



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 04:12 PM UTC

PDB ID : 6D0Y / pdb_00006d0y
Title : X-ray Crystal Structure of PGC-1beta C-terminus bound to the CBP80-CBP20 Cap Binding Complex
Authors : Gleghorn, M.L.; Maquat, L.E.
Deposited on : 2018-04-11
Resolution : 2.68 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 ⓘ 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 Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

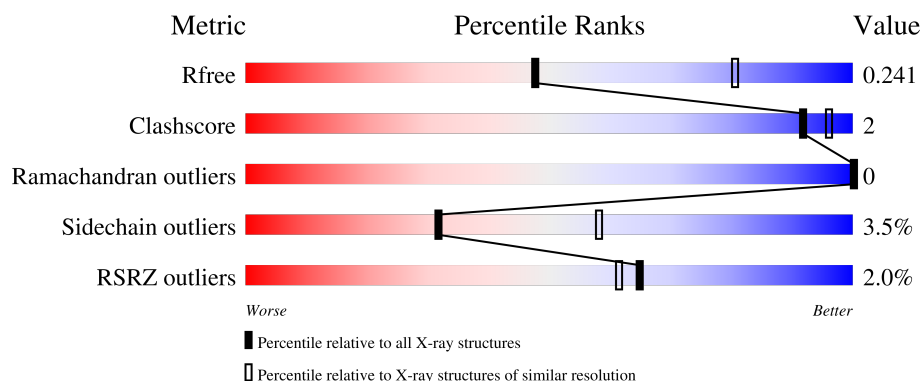
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 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}	180053	5070 (2.70-2.66)
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)

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 $\leq 5\%$. 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	C	782	0% 88% 6% 6%
2	A	176	2% 77% 6% 17%
3	B	31	19% 42% 19% 39%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14749 atoms, of which 7105 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 Nuclear cap-binding protein subunit 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	C	737	Total	C	H	N	O	S	0	0	0
			11855	3864	5865	1011	1077	38			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	23	MET	-	initiating methionine	UNP Q09161
C	479	VAL	ALA	conflict	UNP Q09161
C	791	ALA	-	expression tag	UNP Q09161
C	792	GLU	-	expression tag	UNP Q09161
C	793	ASN	-	expression tag	UNP Q09161
C	794	LEU	-	expression tag	UNP Q09161
C	795	TYR	-	expression tag	UNP Q09161
C	796	PHE	-	expression tag	UNP Q09161
C	797	GLN	-	expression tag	UNP Q09161
C	798	GLY	-	expression tag	UNP Q09161
C	799	HIS	-	expression tag	UNP Q09161
C	800	HIS	-	expression tag	UNP Q09161
C	801	HIS	-	expression tag	UNP Q09161
C	802	HIS	-	expression tag	UNP Q09161
C	803	HIS	-	expression tag	UNP Q09161
C	804	HIS	-	expression tag	UNP Q09161

- Molecule 2 is a protein called Nuclear cap-binding protein subunit 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	A	146	Total	C	H	N	O	S	0	0	0
			2318	740	1127	213	232	6			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	GLY	-	expression tag	UNP P52298
A	-18	PRO	-	expression tag	UNP P52298
A	-17	LEU	-	expression tag	UNP P52298
A	-16	HIS	-	expression tag	UNP P52298
A	-15	MET	-	expression tag	UNP P52298
A	-14	GLY	-	expression tag	UNP P52298
A	-13	GLY	-	expression tag	UNP P52298
A	-12	SER	-	expression tag	UNP P52298
A	-11	PRO	-	expression tag	UNP P52298
A	-10	GLU	-	expression tag	UNP P52298
A	-9	PHE	-	expression tag	UNP P52298
A	-8	PRO	-	expression tag	UNP P52298
A	-7	GLY	-	expression tag	UNP P52298
A	-6	ARG	-	expression tag	UNP P52298
A	-5	LEU	-	expression tag	UNP P52298
A	-4	GLU	-	expression tag	UNP P52298
A	-3	ALA	-	expression tag	UNP P52298
A	-2	ALA	-	expression tag	UNP P52298
A	-1	ASP	-	expression tag	UNP P52298
A	0	PRO	-	expression tag	UNP P52298

- Molecule 3 is a protein called Peroxisome proliferator-activated receptor gamma coactivator 1-beta.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
3	B	19	Total	C	H	I	N	O	S	0	0	0
			223	86	86	1	21	28	1			

