



# Full wwPDB X-ray Structure Validation Report ⓘ

Jul 22, 2025 – 01:17 PM JST

PDB ID : 9MBE / pdb\_00009mbe  
Title : Neutron crystal structure of human MTH1(G2K/C87A/C104S mutant) in complex with 8-oxo-dGTP  
Authors : Fujimiya, K.; Hirata, K.; Ostermann, A.; Schrader, T.E.; Hiromoto, T.; Ari-mori, T.; Hirano, Y.; Tamada, T.; Nakamura, T.  
Deposited on : 2025-03-17  
Resolution : 1.40 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>.

---

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

MolProbity : 4-5-2 with Phenix2.0rc1  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 2.0rc1  
EDS : **FAILED**  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.44

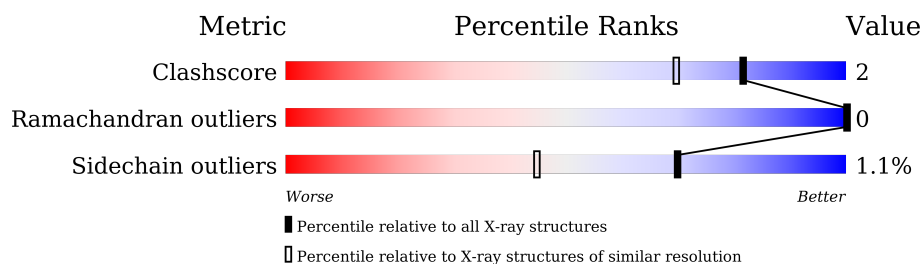
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION, NEUTRON DIFFRACTION*

The reported resolution of this entry is 1.40 Å.

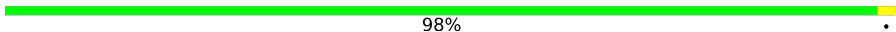
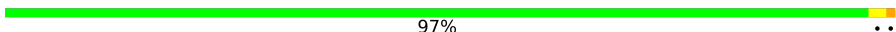
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	180529	2446 (1.40-1.40)
Ramachandran outliers	177936	2398 (1.40-1.40)
Sidechain outliers	177891	2397 (1.40-1.40)

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  $\leq 5\%$ .

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	156	 98% .
1	B	156	 97% ..

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6178 atoms, of which 2603 are hydrogens and 565 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.

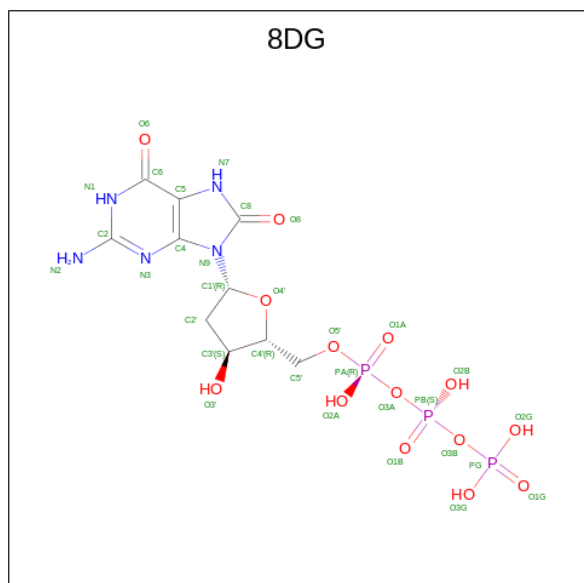
- Molecule 1 is a protein called 7,8-dihydro-8-oxoguanine triphosphatase.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	156	Total	C	D	H	N	O	S	0	132	0
			2842	848	229	1295	218	246	6			
1	B	156	Total	C	D	H	N	O	S	0	133	0
			2848	847	242	1289	219	244	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	LYS	GLY	engineered mutation	UNP P36639
A	87	ALA	CYS	engineered mutation	UNP P36639
A	104	SER	CYS	engineered mutation	UNP P36639
B	2	LYS	GLY	engineered mutation	UNP P36639
B	87	ALA	CYS	engineered mutation	UNP P36639
B	104	SER	CYS	engineered mutation	UNP P36639

