



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 7, 2022 – 06:06 PM EDT

PDB ID : 7MNT
Title : Crystal Structure of the ZnF5 or ZnF6 of Nucleoporin NUP358/RanBP2 in complex with Ran-GDP
Authors : Bley, C.J.; Nie, S.; Mobbs, G.W.; Petrovic, S.; Gres, A.T.; Liu, X.; Mukherjee, S.; Harvey, S.; Huber, F.M.; Lin, D.H.; Brown, B.; Tang, A.W.; Rundlet, E.J.; Correia, A.R.; Chen, S.; Regmi, S.G.; Stevens, T.A.; Jette, C.A.; Dasso, M.; Patke, A.; Palazzo, A.F.; Kossiakoff, A.A.; Hoelz, A.
Deposited on : 2021-05-01
Resolution : 2.45 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.28.1
buster-report : 1.1.7 (2018)
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) : 2.28.1

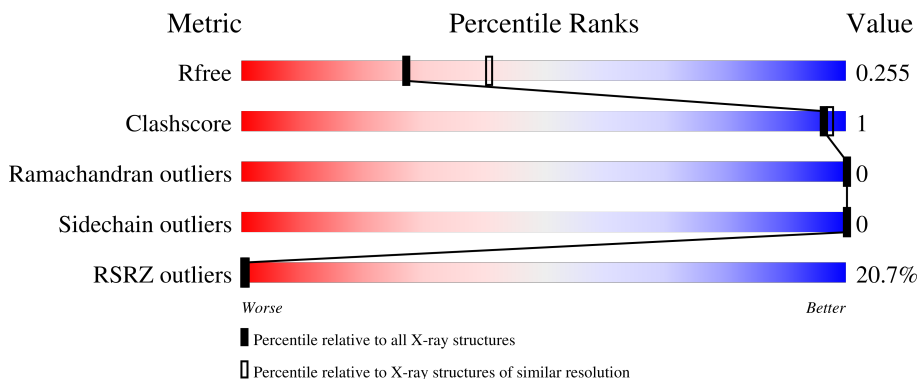
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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	A	236	
1	C	236	
2	B	43	
2	D	43	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7708 atoms, of which 3826 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 GTP-binding nuclear protein Ran.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	201	3230	1037	1623	279	285	6	0	0	0
1	C	202	3245	1042	1631	280	286	6	0	0	0

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P62826
A	-18	GLY	-	expression tag	UNP P62826
A	-17	SER	-	expression tag	UNP P62826
A	-16	SER	-	expression tag	UNP P62826
A	-15	HIS	-	expression tag	UNP P62826
A	-14	HIS	-	expression tag	UNP P62826
A	-13	HIS	-	expression tag	UNP P62826
A	-12	HIS	-	expression tag	UNP P62826
A	-11	HIS	-	expression tag	UNP P62826
A	-10	HIS	-	expression tag	UNP P62826
A	-9	SER	-	expression tag	UNP P62826
A	-8	SER	-	expression tag	UNP P62826
A	-7	GLY	-	expression tag	UNP P62826
A	-6	LEU	-	expression tag	UNP P62826
A	-5	VAL	-	expression tag	UNP P62826
A	-4	PRO	-	expression tag	UNP P62826
A	-3	ARG	-	expression tag	UNP P62826
A	-2	GLY	-	expression tag	UNP P62826
A	-1	SER	-	expression tag	UNP P62826
A	0	HIS	-	expression tag	UNP P62826
A	35	SER	PHE	engineered mutation	UNP P62826
C	-19	MET	-	expression tag	UNP P62826
C	-18	GLY	-	expression tag	UNP P62826
C	-17	SER	-	expression tag	UNP P62826
C	-16	SER	-	expression tag	UNP P62826

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-15	HIS	-	expression tag	UNP P62826
C	-14	HIS	-	expression tag	UNP P62826
C	-13	HIS	-	expression tag	UNP P62826
C	-12	HIS	-	expression tag	UNP P62826
C	-11	HIS	-	expression tag	UNP P62826
C	-10	HIS	-	expression tag	UNP P62826
C	-9	SER	-	expression tag	UNP P62826
C	-8	SER	-	expression tag	UNP P62826
C	-7	GLY	-	expression tag	UNP P62826
C	-6	LEU	-	expression tag	UNP P62826
C	-5	VAL	-	expression tag	UNP P62826
C	-4	PRO	-	expression tag	UNP P62826
C	-3	ARG	-	expression tag	UNP P62826
C	-2	GLY	-	expression tag	UNP P62826
C	-1	SER	-	expression tag	UNP P62826
C	0	HIS	-	expression tag	UNP P62826
C	35	SER	PHE	engineered mutation	UNP P62826

