

# Full wwPDB X-ray Structure Validation Report (i)

#### Jun 14, 2022 – 10:08 am BST

PDB ID	:	70V9
Title	:	Heterodimeric tRNA-Guanine Transglycosylase from mouse, apo-structure
Authors	:	Sebastiani, M.; Heine, A.; Reuter, K.
Deposited on		
Resolution	:	1.90 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

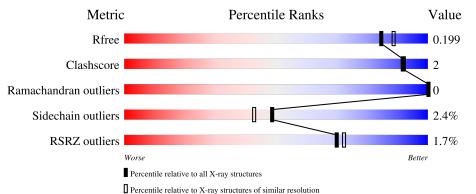
Xtriage (Phenix):1.13EDS:2.28.1buster-report:1.1.7 (2018)Percentile statistics:20191225.v01 (using entries in the PDB archive December 25th 2019)Refmac:5.8.0267CCP4:7.1.010 (Gargrove)Ideal geometry (proteins):Engh & Huber (2001)Ideal geometry (DNA, RNA):Parkinson et al. (1996)	Xtriage (Phenix) EDS buster-report Percentile statistics Refmac CCP4 Ideal geometry (proteins)	:::::::::::::::::::::::::::::::::::::::	<ul> <li>1.8.4, CSD as541be (2020)</li> <li>1.13</li> <li>2.28.1</li> <li>1.1.7 (2018)</li> <li>20191225.v01 (using entries in the PDB archive December 25th 2019)</li> <li>5.8.0267</li> <li>7.1.010 (Gargrove)</li> <li>Engh &amp; Huber (2001)</li> </ul>
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# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 1.90 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
$R_{free}$	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 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 <=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	А	419	80%	6%	13%	1
2	С	395	<sup>2%</sup> 93%		• •	-



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# 2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 6273 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 Queuine tRNA-ribosyltransferase accessory subunit 2.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	364	Total 2812	C 1787	N 471	O 526	S 28	0	10	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	cloning artifact	UNP B8ZXI1
А	1	PRO	-	cloning artifact	UNP B8ZXI1
А	416	ASP	-	expression tag	UNP B8ZXI1
А	417	ASN	-	expression tag	UNP B8ZXI1
А	418	ASN	-	expression tag	UNP B8ZXI1

• Molecule 2 is a protein called Queuine tRNA-ribosyltransferase catalytic subunit 1.

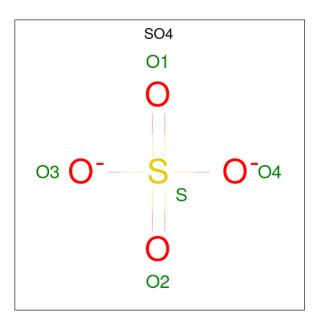
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	С	383	Total 2996	C 1885	N 530	O 553	S 28	0	14	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
С	9	GLY	-	cloning artifact	UNP Q9JMA2
С	10	PRO	-	cloning artifact	UNP Q9JMA2

• Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).





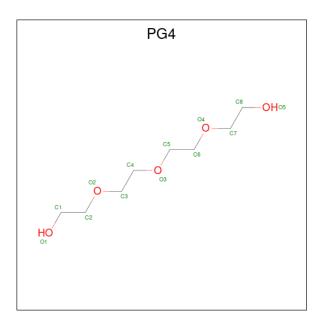
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
3	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
3	С	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
3	С	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
3	С	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
3	С	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0

• Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	Total Zn 1 1	0	0
4	С	1	Total Zn 1 1	0	0

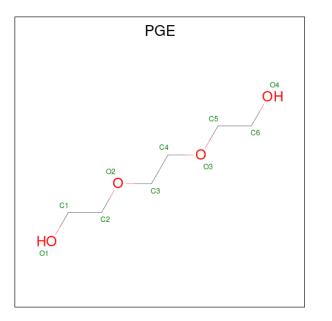
• Molecule 5 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula:  $C_8H_{18}O_5$ ).