- Molecule 2 is 8-OXO-2'-DEOXYGUANOSINE-5'-TRIPHOSPHATE (CCD ID: 8DG) (formula:  $C_{10}H_{16}N_5O_{14}P_3$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
2	A	1	Total	C	D	H	N	O	P	0	1
			49	10	5	12	5	14	3		
2	B	1	Total	C	H	N	O	P		0	0
			39	10	7	5	14	3			

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Na	0	0
			2	2		
3	B	1	Total	Na	0	1
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	159	Total	D	O	0	4
			223	60	163		
4	B	141	Total	D	O	0	3
			173	29	144		

### 3 Residue-property plots

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: 7,8-dihydro-8-oxoguanine triphosphatase

Chain A:  98% .



- Molecule 1: 7,8-dihydro-8-oxoguanine triphosphatase

Chain B:  97% ..



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.85Å 48.16Å 124.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.08 – 1.40	Depositor
% Data completeness (in resolution range)	97.8 (32.08-1.40)	Depositor
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	7.64 (at 1.40Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.174 , 0.197	Depositor
Wilson B-factor (Å <sup>2</sup> )	9.8	Xtriage
Anisotropy	0.360	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for k,h,-l	Xtriage
Total number of atoms	6178	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.81% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 8DG, NA

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	0.73	0/2473	0.74	0/3328
1	B	0.68	0/2478	0.80	0/3333
All	All	0.70	0/4951	0.77	0/6661

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

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	1547	1295	315	1	0
1	B	1559	1289	279	6	0
2	A	37	12	0	0	0
2	B	32	7	12	2	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	223	0	0	1	0
4	B	173	0	0	2	0
All	All	3575	2603	606	11	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 2.

All (11) 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:149[D]:THR:HG23	4:A:345:HOH:O	2.09	0.48
1:B:124:PHE:N	1:B:125:PRO:CD	2.85	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	292/156 (187%)	290 (99%)	2 (1%)	0	100	100
1	B	293/156 (188%)	291 (99%)	2 (1%)	0	100	100
All	All	585/312 (188%)	581 (99%)	4 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	256/136 (188%)	252 (98%)	4 (2%)	58	29
1	B	258/136 (190%)	257 (100%)	1 (0%)	89	76
All	All	514/272 (189%)	509 (99%)	5 (1%)	70	49



All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16[A]	GLN
1	A	16[B]	GLN
1	A	25[A]	ARG
1	A	25[B]	ARG
1	B	124	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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, 4 are monoatomic - leaving 3 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	8DG	B	201	3	30,34,34	1.46	3 (10%)	43,54,54	2.53	13 (30%)

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	8DG	B	201	3	-	6/22/34/34	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	201	8DG	C8-N9	-4.54	1.32	1.40
2	B	201	8DG	C4-N9	-3.98	1.31	1.39
2	B	201	8DG	O8-C8	2.48	1.27	1.23

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	8DG	N7-C8-N9	7.68	115.69	106.58
2	B	201	8DG	C5-C4-N3	-6.55	121.04	127.80
2	B	201	8DG	C5-N7-C8	-5.42	101.67	109.47
2	B	201	8DG	C2-N3-C4	4.72	120.71	112.30
2	B	201	8DG	O6-C6-C5	-4.45	117.03	127.24
2	B	201	8DG	N9-C4-N3	3.88	130.24	125.81
2	B	201	8DG	C2-N1-C6	-3.49	118.74	125.10
2	B	201	8DG	O8-C8-N9	-3.44	121.19	125.99
2	B	201	8DG	C2'-C1'-N9	-2.81	112.69	116.01
2	B	201	8DG	PA-O3A-PB	-2.40	124.58	132.83
2	B	201	8DG	C5-C6-N1	2.39	119.44	112.31
2	B	201	8DG	N2-C2-N1	2.25	121.51	116.71
2	B	201	8DG	N2-C2-N3	-2.01	115.82	119.73

There are no chirality outliers.