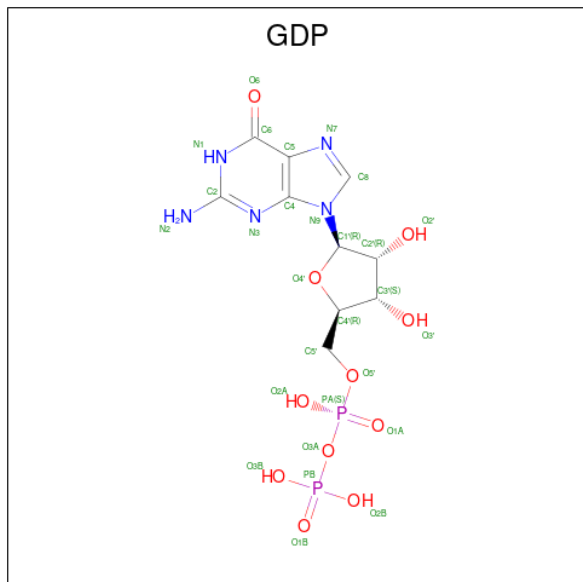
- Molecule 2 is a protein called E3 SUMO-protein ligase RanBP2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
2	B	38	561	176	274	50	55	6	0	0	0
2	D	38	561	176	274	50	55	6	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1592	GLY	-	expression tag	UNP P49792
B	1593	PRO	-	expression tag	UNP P49792
B	1594	LEU	-	expression tag	UNP P49792
B	1595	GLY	-	expression tag	UNP P49792
B	1596	SER	-	expression tag	UNP P49792
B	1597	MET	-	expression tag	UNP P49792
D	1592	GLY	-	expression tag	UNP P49792
D	1593	PRO	-	expression tag	UNP P49792
D	1594	LEU	-	expression tag	UNP P49792
D	1595	GLY	-	expression tag	UNP P49792
D	1596	SER	-	expression tag	UNP P49792
D	1597	MET	-	expression tag	UNP P49792

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			P
3	A	1	Total	C	H	N	O	P	0	0
			40	10	12	5	11	2		
3	C	1	Total	C	H	N	O	P	0	0
			40	10	12	5	11	2		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
5	B	1	Total	Zn	0	0
			1	1		
5	D	1	Total	Zn	0	0
			1	1		

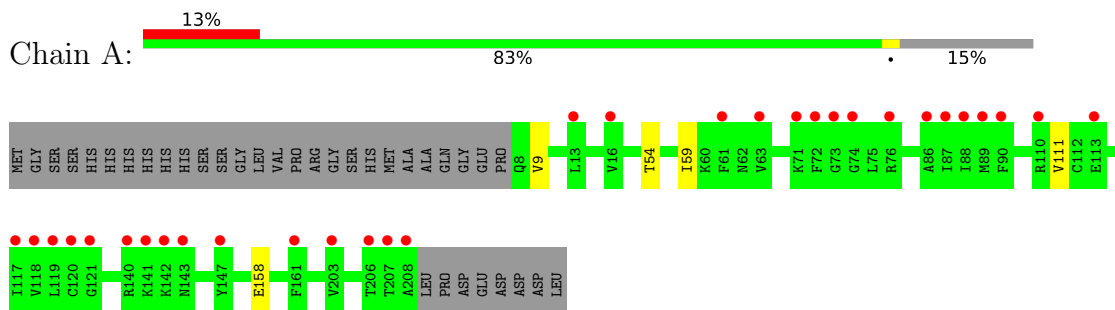
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	15	Total O 15 15	0	0
6	B	5	Total O 5 5	0	0
6	C	6	Total O 6 6	0	0
6	D	1	Total O 1 1	0	0

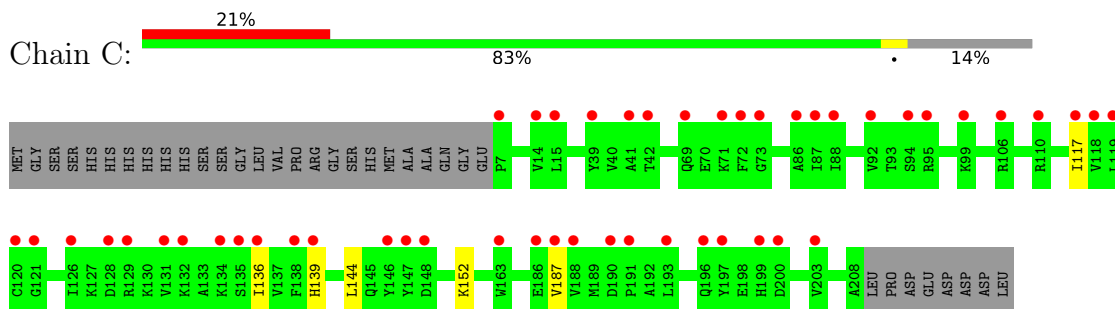
3 Residue-property plots [i](#)

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.