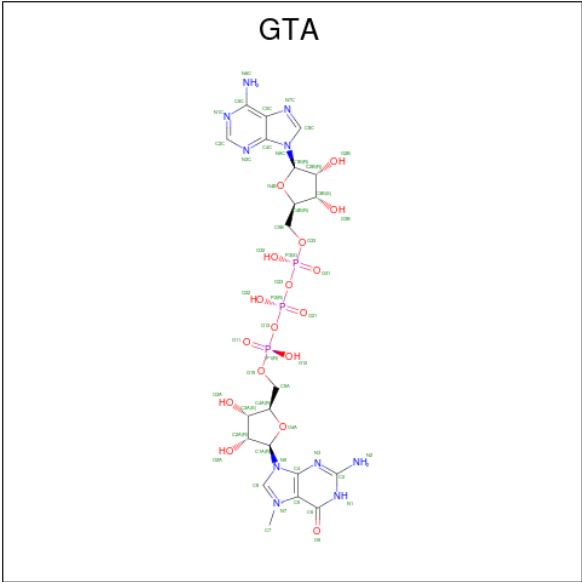
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	ACE	-	expression tag	UNP Q86YN6

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is P1-7-METHYLGUANOSINE-P3-ADENOSINE-5',5'-TRIPHOSPHATE (CCD ID: GTA) (formula: C₂₁H₃₀N₁₀O₁₇P₃).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	P	0	0
			78	21	27	10	17	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	230	Total	O	0	0
			230	230		
6	A	43	Total	O	0	0
			43	43		
6	B	1	Total	O	0	0
			1	1		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.11Å 111.90Å 124.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.05 – 2.68 39.05 – 2.68	Depositor EDS
% Data completeness (in resolution range)	97.7 (39.05-2.68) 92.6 (39.05-2.68)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.181 , 0.240 0.183 , 0.241	Depositor DCC
R_{free} test set	2000 reflections (6.50%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14749	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: GTA, MG, IYR

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	C	0.17	1/6141 (0.0%)	0.30	0/8336
2	A	0.16	0/1211	0.32	0/1616
3	B	0.24	0/123	0.29	0/163
All	All	0.17	1/7475 (0.0%)	0.31	0/10115

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	480	ASN	C-N	5.42	1.46	1.33

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	C	5990	5865	5956	16	0
2	A	1191	1127	1138	5	0
3	B	137	86	112	3	0
4	C	1	0	0	0	0
5	A	51	27	25	2	0
6	A	43	0	0	0	0
6	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	C	230	0	0	2	0
All	All	7644	7105	7231	24	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.

The worst 5 of 24 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)
5:A:201:GTA:C1A	5:A:201:GTA:O4A	1.64	1.15
1:C:403:ASP:O	1:C:741:LYS:NZ	2.24	0.68
1:C:401:ARG:NH1	6:C:1004:HOH:O	2.29	0.66
1:C:606:ASP:OD1	1:C:610:ARG:NH2	2.36	0.58
3:B:7:ASP:OD1	3:B:8:PHE:N	2.37	0.57

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	C	731/782 (94%)	723 (99%)	8 (1%)	0	100	100
2	A	144/176 (82%)	142 (99%)	2 (1%)	0	100	100
3	B	16/31 (52%)	15 (94%)	1 (6%)	0	100	100
All	All	891/989 (90%)	880 (99%)	11 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	C	665/717 (93%)	641 (96%)	24 (4%)	31	57
2	A	122/144 (85%)	119 (98%)	3 (2%)	42	69
3	B	10/24 (42%)	9 (90%)	1 (10%)	7	16
All	All	797/885 (90%)	769 (96%)	28 (4%)	32	58

5 of 28 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	425	GLN
3	B	14	GLU
1	C	717	PHE
2	A	12	ASP
1	C	627	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	439	GLN
1	C	561	HIS
3	B	16	GLN
1	C	724	HIS
1	C	777	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	IYR	B	3	3	12,13,14	1.23	1 (8%)	12,17,19	2.07	2 (16%)

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
3	IYR	B	3	3	-	0/5/6/8	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	3	IYR	CE-IE	-3.34	2.03	2.10

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	3	IYR	OF-CF-CE	5.10	125.06	119.16
3	B	3	IYR	CF-CE-IE	4.06	123.67	119.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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
5	GTA	A	201	-	56,56,56	3.13	22 (39%)	83,88,88	2.05	19 (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
5	GTA	A	201	-	-	4/32/64/64	0/6/6/6

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	201	GTA	O4A-C1A	9.66	1.64	1.42
5	A	201	GTA	O4B-C1B	8.42	1.61	1.42
5	A	201	GTA	C2B-C1B	-7.24	1.30	1.53
5	A	201	GTA	C2A-C1A	-6.56	1.32	1.53
5	A	201	GTA	O4A-C4A	-6.34	1.30	1.45

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	201	GTA	N6C-C6C-N1C	-6.36	104.21	118.38
5	A	201	GTA	N3C-C2C-N1C	-5.69	119.97	128.58
5	A	201	GTA	C5C-C4C-N3C	-5.26	119.48	126.72
5	A	201	GTA	C7-N7-C5	4.99	133.01	126.80
5	A	201	GTA	C5C-C6C-N6C	4.94	135.53	123.29

There are no chirality outliers.