All (6) torsion outliers are listed below:

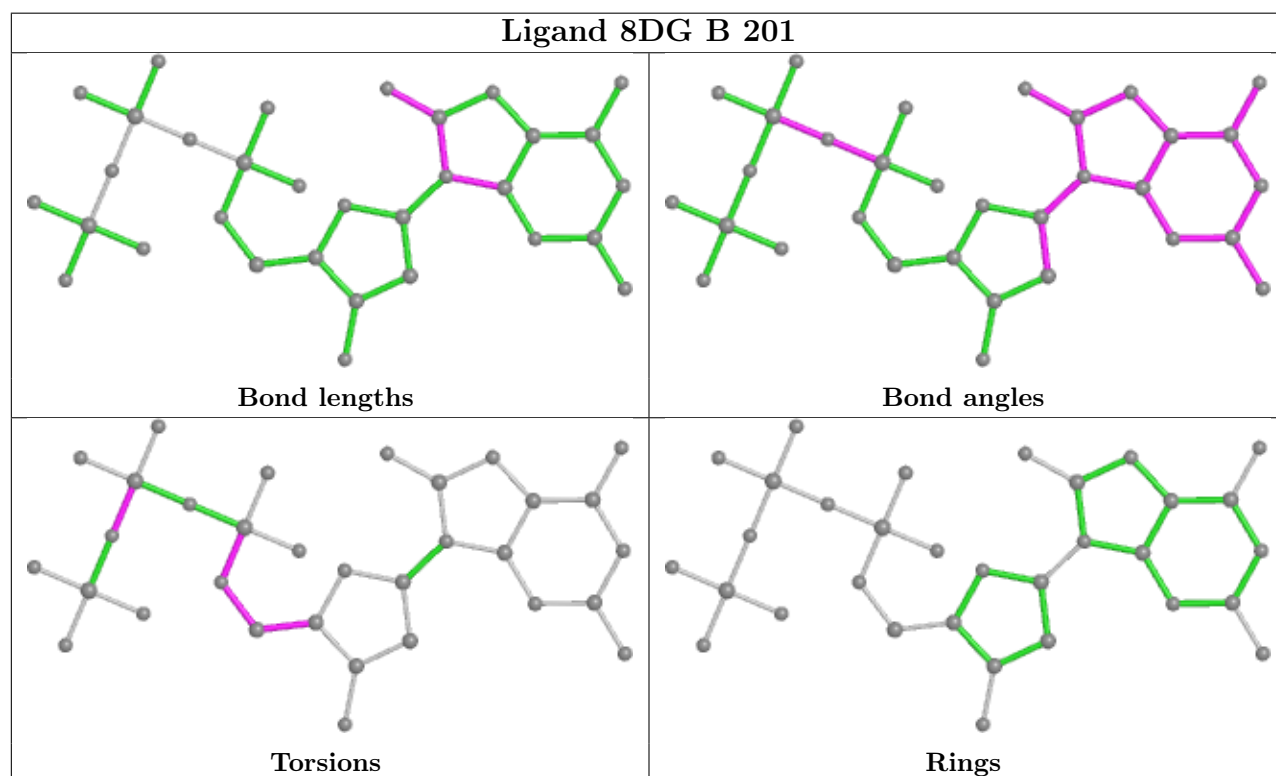
Mol	Chain	Res	Type	Atoms
2	B	201	8DG	C5'-O5'-PA-O1A
2	B	201	8DG	PG-O3B-PB-O2B
2	B	201	8DG	C4'-C5'-O5'-PA
2	B	201	8DG	PG-O3B-PB-O1B
2	B	201	8DG	C5'-O5'-PA-O3A
2	B	201	8DG	O4'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	201	8DG	2	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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

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

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

### 6.3 Carbohydrates [i](#)

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

### 6.4 Ligands [i](#)

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

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

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