



# Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 03:32 AM GMT

PDB ID : 1QK5  
Title : TOXOPLASMA GONDII HYPOXANTHINE-GUANINE PHOSPHORIBOSYLTRANSFERASE WITH XMP, PYROPHOSPHATE AND TWO MG<sup>2+</sup> IONS  
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Deposited on : 1999-07-09  
Resolution : 1.60 Å (reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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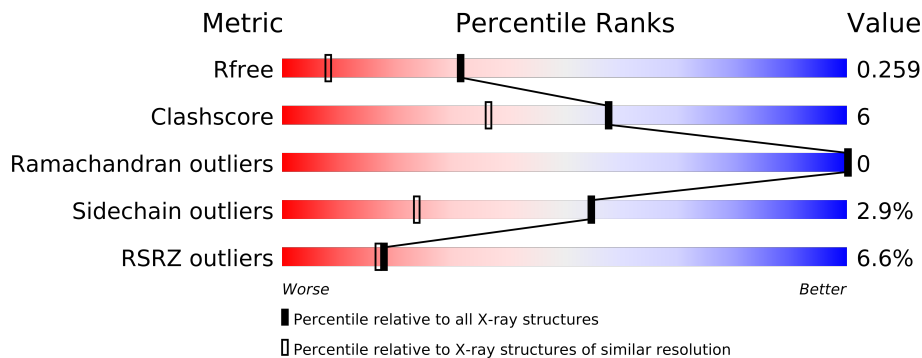
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

# 1 Overall quality at a glance

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))
$R_{free}$	66092	1872 (1.60-1.60)
Clashscore	79885	2199 (1.60-1.60)
Ramachandran outliers	78287	2126 (1.60-1.60)
Sidechain outliers	78261	2125 (1.60-1.60)
RSRZ outliers	66119	1872 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

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

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	MG	A	303	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3724 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

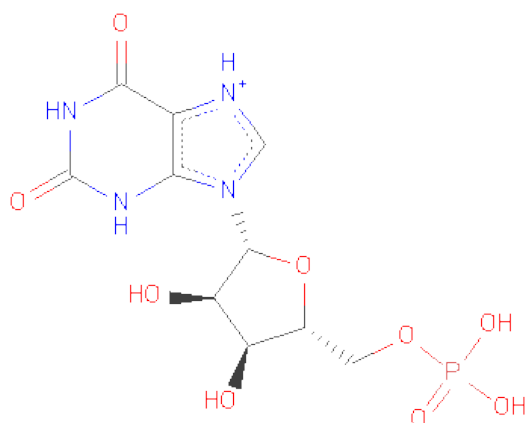
- Molecule 1 is a protein called HYPOXANTHINE-GUANINE PHOSPHORIBOSYLTRANSFERASE.

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:

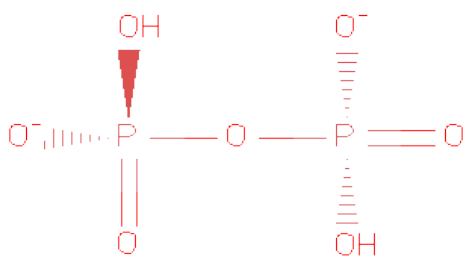
Chain	Residue	Modelled	Actual	Comment	Reference
A	0A	GLY	-	CLONING ARTIFACT	UNP Q26997
A	0B	SER	-	CLONING ARTIFACT	UNP Q26997
A	0C	HIS	-	CLONING ARTIFACT	UNP Q26997
A	150	ALA	ASP	ENGINEERED MUTATION	UNP Q26997
B	0A	GLY	-	CLONING ARTIFACT	UNP Q26997
B	0B	SER	-	CLONING ARTIFACT	UNP Q26997
B	0C	HIS	-	CLONING ARTIFACT	UNP Q26997
B	150	ALA	ASP	ENGINEERED MUTATION	UNP Q26997

- Molecule 2 is XANTHOSINE-5'-MONOPHOSPHATE (three-letter code: XMP) (formula: C<sub>10</sub>H<sub>14</sub>N<sub>4</sub>O<sub>9</sub>P).



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- (three-letter code: POP) (formula:  $\text{H}_2\text{O}_7\text{P}_2$ ).



