



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 3, 2023 – 09:34 pm GMT

PDB ID : 2C46  
Title : CRYSTAL STRUCTURE OF THE HUMAN RNA guanylyltransferase and 5'- phosphatase  
Authors : Debreczeni, J.; Johansson, C.; Longman, E.; Gileadi, O.; SavitskySmee, P.; Smee, C.; Bunkoczi, G.; Ugochukwu, E.; von Delft, F.; Sundstrom, M.; Weigelt, J.; Arrowsmith, C.; Edwards, A.; Knapp, S.  
Deposited on : 2005-10-15  
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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

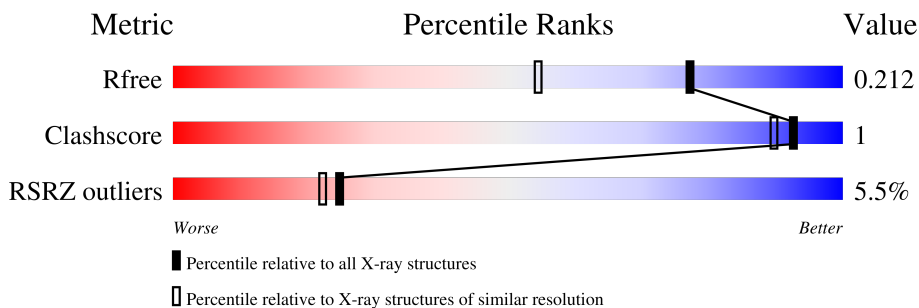
# 1 Overall quality at a glance

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))
$R_{free}$	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
RSRZ outliers	127900	3321 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	241	 2%                      79%                      17%
1	B	241	 2%                      80%                      17%
1	C	241	 3%                      80%                      18%
1	D	241	 10%                      74%                      5%                      22%

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 6719 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 MRNA CAPPING ENZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	200	Total 1603	C 1030	N 274	O 287	S 12	0	2	0
1	B	201	Total 1628	C 1041	N 274	O 301	S 12	0	5	0
1	C	197	Total 1553	C 995	N 262	O 285	S 11	0	0	0
1	D	189	Total 1475	C 946	N 244	O 274	S 11	0	0	0

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	expression tag	UNP O60942
A	-20	HIS	-	expression tag	UNP O60942
A	-19	HIS	-	expression tag	UNP O60942
A	-18	HIS	-	expression tag	UNP O60942
A	-17	HIS	-	expression tag	UNP O60942
A	-16	HIS	-	expression tag	UNP O60942
A	-15	HIS	-	expression tag	UNP O60942
A	-14	SER	-	expression tag	UNP O60942
A	-13	SER	-	expression tag	UNP O60942
A	-12	GLY	-	expression tag	UNP O60942
A	-11	VAL	-	expression tag	UNP O60942
A	-10	ASP	-	expression tag	UNP O60942
A	-9	LEU	-	expression tag	UNP O60942
A	-8	GLY	-	expression tag	UNP O60942
A	-7	THR	-	expression tag	UNP O60942
A	-6	GLU	-	expression tag	UNP O60942
A	-5	ASN	-	expression tag	UNP O60942
A	-4	LEU	-	expression tag	UNP O60942
A	-3	TYR	-	expression tag	UNP O60942
A	-2	PHE	-	expression tag	UNP O60942
A	-1	GLN	-	expression tag	UNP O60942