All (4) torsion outliers are listed below:

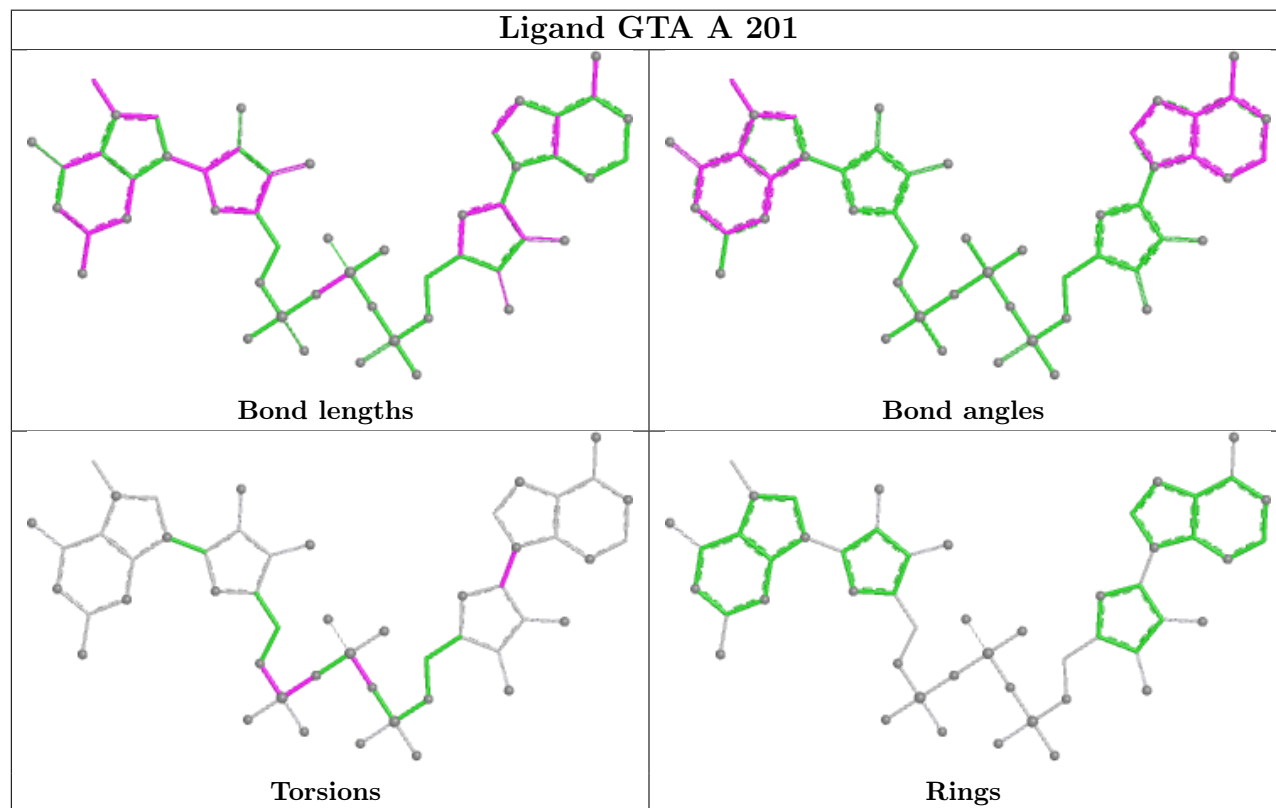
Mol	Chain	Res	Type	Atoms
5	A	201	GTA	C5A-O15-P1-O11
5	A	201	GTA	P2-O13-P1-O15
5	A	201	GTA	P3-O23-P2-O22
5	A	201	GTA	C2B-C1B-N9C-C8C

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	201	GTA	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](#)

