



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 15, 2017 – 06:29 AM EDT

PDB ID : 5KNY  
Title : Crystal structure of Mycobacterium tuberculosis hypoxanthine guanine phosphoribosyltransferase in complex with (3-((3R,4R)-3-(Guanin-9-yl)-4-((S)-2-hydroxy-2-phosphonoethoxy)pyrrolidin-1-yl)-3-oxopropyl)phosphonic acid  
Authors : Eng, W.S.; Rejman, D.; Keough, D.T.; Guddat, L.W.  
Deposited on : unknown  
Resolution : 2.91 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20029824  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20029824

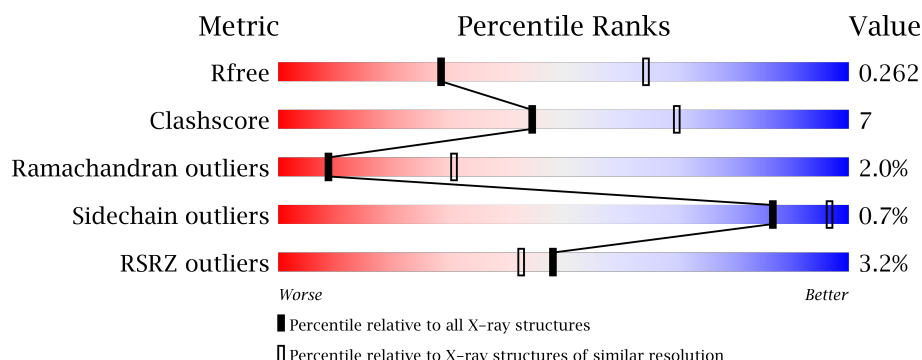
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.91 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))
$R_{free}$	100719	1813 (2.94-2.90)
Clashscore	112137	2045 (2.94-2.90)
Ramachandran outliers	110173	1997 (2.94-2.90)
Sidechain outliers	110143	1999 (2.94-2.90)
RSRZ outliers	101464	1825 (2.94-2.90)

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. 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	207	<div> <div>2%</div> <div>68% 17% 15%</div> </div>
1	B	207	<div> <div>6%</div> <div>75% 14% 10%</div> </div>
1	C	207	<div> <div>%</div> <div>71% 17% 12%</div> </div>
1	D	207	<div> <div>%</div> <div>70% 18% 12%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5698 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	176	Total	C	N	O	S	0	0	0
			1361	870	231	258	2			
1	B	187	Total	C	N	O	S	0	0	0
			1408	899	237	270	2			
1	C	183	Total	C	N	O	S	0	0	0
			1362	873	230	257	2			
1	D	183	Total	C	N	O	S	0	1	0
			1374	877	232	263	2			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	HIS	-	expression tag	UNP P9WHQ9
A	-3	HIS	-	expression tag	UNP P9WHQ9
A	-2	HIS	-	expression tag	UNP P9WHQ9
A	-1	HIS	-	expression tag	UNP P9WHQ9
A	0	HIS	-	expression tag	UNP P9WHQ9
A	1	HIS	-	expression tag	UNP P9WHQ9
B	-4	HIS	-	expression tag	UNP P9WHQ9
B	-3	HIS	-	expression tag	UNP P9WHQ9
B	-2	HIS	-	expression tag	UNP P9WHQ9
B	-1	HIS	-	expression tag	UNP P9WHQ9
B	0	HIS	-	expression tag	UNP P9WHQ9
B	1	HIS	-	expression tag	UNP P9WHQ9
C	-4	HIS	-	expression tag	UNP P9WHQ9
C	-3	HIS	-	expression tag	UNP P9WHQ9
C	-2	HIS	-	expression tag	UNP P9WHQ9
C	-1	HIS	-	expression tag	UNP P9WHQ9
C	0	HIS	-	expression tag	UNP P9WHQ9
C	1	HIS	-	expression tag	UNP P9WHQ9
D	-4	HIS	-	expression tag	UNP P9WHQ9
D	-3	HIS	-	expression tag	UNP P9WHQ9
D	-2	HIS	-	expression tag	UNP P9WHQ9

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	HIS	-	expression tag	UNP P9WHQ9
D	0	HIS	-	expression tag	UNP P9WHQ9
D	1	HIS	-	expression tag	UNP P9WHQ9

