

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

#### Nov 21, 2023 – 10:43 PM JST

PDB ID	:	7XNE
Title	:	Crystal structure of CBP bromodomain liganded with Y08284
Authors	:	Xiang, Q.; Wang, C.; Wu, T.; Zhang, C.; Hu, Q.; Luo, G.; Hu, J.; Zhuang, X.;
		Zou, L.; Shen, H.; Wu, X.; Zhang, Y.; Kong, X.; Xu, Y.
Deposited on	:	2022-04-28
Resolution	:	2.17  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 types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

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

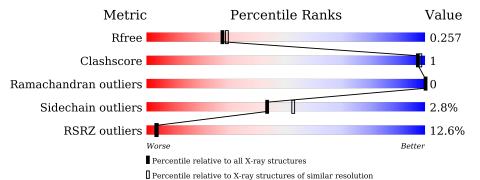
Xtriage (Phenix) EDS buster-report Percentile statistics Refmac CCP4 Ideal geometry (proteins) Ideal geometry (DNA, RNA)	:::::::::::::::::::::::::::::::::::::::	20191225.v01 (using entries in the PDB archive December 25th 2019) 5.8.0158 7.0.044 (Gargrove) Engh & Huber (2001) Parkinson et al. (1996)
Ideal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

# 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 2.17 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R <sub>free</sub>	130704	6864 (2.20-2.16)
Clashscore	141614	7689 (2.20-2.16)
Ramachandran outliers	138981	7564 (2.20-2.16)
Sidechain outliers	138945	7564 (2.20-2.16)
RSRZ outliers	127900	6738 (2.20-2.16)

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	133	9%	••	14%
1	В	133	83%	•	14%



#### 7XNE

# 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2058 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 CREB-binding protein.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	115	Total 963	C 626		0 177	${ m S}{ m 5}$	0	0	0
1	В	115	Total 963	C 626		O 177	${ m S}{ m 5}$	0	0	0

Chain	Residue	Modelled	Actual	Comment	Reference
А	1065	MET	-	expression tag	UNP Q92793
А	1066	LYS	-	expression tag	UNP Q92793
А	1067	LYS	-	expression tag	UNP Q92793
А	1068	GLY	-	expression tag	UNP Q92793
А	1069	HIS	-	expression tag	UNP Q92793
A	1070	HIS	-	expression tag	UNP Q92793
А	1071	HIS	-	expression tag	UNP Q92793
А	1072	HIS	-	expression tag	UNP Q92793
А	1073	HIS	-	expression tag	UNP Q92793
А	1074	HIS	-	expression tag	UNP Q92793
А	1075	LEU	-	expression tag	UNP Q92793
А	1076	VAL	-	expression tag	UNP Q92793
А	1077	PRO	-	expression tag	UNP Q92793
А	1078	ARG	-	expression tag	UNP Q92793
А	1079	GLY	-	expression tag	UNP Q92793
А	1080	SER	-	expression tag	UNP Q92793
В	1065	MET	-	expression tag	UNP Q92793
В	1066	LYS	-	expression tag	UNP Q92793
В	1067	LYS	-	expression tag	UNP Q92793
В	1068	GLY	-	expression tag	UNP Q92793
В	1069	HIS	-	expression tag	UNP Q92793
В	1070	HIS	-	expression tag	UNP Q92793
В	1071	HIS	-	expression tag	UNP Q92793
В	1072	HIS	-	expression tag	UNP Q92793
В	1073	HIS	-	expression tag	UNP Q92793

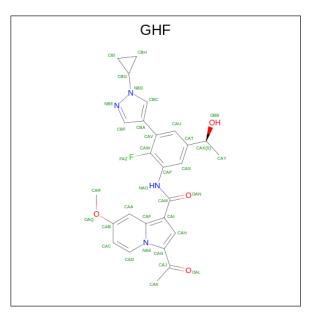
There are 32 discrepancies between the modelled and reference sequences:



Contentia	Contentaca from precioas page									
Chain	Residue	Modelled	Actual	Comment	Reference					
В	1074	HIS	-	expression tag	UNP Q92793					
В	1075	LEU	-	expression tag	UNP Q92793					
В	1076	VAL	-	expression tag	UNP Q92793					
В	1077	PRO	-	expression tag	UNP Q92793					
В	1078	ARG	-	expression tag	UNP Q92793					
В	1079	GLY	-	expression tag	UNP Q92793					
В	1080	SER	-	expression tag	UNP Q92793					