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, 95th 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 ²)	Q<0.9
1	C	737/782 (94%)	-0.06	9 (1%) 76 74	19, 49, 95, 145	0
2	A	146/176 (82%)	0.16	3 (2%) 63 60	25, 55, 110, 176	0
3	B	18/31 (58%)	1.74	6 (33%) 1 0	85, 101, 138, 143	0
All	All	901/989 (91%)	0.01	18 (1%) 65 61	19, 50, 102, 176	0

The worst 5 of 18 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	6	LEU	6.4
3	B	1	SER	4.9
1	C	670	HIS	4.2
2	A	7	LYS	3.9
2	A	8	ALA	3.6

6.2 Non-standard residues in protein, DNA, RNA chains [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, 95th 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
3	IYR	B	3	13/14	0.58	0.18	108,131,156,228	0

6.3 Carbohydrates [i](#)

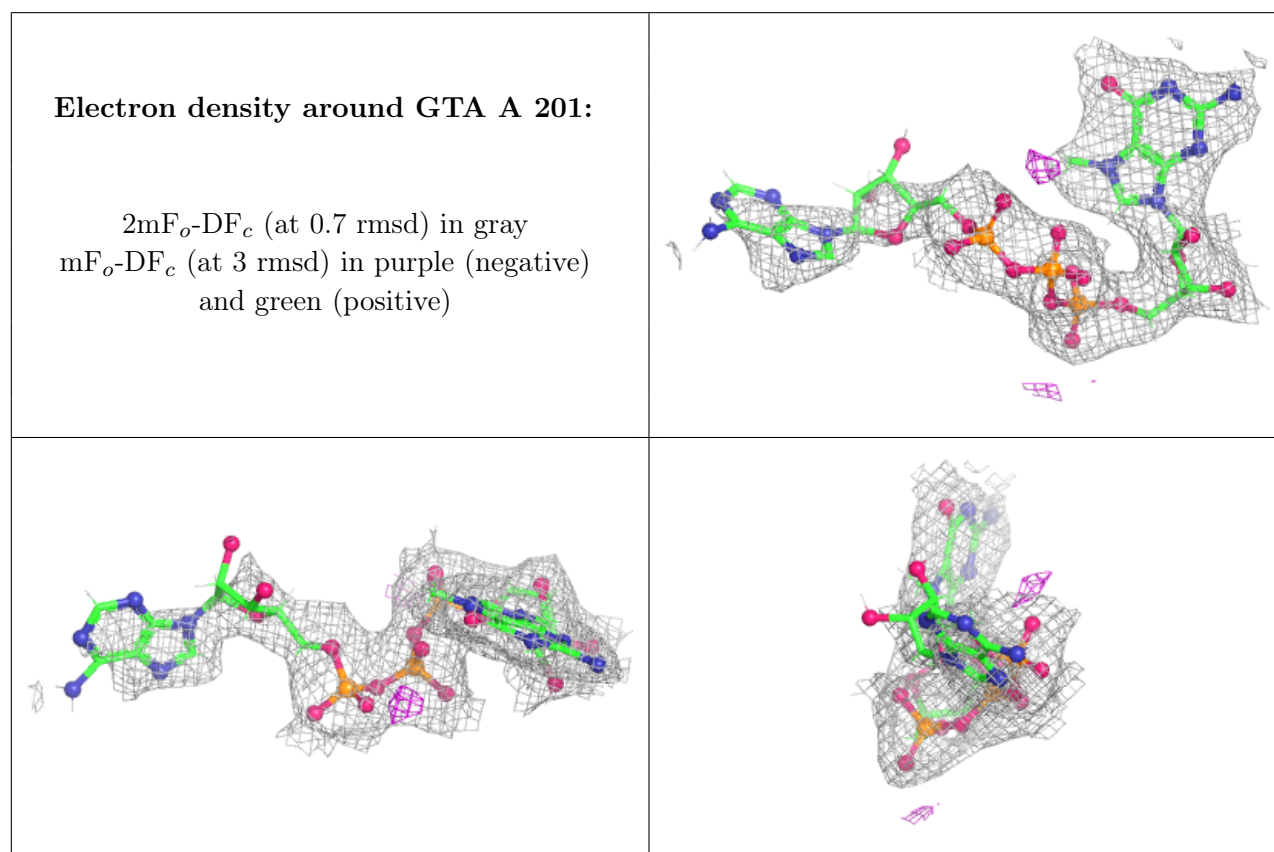
There are no oligosaccharides in this entry.

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, 95th 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	GTA	A	201	51/51	0.93	0.10	34,62,158,168	0
4	MG	C	901	1/1	0.98	0.04	32,32,32,32	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.