:300:XMP:H5'2	1.35	1.06
2:B:300:XMP:H5'2	2:B:300:XMP:H8	1.28	1.05
1:A:49:GLY:H	1:A:216:HIS:HD2	1.18	0.91
1:B:49:GLY:H	1:B:216:HIS:HD2	1.17	0.90
2:A:300:XMP:H8	2:A:300:XMP:C5'	2.04	0.87
2:A:300:XMP:H5'2	2:A:300:XMP:C8	2.07	0.85
1:B:49:GLY:H	1:B:216:HIS:CD2	2.01	0.78
2:B:300:XMP:H5'2	2:B:300:XMP:C8	2.12	0.76
1:A:49:GLY:H	1:A:216:HIS:CD2	2.03	0.75
1:A:112:TYR:OH	1:B:82:ARG:HD3	1.94	0.66
1:A:161:ARG:HG3	1:A:161:ARG:HH11	1.63	0.64
1:A:96:GLN:HG2	5:A:2048:HOH:O	2.02	0.59
1:B:178:LYS:HE3	2:B:300:XMP:H7	1.70	0.57
1:A:41:TYR:OH	1:A:227:LYS:HD3	2.06	0.56
1:A:116:LYS:C	5:A:2101:HOH:O	2.48	0.56
1:B:127:LEU:O	1:B:161:ARG:NH2	2.39	0.56
1:A:26:PHE:CE2	1:A:219:VAL:HG11	2.42	0.54
2:A:300:XMP:H5'1	5:A:2100:HOH:O	2.06	0.54
1:A:82:ARG:HG3	1:B:112:TYR:OH	2.08	0.53
1:A:229:GLU:O	1:A:230:LYS:C	2.53	0.52
1:B:203:CYS:O	1:B:204:CYS:HB2	2.09	0.51
1:A:5:PRO:HG2	1:A:8:ASP:OD2	2.11	0.50
1:A:8:ASP:HA	1:A:11:LYS:HD2	1.95	0.49
1:A:180:THR:HG22	1:A:197:ASP:CG	2.39	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:LEU:O	1:B:165:VAL:HG22	2.14	0.48
1:B:146:GLU:O	1:B:174:THR:HA	2.15	0.47
1:A:9:TYR:HA	5:A:2070:HOH:O	2.14	0.47
1:A:44:LYS:HB2	1:A:219:VAL:HG22	1.96	0.47
2:A:300:XMP:C5'	2:A:300:XMP:C8	2.81	0.46
1:A:26:PHE:CE2	1:A:219:VAL:CG1	2.98	0.45
1:B:49:GLY:N	1:B:216:HIS:HD2	1.99	0.45
1:B:113:VAL:HG22	1:B:131:SER:HB2	1.98	0.44
1:A:35:PRO:HA	1:A:36:PRO:HD3	1.87	0.44
1:A:9:TYR:CD1	1:A:170:MET:HE2	2.52	0.44
1:A:142:VAL:HG11	1:A:162:LEU:HD21	2.00	0.44
1:B:155:LEU:HD23	1:B:172:ILE:HD13	2.00	0.44
1:B:127:LEU:HD21	1:B:158:PHE:HB2	1.99	0.44
1:A:161:ARG:HH11	1:A:161:ARG:CG	2.29	0.44
1:A:147:ASP:HB2	1:A:205:TYR:CE2	2.54	0.42
1:B:142:VAL:HG11	1:B:162:LEU:HD21	2.01	0.42
1:A:72:LEU:HD21	1:A:74:ILE:HD11	2.02	0.42
1:A:105:VAL:HG23	1:A:106:PRO:HD2	2.03	0.41
1:B:114:ARG:HB3	1:B:130:LEU:HB3	2.03	0.41
1:A:72:LEU:CD2	1:A:74:ILE:HD11	2.51	0.41
1:B:155:LEU:HD12	1:B:155:LEU:HA	1.87	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	209/233 (90%)	206 (99%)	3 (1%)	0	100	100
1	B	211/233 (91%)	205 (97%)	6 (3%)	0	100	100
All	All	420/466 (90%)	411 (98%)	9 (2%)	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	187/206 (91%)	182 (97%)	5 (3%)	39	16
1	B	188/206 (91%)	183 (97%)	5 (3%)	39	16
All	All	375/412 (91%)	365 (97%)	10 (3%)	39	16

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	102	GLU
1	A	105	VAL
1	A	114	ARG
1	A	161	ARG
1	B	4	LYS
1	B	30	ASP
1	B	51	LEU
1	B	155	LEU
1	B	203	CYS

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

Mol	Chain	Res	Type
1	A	86	ASN
1	A	216	HIS
1	B	37	HIS
1	B	86	ASN
1	B	96	GLN
1	B	216	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 4 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	XMP	A	300	-	26,26,26	1.00	1 (3%)	35,40,40	1.73	7 (20%)
3	POP	B	301	4	6,8,8	1.15	0	13,13,13	1.53	3 (23%)
3	POP	A	301	4	6,8,8	1.55	1 (16%)	13,13,13	1.81	3 (23%)
2	XMP	B	300	-	26,26,26	1.01	2 (7%)	35,40,40	1.50	5 (14%)