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP O60942
A	212	THR	PRO	conflict	UNP O60942
B	-21	MET	-	expression tag	UNP O60942
B	-20	HIS	-	expression tag	UNP O60942
B	-19	HIS	-	expression tag	UNP O60942
B	-18	HIS	-	expression tag	UNP O60942
B	-17	HIS	-	expression tag	UNP O60942
B	-16	HIS	-	expression tag	UNP O60942
B	-15	HIS	-	expression tag	UNP O60942
B	-14	SER	-	expression tag	UNP O60942
B	-13	SER	-	expression tag	UNP O60942
B	-12	GLY	-	expression tag	UNP O60942
B	-11	VAL	-	expression tag	UNP O60942
B	-10	ASP	-	expression tag	UNP O60942
B	-9	LEU	-	expression tag	UNP O60942
B	-8	GLY	-	expression tag	UNP O60942
B	-7	THR	-	expression tag	UNP O60942
B	-6	GLU	-	expression tag	UNP O60942
B	-5	ASN	-	expression tag	UNP O60942
B	-4	LEU	-	expression tag	UNP O60942
B	-3	TYR	-	expression tag	UNP O60942
B	-2	PHE	-	expression tag	UNP O60942
B	-1	GLN	-	expression tag	UNP O60942
B	0	SER	-	expression tag	UNP O60942
B	212	THR	PRO	conflict	UNP O60942
C	-21	MET	-	expression tag	UNP O60942
C	-20	HIS	-	expression tag	UNP O60942
C	-19	HIS	-	expression tag	UNP O60942
C	-18	HIS	-	expression tag	UNP O60942
C	-17	HIS	-	expression tag	UNP O60942
C	-16	HIS	-	expression tag	UNP O60942
C	-15	HIS	-	expression tag	UNP O60942
C	-14	SER	-	expression tag	UNP O60942
C	-13	SER	-	expression tag	UNP O60942
C	-12	GLY	-	expression tag	UNP O60942
C	-11	VAL	-	expression tag	UNP O60942
C	-10	ASP	-	expression tag	UNP O60942
C	-9	LEU	-	expression tag	UNP O60942
C	-8	GLY	-	expression tag	UNP O60942
C	-7	THR	-	expression tag	UNP O60942
C	-6	GLU	-	expression tag	UNP O60942
C	-5	ASN	-	expression tag	UNP O60942

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	LEU	-	expression tag	UNP O60942
C	-3	TYR	-	expression tag	UNP O60942
C	-2	PHE	-	expression tag	UNP O60942
C	-1	GLN	-	expression tag	UNP O60942
C	0	SER	-	expression tag	UNP O60942
C	212	THR	PRO	conflict	UNP O60942
D	-21	MET	-	expression tag	UNP O60942
D	-20	HIS	-	expression tag	UNP O60942
D	-19	HIS	-	expression tag	UNP O60942
D	-18	HIS	-	expression tag	UNP O60942
D	-17	HIS	-	expression tag	UNP O60942
D	-16	HIS	-	expression tag	UNP O60942
D	-15	HIS	-	expression tag	UNP O60942
D	-14	SER	-	expression tag	UNP O60942
D	-13	SER	-	expression tag	UNP O60942
D	-12	GLY	-	expression tag	UNP O60942
D	-11	VAL	-	expression tag	UNP O60942
D	-10	ASP	-	expression tag	UNP O60942
D	-9	LEU	-	expression tag	UNP O60942
D	-8	GLY	-	expression tag	UNP O60942
D	-7	THR	-	expression tag	UNP O60942
D	-6	GLU	-	expression tag	UNP O60942
D	-5	ASN	-	expression tag	UNP O60942
D	-4	LEU	-	expression tag	UNP O60942
D	-3	TYR	-	expression tag	UNP O60942
D	-2	PHE	-	expression tag	UNP O60942
D	-1	GLN	-	expression tag	UNP O60942
D	0	SER	-	expression tag	UNP O60942
D	212	THR	PRO	conflict	UNP O60942

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	146	Total O 146 146	0	0
2	B	147	Total O 147 147	0	0
2	C	83	Total O 83 83	0	0
2	D	84	Total O 84 84	0	0





## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	60.40Å 161.60Å 60.70Å 90.00° 104.70° 90.00°	Depositor
Resolution (Å)	80.85 – 1.60 47.94 – 1.60	Depositor EDS
% Data completeness (in resolution range)	90.4 (80.85-1.60) 89.1 (47.94-1.60)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.07 (at 1.60Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.204 , 0.234 0.202 , 0.212	Depositor DCC
$R_{free}$ test set	6714 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	19.9	Xtrriage
Anisotropy	0.098	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 48.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtrriage
Estimated twinning fraction	0.116 for l,-k,h	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6719	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% 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](#)

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	0.58	0/1654	0.70	1/2242 (0.0%)
1	B	0.63	1/1683 (0.1%)	0.71	2/2281 (0.1%)
1	C	0.47	0/1594	0.61	0/2163
1	D	0.43	0/1515	0.59	0/2060
All	All	0.54	1/6446 (0.0%)	0.66	3/8746 (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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	138	CYS	CB-SG	-5.03	1.73	1.81

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	175	ARG	NE-CZ-NH2	-5.83	117.38	120.30
1	B	175	ARG	NE-CZ-NH1	5.73	123.16	120.30
1	A	199	ASP	CB-CG-OD2	5.13	122.92	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	113	PHE	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	1603	0	1542	5	0
1	B	1628	0	1563	4	0
1	C	1553	0	1464	2	0
1	D	1475	0	1373	7	0
2	A	146	0	0	0	0
2	B	147	0	0	1	0
2	C	83	0	0	0	0
2	D	84	0	0	0	0
All	All	6719	0	5942	18	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 1.