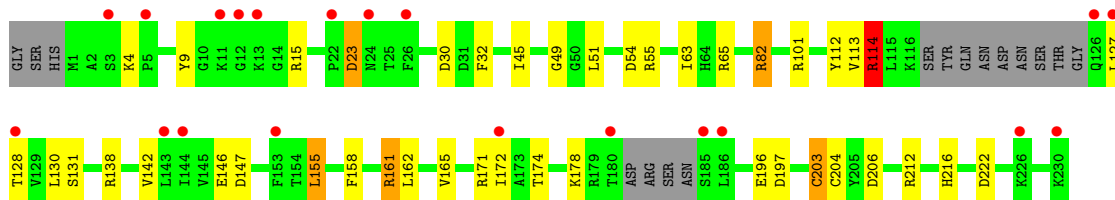
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			9	7	2		
3	B	1	Total	O	P	0	0
			9	7	2		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	55.21Å 112.25Å 144.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	13.00 – 1.60 24.41 – 1.60	Depositor EDS
% Data completeness (in resolution range)	98.0 (13.00-1.60) 97.9 (24.41-1.60)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.14 (at 1.60Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.234 , 0.261 0.231 , 0.259	Depositor DCC
$R_{free}$ test set	2957 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	21.9	Xtriage
Anisotropy	0.337	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 38.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 58200 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3724	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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.*

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

## 5 Model quality

### 5.1 Standard geometry

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

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	0.76	0/1771	1.59	30/2388 (1.3%)
1	B	0.77	0/1783	1.90	38/2405 (1.6%)
All	All	0.76	0/3554	1.76	68/4793 (1.4%)

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

There are no bond length outliers.