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	XMP	A	300	-	-	4/10/26/26	0/3/3/3
3	POP	B	301	4	-	0/6/6/6	-
3	POP	A	301	4	-	0/6/6/6	-
2	XMP	B	300	-	-	5/10/26/26	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	301	POP	P2-O6	-2.50	1.45	1.54
2	B	300	XMP	P-OP2	-2.18	1.46	1.54
2	B	300	XMP	C4-N3	-2.04	1.33	1.37
2	A	300	XMP	C4-N3	-2.01	1.33	1.37

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	300	XMP	C5-C4-N9	-5.99	103.43	106.23
3	A	301	POP	O2-P1-O	4.58	119.98	104.64
2	B	300	XMP	OP2-P-OP3	3.56	121.24	107.64
2	B	300	XMP	C2'-C3'-C4'	-3.54	95.77	102.64
3	B	301	POP	O2-P1-O	3.06	114.91	104.64
2	B	300	XMP	C5-C4-N9	-3.00	104.83	106.23
2	B	300	XMP	O4'-C1'-N9	-2.96	101.60	108.36
2	A	300	XMP	C8-N9-C4	2.77	109.02	106.71
2	A	300	XMP	P-O5'-C5'	2.70	125.74	118.30
3	A	301	POP	O3-P1-O	-2.56	96.05	104.64
3	B	301	POP	P2-O-P1	-2.48	124.33	132.83
3	B	301	POP	O3-P1-O	-2.22	97.18	104.64
2	A	300	XMP	N1-C2-N3	2.19	119.33	115.80
2	A	300	XMP	C4-C5-N7	2.14	112.01	109.23
2	A	300	XMP	O4'-C1'-N9	-2.12	103.53	108.36
2	B	300	XMP	C4'-O4'-C1'	-2.09	104.86	109.47
3	A	301	POP	O6-P2-O5	2.03	115.39	107.64
2	A	300	XMP	O2-C2-N1	-2.01	118.03	121.82

There are no chirality outliers.

All (9) torsion outliers are listed below:

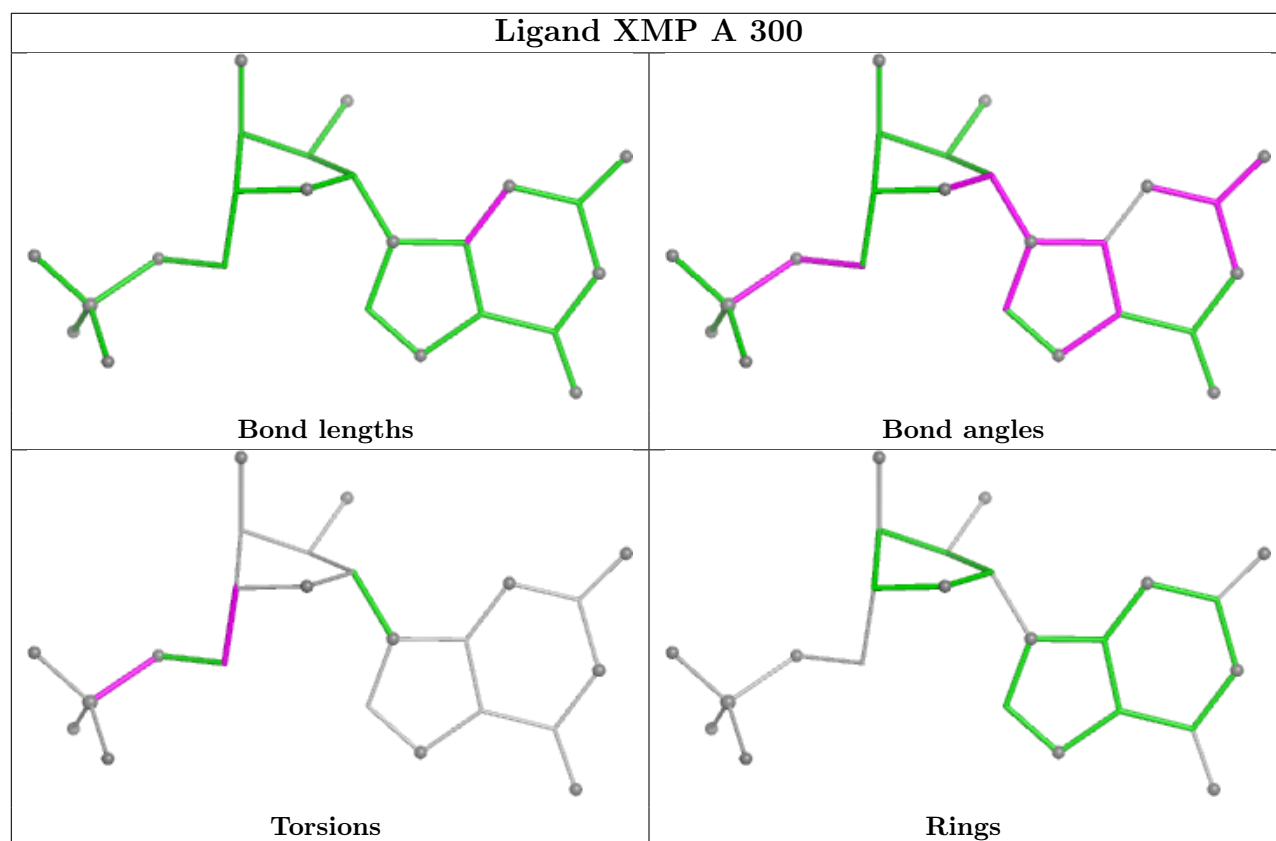
Mol	Chain	Res	Type	Atoms
2	A	300	XMP	C5'-O5'-P-OP3
2	A	300	XMP	C5'-O5'-P-OP2
2	A	300	XMP	C5'-O5'-P-OP1
2	B	300	XMP	C5'-O5'-P-OP3
2	B	300	XMP	C5'-O5'-P-OP2
2	B	300	XMP	O4'-C4'-C5'-O5'
2	B	300	XMP	C5'-O5'-P-OP1
2	B	300	XMP	C3'-C4'-C5'-O5'
2	A	300	XMP	O4'-C4'-C5'-O5'