- 
- The chemical structure of YPG is a complex molecule. It features a purine-like core (a fused bicyclic system with nitrogen atoms N1, N2, N3, N7, N8, N9). The core is substituted with several groups: a phosphate group (OAF, OAT, OAH, OAD, OAE, OAJ, OAK, OAL, OAM, OAN, OAO, OAP, OAQ, OAR, OAS, OAT, OAU, OAV, OAW, OAX, OAY, OAZ) attached to the N9 position; a carbamate group (CBA(F)) attached to the N7 position; a phosphonate group (PBF) attached to the N2 position; and a phosphonate group (PBE) attached to the N1 position. The structure is highly detailed, showing various atoms (O, N, C, P, S, F) and their connectivity, with labels for each atom and bond type (single, double, triple, aromatic, etc.).

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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



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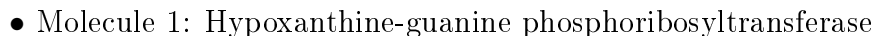
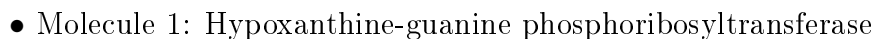
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total	Mg	0	0
			1	1		

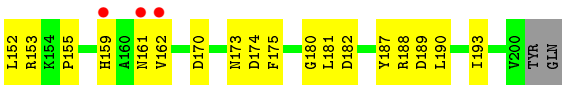
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	7	Total	O	0	0
			7	7		
4	B	10	Total	O	0	0
			10	10		
4	C	6	Total	O	0	0
			6	6		
4	D	6	Total	O	0	0
			6	6		

**i**

- Molecule 1: Hypoxanthine-guanine phosphoribosyltransferase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	54.52Å 86.08Å 79.82Å 90.00° 105.95° 90.00°	Depositor
Resolution (Å)	44.77 – 2.91 44.77 – 2.91	Depositor EDS
% Data completeness (in resolution range)	99.7 (44.77-2.91) 92.8 (44.77-2.91)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.00 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.201 , 0.262 0.203 , 0.262	Depositor DCC
$R_{free}$ test set	1428 reflections (9.83%)	DCC
Wilson B-factor (Å <sup>2</sup> )	38.9	Xtriage
Anisotropy	0.366	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 51.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	5698	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.31% 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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.22	0/1383	0.41	0/1882
1	B	0.24	0/1432	0.51	0/1957
1	C	0.26	0/1384	0.48	0/1891
1	D	0.24	0/1397	0.50	0/1910
All	All	0.24	0/5596	0.48	0/7640

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 [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1361	0	1354	21	0
1	B	1408	0	1370	19	0
1	C	1362	0	1338	25	0
1	D	1374	0	1344	21	0
2	A	64	0	0	3	0
2	B	32	0	0	1	0
2	C	32	0	0	6	0
2	D	32	0	0	1	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	7	0	0	0	0
4	B	10	0	0	1	0
4	C	6	0	0	0	0
4	D	6	0	0	1	0
All	All	5698	0	5406	83	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 7.