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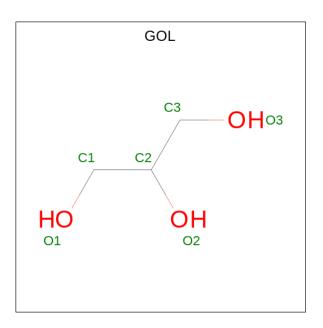
• Molecule 2 is N-[3-(1-cyclopropylpyrazol-4-yl)-2-fluoranyl-5-[(1S)-1-oxidanylethyl]phe nyl]-3-ethanoyl-7-methoxy-indolizine-1-carboxamide (three-letter code: GHF) (formula:  $C_{26}H_{25}FN_4O_4$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
9	Λ	1	Total	С	F	Ν	0	0	0	
Z	Z A	L	35	26	1	4	4	0	0	
0	D	1	Total	С	F	Ν	0	0	0	
Ζ	D		35	26	1	4	4	0	0	

• Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	В	1	Total 6	${ m C} { m 3}$	O 3	0	0

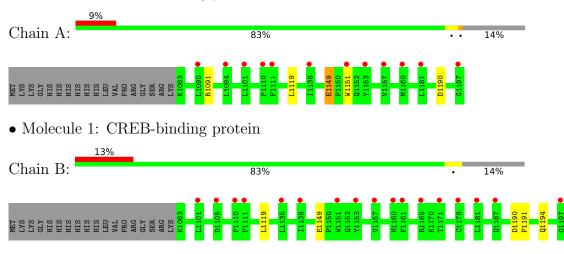
• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	27	TotalO2727	0	0
4	В	29	TotalO2929	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: CREB-binding protein



# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	47.38Å 63.27Å 97.29Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	53.04 - 2.17	Depositor
Resolution (A)	19.35 - 2.17	EDS
% Data completeness	99.7 (53.04-2.17)	Depositor
(in resolution range)	99.8(19.35-2.17)	EDS
R <sub>merge</sub>	0.06	Depositor
R <sub>sym</sub>	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.61 (at 2.17 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
D D.	0.199 , $0.259$	Depositor
$R, R_{free}$	0.210 , $0.257$	DCC
$R_{free}$ test set	785 reflections $(4.90\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	48.0	Xtriage
Anisotropy	0.336	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.36 , $46.2$	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	2058	wwPDB-VP
Average B, all atoms $(Å^2)$	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The analyses of the Patterson function reveals a significant off-origin peak that is 33.95 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.3678e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

<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: GHF, GOL

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 RMSZ	lengths	Bond angles		
			# Z  > 5	RMSZ	# Z  > 5	
1	А	0.55	0/990	0.72	0/1344	
1	В	0.55	0/990	0.70	0/1344	
All	All	0.55	0/1980	0.71	0/2688	

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	963	0	948	2	0
1	В	963	0	948	1	0
2	А	35	0	0	0	0
2	В	35	0	0	0	0
3	В	6	0	8	0	0
4	А	27	0	0	1	0
4	В	29	0	0	0	0
All	All	2058	0	1904	3	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.



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)	
1:A:1149:GLU:OE2	1:A:1151:TRP:HB2	2.09	0.52	
1:A:1091:ARG:NE	4:A:1301:HOH:O	2.42	0.52	
1:B:1190:ASP:CB	1:B:1191:PRO:CD	2.98	0.42	

All (3) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

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	А	113/133~(85%)	113 (100%)	0	0	100	100
1	В	113/133~(85%)	111 (98%)	2(2%)	0	100	100
All	All	226/266~(85%)	224 (99%)	2(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	Rotameric Outliers	
1	А	107/124~(86%)	104 (97%)	3~(3%)	43 53
1	В	107/124~(86%)	104 (97%)	3(3%)	43 53
All	All	214/248~(86%)	208~(97%)	6 (3%)	43 53



All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	1119	LEU
1	А	1149	GLU
1	А	1190	ASP
1	В	1119	LEU
1	В	1149	GLU
1	В	1194	GLN

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such side chains are listed below:

Mol	Chain	Res	Type
1	А	1092	GLN
1	А	1194	GLN
1	В	1092	GLN
1	В	1131	ASN

#### 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)

3 ligands are modelled in this entry.