All (18) 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:A:99:THR:O	1:A:103:THR:HG23	1.85	0.76
1:D:75:ASP:OD1	1:D:77:ASN:ND2	2.39	0.54
1:D:46:HIS:ND1	1:D:46:HIS:N	2.44	0.54
1:B:113:PHE:HB2	1:B:122:ILE:HD11	1.91	0.53
1:D:174:PHE:CE1	1:D:184:PRO:HD3	2.46	0.51
1:A:1:MET:HE2	1:A:34:ARG:CB	2.42	0.50
1:A:27:LEU:HG	1:A:125:HIS:HB3	1.93	0.50
1:B:96:GLU:OE1	2:B:2074:HOH:O	2.19	0.49
1:D:63:LEU:HB2	1:D:113:PHE:CZ	2.49	0.48
1:A:1:MET:CE	1:A:34:ARG:CB	2.92	0.47
1:C:99:THR:O	1:C:103:THR:HG23	2.17	0.45
1:B:87[B]:ILE:HG21	1:B:109:LEU:HD21	1.99	0.44
1:D:79:ILE:HG22	1:D:84:ILE:HB	2.00	0.43
1:C:85:LYS:HZ3	1:C:87:ILE:CG1	2.32	0.42
1:A:113:PHE:HB2	1:A:122:ILE:HD11	2.03	0.41
1:D:179:ASP:HB3	1:D:182:GLU:HG3	2.02	0.41
1:B:63:LEU:HD11	1:B:87[B]:ILE:HG22	2.03	0.40
1:D:179:ASP:HB3	1:D:182:GLU:CG	2.51	0.40

There are no symmetry-related clashes.

### 5.3 Torsion angles [i](#)

#### 5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

#### 5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

#### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 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	200/241 (82%)	0.48	5 (2%) 57 55	21, 30, 42, 47	0
1	B	201/241 (83%)	0.47	6 (2%) 50 48	21, 30, 42, 52	0
1	C	197/241 (81%)	0.49	7 (3%) 42 40	21, 31, 39, 45	0
1	D	189/241 (78%)	1.00	25 (13%) 3 2	22, 32, 39, 46	0
All	All	787/964 (81%)	0.60	43 (5%) 25 22	21, 31, 41, 52	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	ALA	6.0
1	D	6	ILE	5.7
1	A	1	MET	4.9
1	C	0	SER	4.4
1	A	3	HIS	4.4
1	D	32	GLY	4.4
1	D	79	ILE	4.3
1	D	4	ASN	4.2
1	D	22	GLY	4.2
1	D	46	HIS	4.2
1	B	0	SER	4.2
1	D	31	LEU	4.1
1	D	21	ALA	4.0
1	D	33	PRO	3.7
1	D	53	TYR	3.7
1	D	109	LEU	3.6
1	D	7	PRO	3.5
1	B	53	TYR	3.2
1	A	0	SER	3.2
1	D	9	ARG	3.0
1	D	10	TRP	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	81	LYS	3.0
1	C	4	ASN	2.9
1	D	77	ASN	2.9
1	D	73	PHE	2.7
1	D	87	ILE	2.7
1	C	-1	GLN	2.6
1	B	109	LEU	2.6
1	C	198	GLU	2.6
1	C	57	LEU	2.5
1	D	200	GLU	2.5
1	B	59	VAL	2.5
1	D	11	LEU	2.4
1	A	109	LEU	2.4
1	D	78	ASP	2.3
1	D	23	ARG	2.3
1	C	110	CYS	2.3
1	D	107	ILE	2.3
1	C	53	TYR	2.2
1	B	-6	GLU	2.2
1	D	26	PRO	2.1
1	D	5	LYS	2.1
1	B	-1	GLN	2.0

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

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

## 6.3 Carbohydrates [\(i\)](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [\(i\)](#)

There are no ligands in this entry.

## 6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.