All (83) 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:B:57:GLN:HE21	1:B:116:ARG:HD2	1.50	0.77
1:B:185:GLU:OE2	1:B:188:ARG:NH2	2.23	0.72
1:D:153:ARG:NH1	1:D:170:ASP:OD1	2.27	0.67
1:A:182:ASP:O	2:A:301[A]:YPG:N2	2.29	0.66
1:A:152:LEU:HD11	1:A:193:ILE:HD13	1.78	0.64
1:B:102:VAL:HG11	1:C:141:ARG:HA	1.78	0.64
1:A:124:VAL:HG22	1:A:152:LEU:HB2	1.79	0.64
1:A:180:GLY:O	1:A:188:ARG:NH1	2.31	0.64
1:D:152:LEU:HD11	1:D:193:ILE:HD13	1.80	0.63
1:B:152:LEU:HD11	1:B:193:ILE:HD13	1.81	0.63
1:C:94:GLY:CA	1:C:96:SER:H	2.11	0.62
1:A:129:LEU:N	2:A:301[A]:YPG:OAG	2.32	0.62
1:C:176:VAL:O	2:C:301:YPG:N1	2.33	0.61
1:B:150:THR:HG22	1:B:164:ILE:HG21	1.82	0.60
1:C:130:THR:OG1	2:C:301:YPG:OAG	2.13	0.58
1:C:127:SER:OG	2:C:301:YPG:OAD	2.15	0.58
1:D:174:ASP:OD1	1:D:175:PHE:N	2.37	0.57
1:D:182:ASP:O	2:D:301:YPG:N2	2.38	0.57
1:B:27:LEU:HD22	1:B:176:VAL:HG11	1.86	0.57
1:D:29:THR:HG23	1:D:32:GLN:H	1.71	0.55
1:C:186:ARG:HG2	1:D:84:GLN:HE21	1.71	0.55
1:C:161:ASN:OD1	1:C:161:ASN:N	2.38	0.55
1:C:94:GLY:HA3	1:C:96:SER:H	1.72	0.55
1:B:65:LEU:HD11	1:B:87:PHE:HB3	1.89	0.54
1:A:178:GLY:HA2	1:A:190:LEU:HB2	1.89	0.54
1:D:27:LEU:HD22	1:D:28:LEU:HG	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:105:ILE:HD11	1:D:133:TRP:HH2	1.72	0.53
1:D:180:GLY:O	1:D:188:ARG:NH1	2.41	0.53
1:D:97:THR:HG22	1:D:102:VAL:O	2.09	0.53
1:C:186:ARG:NH2	1:C:187:TYR:OH	2.42	0.53
1:D:141:ARG:NH1	4:D:402:HOH:O	2.43	0.52
1:C:63:THR:OG1	1:C:72:VAL:HG11	2.10	0.52
1:D:93:TYR:H	1:D:97:THR:CG2	2.23	0.51
1:A:186:ARG:HG2	1:B:84:GLN:HE21	1.76	0.51
1:C:128:GLY:N	2:C:301:YPG:OAD	2.41	0.51
1:D:63:THR:OG1	1:D:72:VAL:HG11	2.12	0.50
1:A:69:VAL:HG23	1:A:70:LEU:HD12	1.92	0.50
1:C:128:GLY:H	2:C:301:YPG:PBE	2.35	0.49
1:B:57:GLN:NE2	1:B:116:ARG:HD2	2.23	0.49
1:C:60:LEU:HD11	1:C:86:GLU:HB2	1.96	0.48
1:A:152:LEU:HD13	1:A:171:ILE:HD13	1.96	0.47
1:B:185:GLU:CD	1:B:188:ARG:HH22	2.18	0.47
1:C:182:ASP:O	2:C:301:YPG:N2	2.47	0.47
1:A:198:PRO:HA	1:A:201:TYR:CD2	2.49	0.47
1:B:65:LEU:HA	1:B:66:LYS:HA	1.58	0.46
1:B:17:GLU:O	1:B:18:LEU:HB2	2.15	0.46
1:D:92:SER:HA	1:D:97:THR:HG21	1.97	0.46
1:A:27:LEU:H	1:A:194:GLY:HA2	1.80	0.46
1:A:197:ASP:OD1	1:A:198:PRO:HD2	2.15	0.46
1:B:79:ILE:HG22	1:B:81:VAL:H	1.81	0.46
1:D:65:LEU:HB3	1:D:66:LYS:HG2	1.98	0.45
1:C:72:VAL:HG12	1:C:121:VAL:HG11	1.99	0.45
1:B:102:VAL:HA	1:B:133:TRP:CD1	2.53	0.44
1:C:132:SER:OG	1:C:161:ASN:ND2	2.50	0.44
1:C:180:GLY:O	1:C:188:ARG:NH1	2.50	0.44
1:D:65:LEU:HA	1:D:66:LYS:HA	1.63	0.44
1:A:59:LEU:HB3	1:A:83:THR:HG22	1.99	0.44
1:D:189:ASP:OD1	1:D:189:ASP:N	2.51	0.44
1:C:65:LEU:HA	1:C:66:LYS:HA	1.62	0.44
1:A:86:GLU:HG2	1:A:109:LEU:HD22	1.99	0.44
1:A:65:LEU:HA	1:A:66:LYS:HA	1.59	0.44
1:C:22:ASP:O	1:C:196:LEU:HA	2.18	0.43
1:A:137:ASN:O	1:A:140:SER:OG	2.26	0.43
1:B:59:LEU:HB3	1:B:83:THR:HG22	2.01	0.43
1:C:174:ASP:O	1:C:176:VAL:HG23	2.19	0.42
1:C:174:ASP:OD1	1:C:175:PHE:N	2.52	0.42
1:A:18:LEU:HD23	1:A:190:LEU:HD23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:102:VAL:HA	1:C:133:TRP:CD1	2.55	0.42
1:D:152:LEU:HD12	1:D:181:LEU:HD11	2.02	0.42
1:A:128:GLY:N	2:A:301[B]:YPG:OAJ	2.46	0.41
1:A:68:ALA:HB2	1:A:123:ASP:HB2	2.02	0.41
1:C:18:LEU:HB3	1:C:19:TYR:H	1.75	0.41
1:A:125:VAL:HG11	1:A:164:ILE:HD11	2.02	0.41
1:B:107:LYS:NZ	4:B:403:HOH:O	2.41	0.41
1:C:45:GLY:HA3	1:C:80:PRO:HD2	2.01	0.41
1:C:94:GLY:HA3	1:C:96:SER:N	2.34	0.41
1:D:72:VAL:HG12	1:D:121:VAL:HG11	2.03	0.41
1:D:187:TYR:HA	1:D:190:LEU:HD13	2.03	0.41
1:B:86:GLU:HG2	1:B:109:LEU:HD22	2.02	0.41
1:B:176:VAL:O	2:B:301:YPG:N1	2.53	0.41
1:D:155:PRO:HG2	1:D:173:ASN:HB2	2.02	0.41
1:B:95:SER:C	1:B:97:THR:H	2.23	0.40
1:A:75:LEU:O	1:A:79:ILE:HG13	2.22	0.40