$\mathbf{N}$	ſol	Chain	Residues	Atoms	ZeroOcc	AltConf
	5	А	1	Total         C         O           13         8         5	0	0
	5	С	1	Total         C         O           13         8         5	0	0

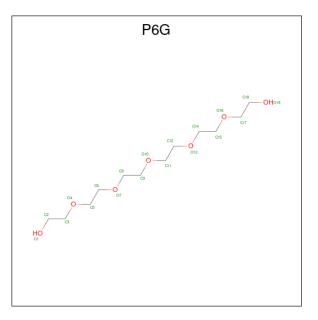
• Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula:  $C_6H_{14}O_4$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	С	1	Total         C         O           10         6         4	0	0
6	С	1	Total         C         O           10         6         4	0	0



• Molecule 7 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula:  $C_{12}H_{26}O_7$ ).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
7	С	1	Total 19	C 12	O 7	0	0

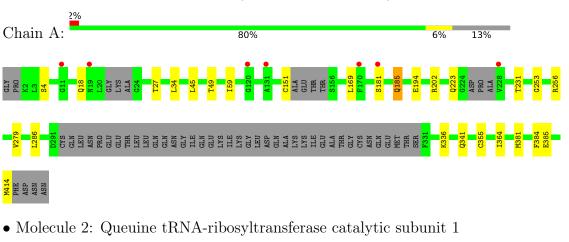
• Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	А	137	Total O 137 137	0	0
8	С	231	Total         O           231         231	0	0

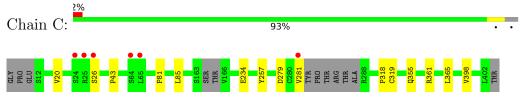


# 3 Residue-property plots (i)

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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). 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.



• Molecule 1: Queuine tRNA-ribosyltransferase accessory subunit 2





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants	100.26Å 100.26Å 202.48Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	48.66 - 1.90	Depositor
Resolution (A)	48.66 - 1.90	EDS
% Data completeness	99.9 (48.66-1.90)	Depositor
(in resolution range)	99.9 (48.66 - 1.90)	EDS
R <sub>merge</sub>	(Not available)	Depositor
R <sub>sym</sub>	0.08	Depositor
$< I/\sigma(I) > 1$	$1.67 (at 1.90 \text{\AA})$	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
D D.	0.175 , $0.199$	Depositor
$R, R_{free}$	0.176 , $0.199$	DCC
$R_{free}$ test set	4102 reflections $(5.00\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	31.5	Xtriage
Anisotropy	0.402	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	(Not available), (Not available)	EDS
L-test for twinning <sup>2</sup>	$ L  > = 0.50, < L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6273	wwPDB-VP
Average B, all atoms $(Å^2)$	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.67% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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: SO4, ZN, PG4, PGE, P6G

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	А	0.34	0/2864	0.51	0/3875	
2	С	0.39	0/3057	0.52	0/4145	
All	All	0.37	0/5921	0.52	0/8020	

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	А	2812	0	2653	11	0
2	С	2996	0	2937	6	0
3	А	10	0	0	0	0
3	С	20	0	0	0	0
4	А	1	0	0	0	0
4	С	1	0	0	0	0
5	А	13	0	18	1	0
5	С	13	0	18	0	0
6	Ċ	20	0	28	2	0
7	С	19	0	26	0	0
8	А	137	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	С	231	0	0	1	0
All	All	6273	0	5680	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 2.

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)
2:C:81:PRO:HB2	2:C:85[A]:LEU:HD23	1.70	0.74
2:C:43:PRO:HD2	2:C:361[B]:ARG:HG2	1.74	0.69
1:A:181:SER:O	1:A:185:GLN:NE2	2.27	0.68
1:A:45:LEU:HD22	1:A:49[B]:THR:HG21	1.84	0.59
1:A:34:LEU:HG	1:A:59[A]:ILE:HG23	1.88	0.56
1:A:223:GLN:HG3	1:A:253:GLY:HA3	1.89	0.53
2:C:318:PRO:HD2	2:C:355:GLN:NE2	2.24	0.53
1:A:194:GLU:O	1:A:202:ARG:HD3	2.09	0.52
1:A:336:LYS:HE3	5:A:504:PG4:H61	1.93	0.51
2:C:234:GLU:O	6:C:507:PGE:H42	2.11	0.50
1:A:45:LEU:HD22	1:A:49[B]:THR:CG2	2.47	0.44
1:A:355:CYS:SG	1:A:364[A]:ILE:HD11	2.58	0.43
2:C:26:SER:HA	2:C:398:VAL:HB	2.01	0.42
1:A:279:VAL:HG21	1:A:286:LEU:HD21	2.01	0.42
6:C:507:PGE:H62	8:C:753:HOH:O	2.18	0.42
1:A:49[B]:THR:HG23	1:A:381[B]:MET:HE3	2.01	0.42
2:C:20:VAL:HG23	2:C:365:LEU:HD21	2.03	0.41
1:A:18:GLN:HG2	1:A:27:THR:HG22	2.01	0.41