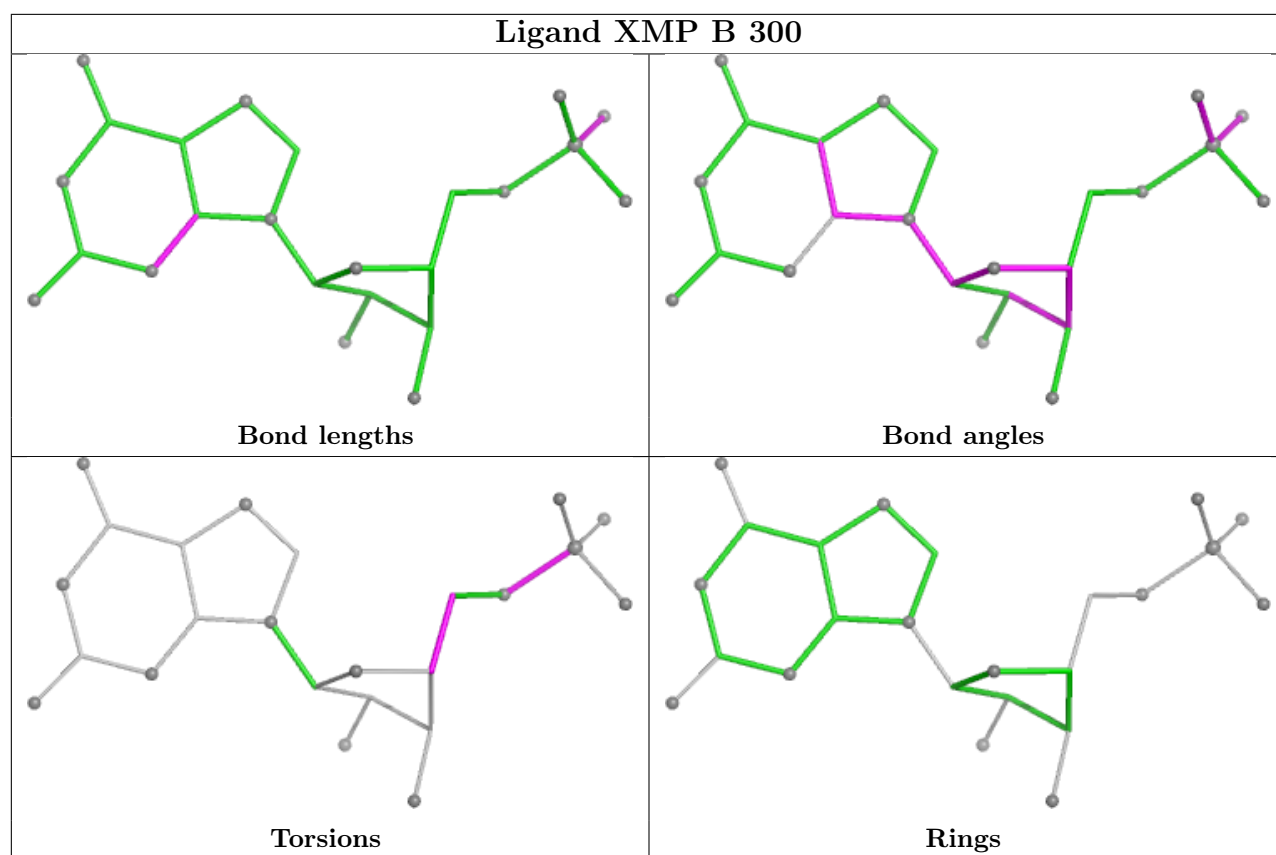
There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	300	XMP	5	0
2	B	300	XMP	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.





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

EDS failed to run properly - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS failed to run properly - this section is therefore empty.

6.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

6.4 Ligands

EDS failed to run properly - this section is therefore empty.

6.5 Other polymers

EDS failed to run properly - this section is therefore empty.