There are no symmetry-related clashes.

## 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	170/207 (82%)	163 (96%)	7 (4%)	0	100	100
1	B	185/207 (89%)	169 (91%)	10 (5%)	6 (3%)	5	18
1	C	181/207 (87%)	172 (95%)	5 (3%)	4 (2%)	8	28
1	D	182/207 (88%)	171 (94%)	7 (4%)	4 (2%)	8	28
All	All	718/828 (87%)	675 (94%)	29 (4%)	14 (2%)	9	32

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	58	ASP
1	C	197	ASP
1	C	198	PRO
1	D	95	SER
1	D	162	VAL
1	B	49	ARG
1	C	99	SER
1	D	161	ASN
1	B	50	GLU
1	B	51	LEU
1	B	102	VAL
1	D	159	HIS
1	B	17	GLU
1	C	98	SER

### 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	147/180 (82%)	146 (99%)	1 (1%)	87	96
1	B	146/180 (81%)	146 (100%)	0	100	100
1	C	140/180 (78%)	139 (99%)	1 (1%)	87	96
1	D	144/180 (80%)	142 (99%)	2 (1%)	71	91
All	All	577/720 (80%)	573 (99%)	4 (1%)	87	96

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	VAL
1	C	103	VAL
1	D	90	VAL
1	D	109	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	57	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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	YPG	A	301[A]	3	29,34,34	2.68	8 (27%)	31,52,52	2.07	8 (25%)
2	YPG	A	301[B]	3	29,34,34	2.81	8 (27%)	31,52,52	2.26	10 (32%)
2	YPG	B	301	3	29,34,34	2.67	8 (27%)	31,52,52	2.17	9 (29%)
2	YPG	C	301	3	29,34,34	2.67	8 (27%)	31,52,52	2.09	9 (29%)
2	YPG	D	301	3	29,34,34	2.69	8 (27%)	31,52,52	2.05	7 (22%)

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	YPG	A	301[A]	3	-	0/15/37/37	0/3/3/3
2	YPG	A	301[B]	3	-	0/15/37/37	0/3/3/3
2	YPG	B	301	3	-	0/15/37/37	0/3/3/3
2	YPG	C	301	3	-	0/15/37/37	0/3/3/3
2	YPG	D	301	3	-	0/15/37/37	0/3/3/3