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	Perce	ntiles
1	А	364/419~(87%)	360 (99%)	4 (1%)	0	100	100
2	С	391/395~(99%)	381 (97%)	10 (3%)	0	100	100
All	All	755/814~(93%)	741 (98%)	14 (2%)	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 side chain 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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles		
1	А	289/363~(80%)	279~(96%)	10 (4%)	36 27		
2	С	320/333~(96%)	316~(99%)	4 (1%)	69 68		
All	All	609/696~(88%)	595~(98%)	14 (2%)	49 45		

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	4	SER
1	А	151	CYS
1	А	169	LEU
1	А	185	GLN
1	А	231	THR
1	А	256	ARG
1	А	341	GLN
1	А	384	PHE
1	А	385	GLU
1	А	414	MET
2	С	257	TYR
2	С	279	ASP
2	С	281	VAL
2	С	319	CYS

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



Mol	Chain	Res	Type
2	С	110	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 monosaccharides in this entry.

### 5.6 Ligand geometry (i)

Of 13 ligands modelled in this entry, 2 are monoatomic - leaving 11 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	Turne	Chain	Res	Tink	Bo	ond leng	ths	В	ond ang	les	
INIOI	Type	Unam	nes	nes	Res Link	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
7	P6G	С	509	-	18,18,18	0.53	0	17,17,17	0.46	0	
3	SO4	А	501	-	4,4,4	0.16	0	6,6,6	0.13	0	
6	PGE	С	506	-	9,9,9	0.15	0	8,8,8	0.11	0	
3	SO4	А	502	-	4,4,4	0.16	0	$6,\!6,\!6$	0.35	0	
5	PG4	A	504	-	$12,\!12,\!12$	0.14	0	11,11,11	0.14	0	
3	SO4	С	504	-	4,4,4	0.15	0	$6,\!6,\!6$	0.12	0	
6	PGE	С	507	-	$9,\!9,\!9$	0.11	0	8,8,8	0.21	0	
5	PG4	С	508	-	12,12,12	0.17	0	11,11,11	0.13	0	
3	SO4	С	503	-	4,4,4	0.14	0	$6,\!6,\!6$	0.10	0	
3	SO4	С	502	-	4,4,4	0.16	0	$6,\!6,\!6$	0.11	0	
3	SO4	С	501	-	4,4,4	0.17	0	$6,\!6,\!6$	0.24	0	

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



Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	P6G	С	509	-	-	7/16/16/16	-
6	PGE	С	506	-	-	2/7/7/7	-
5	PG4	А	504	-	-	4/10/10/10	-
6	PGE	С	507	-	-	3/7/7/7	-
5	PG4	С	508	-	-	5/10/10/10	-

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	А	504	PG4	O3-C5-C6-O4
7	С	509	P6G	O10-C11-C12-O13
7	С	509	P6G	O16-C17-C18-O19
7	С	509	P6G	O1-C2-C3-O4
7	С	509	P6G	O7-C8-C9-O10
5	А	504	PG4	C3-C4-O3-C5
5	А	504	PG4	C5-C6-O4-C7
7	С	509	P6G	C9-C8-O7-C6
5	С	508	PG4	C3-C4-O3-C5
7	С	509	P6G	C8-C9-O10-C11
5	С	508	PG4	C1-C2-O2-C3
6	С	507	PGE	C6-C5-O3-C4
5	С	508	PG4	C6-C5-O3-C4
6	С	506	PGE	O3-C5-C6-O4
5	С	508	PG4	O2-C3-C4-O3
5	А	504	PG4	C8-C7-O4-C6
6	С	506	PGE	C3-C4-O3-C5
5	С	508	PG4	C4-C3-O2-C2
6	С	507	PGE	O3-C5-C6-O4
6	С	507	PGE	C4-C3-O2-C2
7	С	509	P6G	C14-C15-O16-C17