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	212	ARG	NE-CZ-NH1	26.82	133.71	120.30
1	B	82	ARG	CD-NE-CZ	26.21	160.30	123.60
1	B	171	ARG	NE-CZ-NH1	18.86	129.73	120.30
1	B	82	ARG	NE-CZ-NH2	13.59	127.09	120.30
1	B	212	ARG	NE-CZ-NH2	-12.18	114.21	120.30
1	B	101	ARG	NE-CZ-NH2	-10.51	115.05	120.30
1	B	23	ASP	CB-CG-OD1	10.02	127.32	118.30
1	B	101	ARG	NE-CZ-NH1	9.86	125.23	120.30
1	B	82	ARG	CG-CD-NE	9.82	132.42	111.80
1	A	133	ASP	CB-CG-OD2	-9.73	109.54	118.30
1	B	55	ARG	CD-NE-CZ	9.21	136.50	123.60
1	A	171	ARG	NE-CZ-NH2	-8.91	115.85	120.30
1	B	32	PHE	CB-CG-CD1	8.54	126.78	120.80
1	B	171	ARG	CD-NE-CZ	8.51	135.51	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	114	ARG	NE-CZ-NH1	-8.19	116.20	120.30
1	B	9	TYR	CB-CG-CD2	8.12	125.88	121.00
1	A	171	ARG	NE-CZ-NH1	8.10	124.35	120.30
1	B	55	ARG	NE-CZ-NH1	8.02	124.31	120.30
1	B	171	ARG	NE-CZ-NH2	-7.96	116.32	120.30
1	A	23	ASP	CB-CG-OD2	7.88	125.39	118.30
1	A	55	ARG	CG-CD-NE	-7.84	95.33	111.80
1	A	55	ARG	NE-CZ-NH2	-7.81	116.40	120.30
1	A	147	ASP	CB-CG-OD1	7.66	125.20	118.30
1	B	82	ARG	NE-CZ-NH1	-7.54	116.53	120.30
1	B	222	ASP	CB-CG-OD1	7.46	125.02	118.30
1	B	65	ARG	NE-CZ-NH1	7.38	123.99	120.30
1	B	9	TYR	CB-CG-CD1	-7.24	116.65	121.00
1	A	101	ARG	NE-CZ-NH1	7.16	123.88	120.30
1	A	43	ASP	CB-CG-OD2	-7.14	111.87	118.30
1	A	161	ARG	CD-NE-CZ	6.99	133.38	123.60
1	A	139	ASP	CB-CG-OD2	-6.73	112.25	118.30
1	A	206	ASP	CB-CG-OD2	6.68	124.32	118.30
1	B	212	ARG	NH1-CZ-NH2	-6.65	112.08	119.40
1	B	147	ASP	CB-CG-OD1	6.64	124.28	118.30
1	A	161	ARG	NE-CZ-NH1	6.61	123.60	120.30
1	A	55	ARG	CD-NE-CZ	6.51	132.72	123.60
1	B	114	ARG	NH1-CZ-NH2	6.41	126.45	119.40
1	B	138	ARG	NE-CZ-NH1	6.37	123.49	120.30
1	B	161	ARG	NE-CZ-NH1	-6.23	117.19	120.30
1	B	45	ILE	O-C-N	-6.19	112.79	122.70
1	A	171	ARG	O-C-N	6.17	132.58	122.70
1	A	185	SER	CA-C-O	6.17	133.05	120.10
1	B	114	ARG	NE-CZ-NH2	-6.12	117.24	120.30
1	B	197	ASP	CB-CG-OD2	-6.10	112.81	118.30
1	B	206	ASP	O-C-N	6.09	132.45	122.70
1	A	43	ASP	CB-CG-OD1	5.92	123.63	118.30
1	A	98	TYR	CB-CG-CD2	5.91	124.55	121.00
1	B	32	PHE	CB-CG-CD2	-5.81	116.73	120.80
1	A	61	TYR	CB-CG-CD2	-5.81	117.51	121.00
1	A	61	TYR	CG-CD1-CE1	-5.69	116.75	121.30
1	B	128	THR	O-C-N	5.67	131.77	122.70
1	A	185	SER	O-C-N	-5.64	113.67	122.70
1	A	67	TYR	CG-CD1-CE1	5.62	125.79	121.30
1	A	215	ASP	CB-CG-OD2	-5.42	113.42	118.30
1	A	67	TYR	CB-CG-CD1	5.39	124.23	121.00
1	B	222	ASP	CB-CG-OD2	-5.38	113.45	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	114	ARG	NE-CZ-NH1	-5.28	117.66	120.30
1	A	52	VAL	CA-CB-CG2	-5.27	103.00	110.90
1	B	147	ASP	OD1-CG-OD2	-5.25	113.33	123.30
1	B	63	ILE	O-C-N	-5.20	114.38	122.70
1	B	196	GLU	OE1-CD-OE2	-5.18	117.08	123.30
1	A	101	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	A	219	VAL	CG1-CB-CG2	-5.16	102.64	110.90
1	B	15	ARG	NE-CZ-NH2	5.15	122.88	120.30
1	B	54	ASP	CB-CG-OD2	5.13	122.92	118.30
1	B	23	ASP	CB-CG-OD2	-5.12	113.69	118.30
1	A	139	ASP	CB-CG-OD1	5.11	122.90	118.30
1	A	102	GLU	OE1-CD-OE2	-5.05	117.24	123.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	195	ILE	Mainchain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1730	0	1716	23	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	44	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (44) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:300:XMP:H5'2	2:A:300:XMP:H8	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:C5'	2:A:300:XMP:H8	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:C8	2:B:300:XMP:H5'2	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: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:116:LYS:C	5:A:2101:HOH:O	2.48	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: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:44:LYS:HB2	1:A:219:VAL:HG22	1.96	0.47
1:A:9:TYR:HA	5:A:2070:HOH:O	2.14	0.47
1:A:229:GLU:O	1:A:230:LYS:C	2.53	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: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:CG	1:A:161:ARG:HH11	2.29	0.44
1:A:180:THR:HG22	1:A:197:ASP:CG	2.39	0.43
1:B:142:VAL:HG11	1:B:162:LEU:HD21	2.01	0.42
1:A:147:ASP:HB2	1:A:205:TYR:CE2	2.54	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:HA	1:B:155:LEU:HD12	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%)	183 (98%)	4 (2%)	66	37
1	B	188/206 (91%)	181 (96%)	7 (4%)	45	16
All	All	375/412 (91%)	364 (97%)	11 (3%)	55	24

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	102	GLU
1	A	114	ARG
1	A	161	ARG
1	B	4	LYS
1	B	23	ASP
1	B	30	ASP
1	B	51	LEU
1	B	114	ARG
1	B	155	LEU
1	B	203	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	ASN
1	A	96	GLN
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 chains 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 carbohydrates 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	1.17	3 (11%)	33,40,40	2.86	8 (24%)
3	POP	A	301	4	8,8,8	1.26	0	13,13,13	1.97	3 (23%)
2	XMP	B	300	-	26,26,26	1.31	4 (15%)	33,40,40	2.87	8 (24%)
3	POP	B	301	4	8,8,8	1.26	1 (12%)	13,13,13	1.98	6 (46%)