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	YPG	CBA-CBB	-3.30	1.44	1.52
2	C	301	YPG	CBA-CBB	-3.23	1.44	1.52
2	A	301[B]	YPG	CBA-CBB	-3.01	1.45	1.52
2	B	301	YPG	CBA-CBB	-2.83	1.45	1.52
2	A	301[A]	YPG	CBA-CBB	-2.82	1.45	1.52
2	A	301[A]	YPG	C6-N1	2.95	1.38	1.33
2	C	301	YPG	C6-N1	3.00	1.38	1.33
2	A	301[B]	YPG	C2-N1	3.01	1.40	1.35
2	A	301[A]	YPG	C2-N1	3.05	1.40	1.35
2	B	301	YPG	C2-N1	3.06	1.40	1.35
2	D	301	YPG	C2-N1	3.06	1.40	1.35
2	C	301	YPG	C2-N1	3.07	1.40	1.35
2	D	301	YPG	C6-N1	3.10	1.38	1.33
2	A	301[B]	YPG	C6-N1	3.11	1.38	1.33
2	B	301	YPG	C6-N1	3.12	1.38	1.33
2	B	301	YPG	CAO-NBC	3.29	1.54	1.46
2	A	301[A]	YPG	CAO-NBC	3.40	1.54	1.46
2	C	301	YPG	CAO-NBC	3.53	1.54	1.46
2	D	301	YPG	CAO-NBC	3.54	1.54	1.46
2	A	301[B]	YPG	CAO-NBC	3.80	1.55	1.46
2	D	301	YPG	CAU-NBC	4.79	1.45	1.35
2	C	301	YPG	CAU-NBC	4.81	1.45	1.35
2	A	301[A]	YPG	CAU-NBC	4.88	1.45	1.35
2	A	301[B]	YPG	CAU-NBC	4.90	1.45	1.35
2	B	301	YPG	CAU-NBC	4.97	1.46	1.35
2	A	301[A]	YPG	C2-N2	5.01	1.44	1.34
2	C	301	YPG	C2-N2	5.08	1.44	1.34
2	B	301	YPG	C2-N2	5.09	1.44	1.34
2	D	301	YPG	C2-N2	5.09	1.44	1.34
2	A	301[B]	YPG	C2-N2	5.14	1.44	1.34
2	A	301[B]	YPG	O6-C6	6.35	1.40	1.24
2	C	301	YPG	O6-C6	6.36	1.40	1.24
2	A	301[A]	YPG	O6-C6	6.37	1.40	1.24
2	D	301	YPG	O6-C6	6.40	1.40	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	YPG	O6-C6	6.41	1.40	1.24
2	C	301	YPG	PBE-CAM	6.97	1.85	1.78
2	B	301	YPG	PBE-CAM	7.10	1.86	1.78
2	D	301	YPG	PBE-CAM	7.11	1.86	1.78
2	A	301[A]	YPG	PBE-CAM	7.35	1.86	1.78
2	A	301[B]	YPG	PBE-CAM	8.22	1.87	1.78