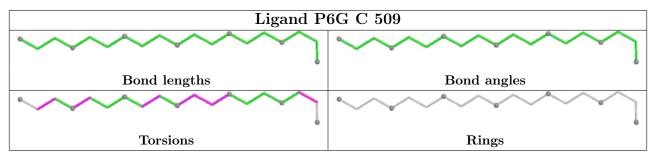
There are no ring outliers.

2 monomers are involved in 3 short contacts:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	А	504	PG4	1	0
6	С	507	PGE	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 then 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)

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,  $95^{th}$  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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	Q < 0.9
1	А	364/419~(86%)	0.04	7 (1%) 66 69	22, 41, 71, 98	0
2	С	383/395~(96%)	-0.07	6 (1%) 72 74	21, 30, 51, 82	0
All	All	747/814~(91%)	-0.02	13 (1%) 70 72	21, 34, 68, 98	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	С	281	VAL	4.5
2	С	24	SER	3.3
2	С	25	ARG	3.0
1	А	19	ASN	2.7
2	С	65	LEU	2.6
2	С	26	SER	2.6
1	А	11	GLY	2.5
2	С	64[A]	SER	2.2
1	А	120	GLY	2.2
1	А	228	VAL	2.2
1	А	170	PHE	2.1
1	А	131	ALA	2.1
1	А	181	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 monosaccharides in this entry.



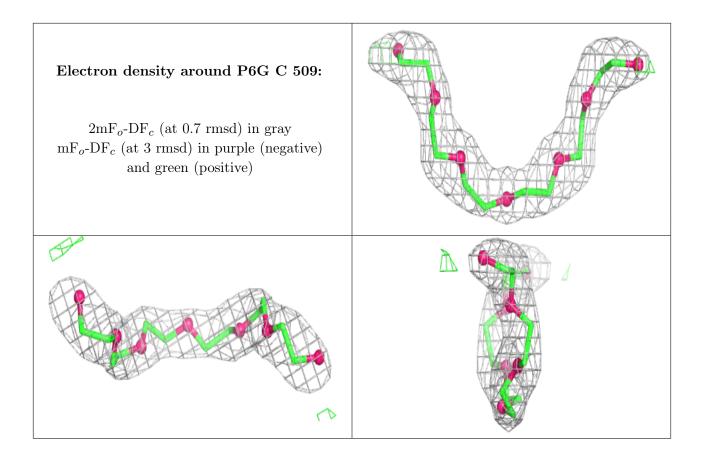
## 6.4 Ligands (i)

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. The B-factors column lists the minimum, median,  $95^{th}$  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	$B-factors(Å^2)$	Q < 0.9
5	PG4	С	508	13/13	0.82	0.18	48,58,63,64	0
6	PGE	С	507	10/10	0.89	0.14	$51,\!56,\!66,\!70$	0
3	SO4	С	502	5/5	0.92	0.12	73,83,90,96	0
5	PG4	А	504	13/13	0.92	0.14	$43,\!51,\!72,\!72$	0
7	P6G	С	509	19/19	0.92	0.13	$47,\!51,\!54,\!55$	19
3	SO4	А	502	5/5	0.93	0.13	$34,\!54,\!60,\!66$	5
6	PGE	С	506	10/10	0.93	0.19	$32,\!34,\!42,\!47$	10
3	SO4	С	504	5/5	0.94	0.10	$58,\!61,\!65,\!77$	5
3	SO4	С	503	5/5	0.96	0.16	66,69,71,72	5
3	SO4	А	501	5/5	0.98	0.08	$67,\!68,\!73,\!74$	0
3	SO4	С	501	5/5	0.98	0.13	$31,\!36,\!37,\!43$	5
4	ZN	С	505	1/1	1.00	0.10	26,26,26,26	0
4	ZN	А	503	1/1	1.00	0.06	29,29,29,29	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





## 6.5 Other polymers (i)

There are no such residues in this entry.