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	-	-	0/10/26/26	0/1/3/3
3	POP	A	301	4	-	0/6/6/6	0/0/0/0
2	XMP	B	300	-	-	0/10/26/26	0/1/3/3
3	POP	B	301	4	-	0/6/6/6	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	301	POP	P2-O	2.91	1.67	1.60
2	B	300	XMP	C4-N3	-2.87	1.33	1.37
2	A	300	XMP	C4-N3	-2.83	1.33	1.37
2	A	300	XMP	C5-C4	-2.31	1.36	1.40
2	B	300	XMP	P-O2P	-2.26	1.46	1.54
2	B	300	XMP	C8-N7	-2.23	1.31	1.34
2	A	300	XMP	O6-C6	-2.03	1.20	1.24
2	B	300	XMP	C8-N9	2.01	1.37	1.34

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	300	XMP	C5-C4-N3	9.13	130.52	121.29
2	B	300	XMP	C5-C4-N3	7.96	129.33	121.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	300	XMP	O4'-C1'-N9	-7.35	101.60	108.44
2	A	300	XMP	C4-C5-N7	7.29	112.01	106.07
2	B	300	XMP	C8-N7-C5	-6.09	104.19	108.52
2	A	300	XMP	C8-N7-C5	-5.91	104.31	108.52
2	B	300	XMP	C4-C5-N7	5.66	110.68	106.07
2	A	300	XMP	O4'-C1'-N9	-5.28	103.53	108.44
3	A	301	POP	O2-P1-O	4.93	119.98	104.29
2	B	300	XMP	C4'-O4'-C1'	-4.50	104.86	109.75
2	B	300	XMP	O2P-P-O1P	3.50	121.24	107.61
2	B	300	XMP	C2'-C3'-C4'	-3.46	95.77	102.65
3	A	301	POP	O3-P1-O	-3.44	96.05	106.65
3	B	301	POP	O2-P1-O	3.34	114.91	104.29
2	B	300	XMP	N7-C8-N9	3.22	112.73	109.18
3	B	301	POP	O3-P1-O	-3.07	97.18	106.65
3	B	301	POP	O-P2-O4	2.96	116.38	107.57
2	A	300	XMP	C4'-O4'-C1'	-2.87	106.63	109.75
3	B	301	POP	O6-P2-O	-2.70	98.31	106.65
2	A	300	XMP	C1'-N9-C4	-2.70	121.87	126.29
2	A	300	XMP	P-O5'-C5'	2.61	125.74	118.19
2	A	300	XMP	C8-N9-C4	2.48	109.02	107.13
3	B	301	POP	P2-O-P1	-2.47	124.33	131.74
3	A	301	POP	O6-P2-O	-2.12	100.11	106.65
3	B	301	POP	O5-P2-O4	2.08	118.34	110.82

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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	215/233 (92%)	0.44	9 (4%) 35 33	12, 21, 39, 50	0
1	B	217/233 (93%)	0.64	20 (9%) 9 8	14, 25, 44, 56	0
All	All	432/466 (92%)	0.54	29 (6%) 18 16	12, 23, 41, 56	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	127	LEU	6.6
1	A	180	THR	4.8
1	A	104	SER	4.0
1	B	3	SER	4.0
1	B	26	PHE	3.9
1	A	7	GLU	3.9
1	B	22	PRO	3.9
1	B	126	GLN	3.6
1	B	128	THR	3.6
1	B	185	SER	3.5
1	A	230	LYS	3.5
1	B	5	PRO	3.5
1	A	43	ASP	3.0
1	B	144	ILE	2.6
1	A	229	GLU	2.6
1	B	11	LYS	2.5
1	B	143	LEU	2.4
1	B	180	THR	2.4
1	B	12	GLY	2.4
1	A	116	LYS	2.4
1	B	153	PHE	2.3
1	B	172	ILE	2.3
1	A	1	MET	2.2
1	B	226	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	224	ALA	2.2
1	B	24	ASN	2.2
1	B	186	LEU	2.1
1	B	13	LYS	2.1
1	B	230	LYS	2.0

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	MG	A	303	1/1	0.15	7.74	21,21,21,21	0
2	XMP	A	300	24/24	0.11	0.29	20,28,34,35	0
3	POP	A	301	9/9	0.11	-0.00	17,20,23,23	0
3	POP	B	301	9/9	0.10	-0.07	21,23,26,27	0
4	MG	A	302	1/1	0.08	-0.39	29,29,29,29	0
2	XMP	B	300	24/24	0.10	-0.79	26,30,34,36	0
4	MG	B	302	1/1	0.07	-1.16	28,28,28,28	0
4	MG	B	303	1/1	0.08	-2.08	31,31,31,31	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.