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301[B]	YPG	N3-C2-N1	-6.48	118.00	127.46
2	B	301	YPG	N3-C2-N1	-6.32	118.23	127.46
2	D	301	YPG	N3-C2-N1	-6.30	118.26	127.46
2	C	301	YPG	N3-C2-N1	-6.25	118.33	127.46
2	A	301[A]	YPG	N3-C2-N1	-6.03	118.65	127.46
2	A	301[A]	YPG	CAP-NBC-CAO	-3.36	105.90	112.79
2	B	301	YPG	C4-C5-N7	-3.32	106.20	109.41
2	A	301[B]	YPG	C4-C5-N7	-3.31	106.21	109.41
2	C	301	YPG	C4-C5-N7	-3.22	106.30	109.41
2	D	301	YPG	C4-C5-N7	-3.21	106.31	109.41
2	A	301[B]	YPG	CAP-NBC-CAO	-3.09	106.45	112.79
2	B	301	YPG	CAP-NBC-CAO	-3.09	106.46	112.79
2	A	301[A]	YPG	C4-C5-N7	-3.03	106.48	109.41
2	C	301	YPG	OAT-CBA-CAO	-2.69	106.50	112.02
2	A	301[A]	YPG	C5-C6-N1	-2.65	119.70	123.48
2	D	301	YPG	C5-C6-N1	-2.62	119.75	123.48
2	C	301	YPG	C5-C6-N1	-2.59	119.80	123.48
2	B	301	YPG	C5-C6-N1	-2.55	119.86	123.48
2	A	301[B]	YPG	C5-C6-N1	-2.39	120.08	123.48
2	B	301	YPG	CAM-CAL-CAU	-2.26	108.12	111.91
2	B	301	YPG	CAP-CBB-N9	-2.24	108.81	113.33
2	C	301	YPG	CAM-CAL-CAU	-2.22	108.18	111.91
2	D	301	YPG	OAT-CBA-CAO	-2.15	107.61	112.02
2	A	301[B]	YPG	CAM-CAL-CAU	-2.09	108.41	111.91
2	C	301	YPG	CAP-NBC-CAO	-2.06	108.57	112.79
2	A	301[A]	YPG	N2-C2-N1	2.05	120.52	117.24
2	C	301	YPG	N2-C2-N1	2.08	120.56	117.24
2	A	301[B]	YPG	CAO-NBC-CAU	2.08	129.29	123.85
2	A	301[A]	YPG	CAN-OAT-CBA	2.11	118.27	115.18
2	B	301	YPG	N2-C2-N1	2.18	120.73	117.24
2	D	301	YPG	N2-C2-N1	2.34	120.98	117.24
2	A	301[B]	YPG	CAL-CAU-NBC	2.37	121.99	117.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301[A]	YPG	C6-N1-C2	2.43	119.55	116.06
2	A	301[B]	YPG	N2-C2-N1	2.45	121.16	117.24
2	A	301[B]	YPG	C6-N1-C2	2.45	119.58	116.06
2	B	301	YPG	C6-N1-C2	2.53	119.70	116.06
2	C	301	YPG	C6-N1-C2	2.55	119.73	116.06
2	D	301	YPG	C6-N1-C2	2.58	119.77	116.06
2	A	301[A]	YPG	C2-N3-C4	5.06	121.07	115.16
2	D	301	YPG	C2-N3-C4	5.19	121.22	115.16
2	C	301	YPG	C2-N3-C4	5.27	121.31	115.16
2	B	301	YPG	C2-N3-C4	5.29	121.34	115.16
2	A	301[B]	YPG	C2-N3-C4	5.84	121.98	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301[A]	YPG	2	0
2	A	301[B]	YPG	1	0
2	B	301	YPG	1	0
2	C	301	YPG	6	0
2	D	301	YPG	1	0

## 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, 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	176/207 (85%)	-0.10	4 (2%) 61 58	23, 39, 79, 105	0
1	B	187/207 (90%)	0.14	13 (6%) 17 13	26, 41, 100, 121	0
1	C	183/207 (88%)	0.25	3 (1%) 72 70	32, 56, 90, 108	0
1	D	183/207 (88%)	0.20	3 (1%) 72 70	32, 63, 96, 123	0
All	All	729/828 (88%)	0.12	23 (3%) 48 42	23, 51, 92, 123	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	52	SER	6.5
1	B	51	LEU	5.3
1	B	20	PRO	4.5
1	B	98	SER	4.4
1	B	100	SER	4.3
1	D	159	HIS	4.0
1	B	99	SER	3.4
1	B	53	ALA	3.1
1	B	50	GLU	3.0
1	B	56	GLY	3.0
1	B	55	THR	2.6
1	D	161	ASN	2.6
1	B	54	THR	2.6
1	C	95	SER	2.4
1	C	200	VAL	2.3
1	A	161	ASN	2.3
1	A	201	TYR	2.3
1	B	200	VAL	2.3
1	D	162	VAL	2.2
1	B	19	TYR	2.1
1	A	20	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	19	TYR	2.1
1	C	98	SER	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 carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	YPG	A	301[B]	32/32	0.90	0.21	0.58	30,41,75,79	32
2	YPG	A	301[A]	32/32	0.90	0.21	0.55	32,40,79,80	32
2	YPG	B	301	32/32	0.93	0.16	-0.58	12,35,86,88	0
2	YPG	C	301	32/32	0.92	0.16	-0.99	38,50,95,96	0
2	YPG	D	301	32/32	0.93	0.15	-1.38	44,49,73,74	9
3	MG	B	302	1/1	0.95	0.07	-	63,63,63,63	0
3	MG	C	302	1/1	0.93	0.08	-	32,32,32,32	0
3	MG	A	302	1/1	0.99	0.08	-	33,33,33,33	0
3	MG	D	302	1/1	0.81	0.17	-	60,60,60,60	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.