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	Link	Bond lengths			Bond angles		
NIOI	ol Type Chain R	nes	LIIIK	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2	
2	GHF	А	1201	-	31,39,39	2.24	9 (29%)	33,58,58	1.49	4 (12%)
2	GHF	В	1201	-	31,39,39	2.36	12 (38%)	33,58,58	1.75	4 (12%)
3	GOL	В	1202	-	$5,\!5,\!5$	0.41	0	$5,\!5,\!5$	0.39	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 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	GHF	А	1201	-	-	1/16/28/28	0/5/5/5
2	GHF	В	1201	-	-	1/16/28/28	0/5/5/5
3	GOL	В	1202	-	-	0/4/4/4	-

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	В	1201	GHF	CAI-CAM	-6.78	1.39	1.50
2	А	1201	GHF	CAI-CAM	-6.57	1.39	1.50
2	А	1201	GHF	CAV-CBA	-6.34	1.37	1.49
2	В	1201	GHF	CAV-CBA	-6.00	1.38	1.49
2	В	1201	GHF	CAD-NAE	3.44	1.43	1.38
2	В	1201	GHF	CAT-CAX	-3.39	1.42	1.52
2	А	1201	GHF	CAD-NAE	3.10	1.43	1.38
2	В	1201	GHF	CAA-CAB	3.06	1.42	1.37
2	А	1201	GHF	CAA-CAB	2.95	1.42	1.37
2	В	1201	GHF	CAH-CAG	-2.88	1.35	1.39
2	В	1201	GHF	CAA-CAF	-2.77	1.35	1.40
2	А	1201	GHF	CAH-CAG	-2.56	1.35	1.39
2	А	1201	GHF	CAT-CAX	-2.51	1.44	1.52
2	А	1201	GHF	CAP-CAW	2.51	1.43	1.38
2	А	1201	GHF	CAP-NAO	-2.39	1.37	1.41
2	В	1201	GHF	CAV-CAW	2.29	1.42	1.39
2	В	1201	GHF	FAZ-CAW	-2.26	1.31	1.35
2	В	1201	GHF	CAH-CAI	-2.19	1.35	1.40
2	В	1201	GHF	CAP-CAW	2.18	1.42	1.38
2	А	1201	GHF	CAA-CAF	-2.13	1.36	1.40
2	В	1201	GHF	CAP-NAO	-2.13	1.37	1.41

All (21) bond length outliers are listed below:

All (8) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
2	В	1201	GHF	CBC-NBD-CBG	7.52	132.14	125.48
2	А	1201	GHF	CBC-NBD-CBG	5.15	130.05	125.48
2	А	1201	GHF	CAU-CAV-CAW	-2.81	115.31	117.26
2	В	1201	GHF	CAI-CAM-NAO	2.79	118.97	114.79
2	А	1201	GHF	CAI-CAM-NAO	2.70	118.82	114.79
2	В	1201	GHF	CAR-OAQ-CAB	2.40	122.72	117.51
2	А	1201	GHF	CAB-CAA-CAF	2.32	122.78	118.10
2	В	1201	GHF	CAU-CAV-CAW	-2.20	115.74	117.26

There are no chirality outliers.

All (2) torsion outliers are listed below:

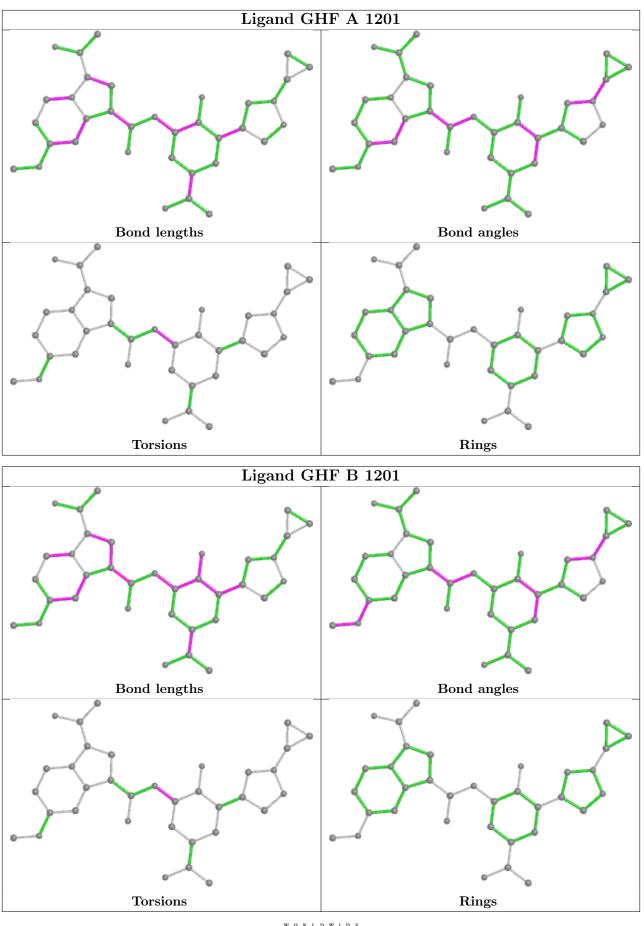
Mol	Chain	Res	Type	Atoms
2	А	1201	GHF	CAW-CAP-NAO-CAM
2	В	1201	GHF	CAW-CAP-NAO-CAM

There are no ring outliers.

No monomer is involved in short contacts.

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	< <b>RSRZ</b> >	#RSRZ>2		$OWAB(Å^2)$	Q<0.9	
1	А	115/133~(86%)	0.49	12 (10%)	6	6	36, 52, 72, 86	0
1	В	115/133~(86%)	0.71	17 (14%)	2	2	38, 57, 84, 111	0
All	All	230/266~(86%)	0.60	29~(12%)	3	4	36, 54, 80, 111	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res Type		RSRZ
1	В	1197	GLY	5.9
1	В	1111	PHE	4.1
1	А	1111	PHE	4.1
1	А	1151	TRP	4.1
1	В	1151	TRP	4.0
1	В	1157	VAL	3.9
1	В	1169	ARG	3.9
1	А	1110	PRO	3.8
1	В	1161	PHE	3.7
1	А	1157	VAL	3.6
1	А	1197	GLY	3.3
1	В	1181	LEU	3.2
1	В	1105	ASP	3.1
1	В	1178	CYS	2.9
1	А	1160	MET	2.9
1	В	1110	PRO	2.9
1	В	1138	ILE	2.9
1	А	1153	TYR	2.8
1	В	1153	TYR	2.8
1	В	1187	GLN	2.7
1	А	1101	LEU	2.4
1	А	1138	ILE	2.3
1	В	1160	MET	2.2
1	В	1101	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	В	1135	LEU	2.1
1	В	1171	THR	2.1
1	А	1094	LEU	2.1
1	А	1181	LEU	2.0
1	А	1090	LEU	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.

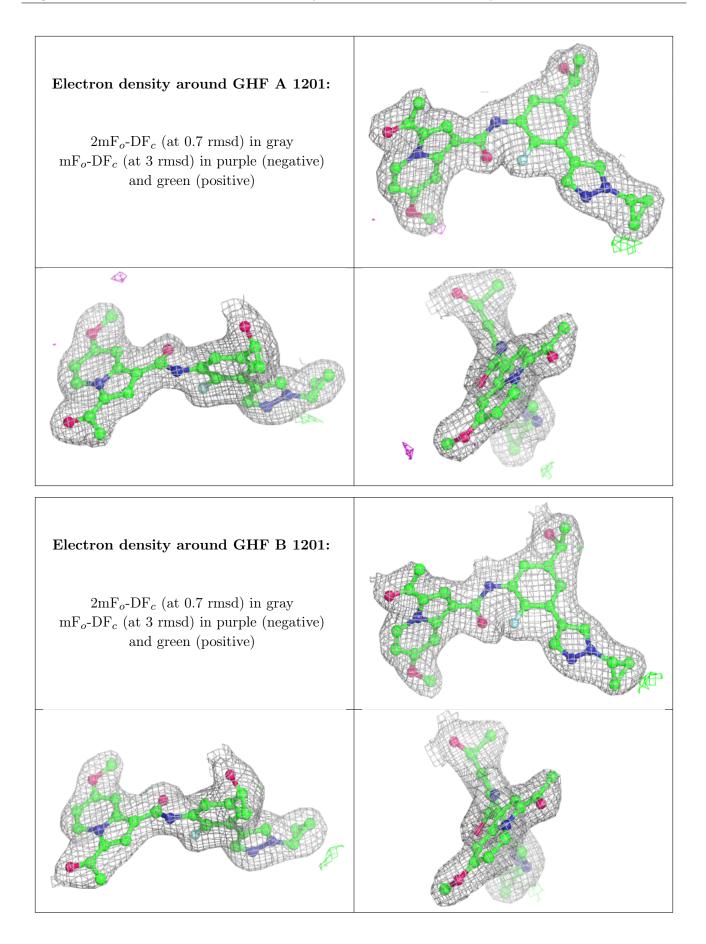
### 6.4 Ligands (i)

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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
3	GOL	В	1202	6/6	0.77	0.17	73,74,75,76	0
2	GHF	А	1201	35/35	0.94	0.14	35,42,46,48	0
2	GHF	В	1201	35/35	0.95	0.15	35,42,48,49	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.