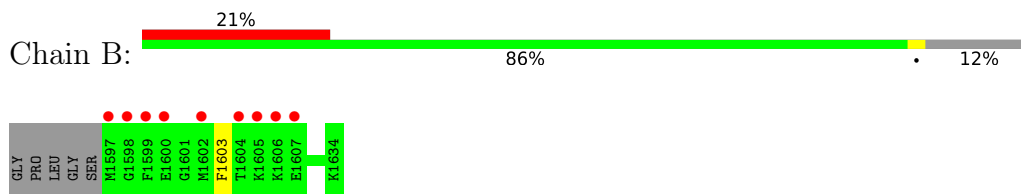
- Molecule 1: GTP-binding nuclear protein Ran



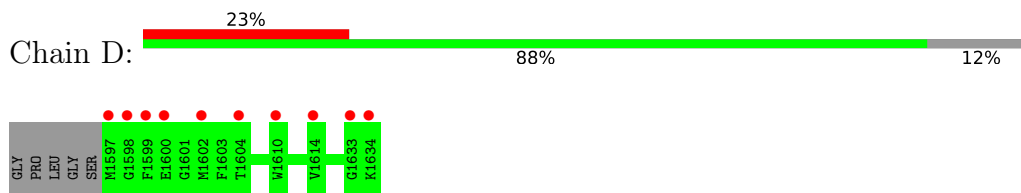
- Molecule 1: GTP-binding nuclear protein Ran



- Molecule 2: E3 SUMO-protein ligase RanBP2



- Molecule 2: E3 SUMO-protein ligase RanBP2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.02Å 94.27Å 108.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.52 – 2.45 29.52 – 2.45	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.52-2.45) 100.0 (29.52-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 2.45Å)	Xtrriage
Refinement program	PHENIX 1.18.2	Depositor
R, R_{free}	0.212 , 0.257 0.212 , 0.255	Depositor DCC
R_{free} test set	1022 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	72.2	Xtrriage
Anisotropy	0.213	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 59.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7708	wwPDB-VP
Average B, all atoms (Å ²)	108.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.68% 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: MG, ZN, GDP

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	0.26	0/1646	0.44	0/2231
1	C	0.26	0/1654	0.44	0/2242
2	B	0.27	0/291	0.43	0/386
2	D	0.25	0/291	0.43	0/386
All	All	0.26	0/3882	0.44	0/5245

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	A	1607	1623	1623	4	0
1	C	1614	1631	1631	4	0
2	B	287	274	274	1	0
2	D	287	274	274	0	0
3	A	28	12	10	0	0
3	C	28	12	11	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	1	0	0	0	0
6	A	15	0	0	0	0
6	B	5	0	0	0	0
6	C	6	0	0	0	0
6	D	1	0	0	0	0
All	All	3882	3826	3823	8	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 (8) 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:C:152:LYS:HB3	1:C:187:VAL:HG22	1.89	0.54
1:C:117:ILE:HB	1:C:144:LEU:HD22	1.91	0.51
1:A:54:THR:HG21	1:A:59:ILE:HD12	1.93	0.49
1:A:9:VAL:HG12	2:B:1603:PHE:HB2	1.94	0.49
1:A:158:GLU:OE1	1:A:158:GLU:N	2.39	0.48
1:C:136:ILE:HG22	1:C:139:HIS:CE1	2.49	0.48
1:C:187:VAL:HG23	1:C:187:VAL:O	2.13	0.47
1:A:111:VAL:O	1:A:111:VAL:HG12	2.18	0.43

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	199/236 (84%)	196 (98%)	3 (2%)	0	100	100
1	C	200/236 (85%)	196 (98%)	4 (2%)	0	100	100
2	B	36/43 (84%)	35 (97%)	1 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	36/43 (84%)	36 (100%)	0	0	100	100
All	All	471/558 (84%)	463 (98%)	8 (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 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	173/202 (86%)	173 (100%)	0	100	100
1	C	174/202 (86%)	174 (100%)	0	100	100
2	B	31/34 (91%)	31 (100%)	0	100	100
2	D	31/34 (91%)	31 (100%)	0	100	100
All	All	409/472 (87%)	409 (100%)	0	100	100

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

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 monosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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
3	GDP	C	301	4	24,30,30	5.71	14 (58%)	31,47,47	2.50	10 (32%)
3	GDP	A	301	4	24,30,30	5.67	13 (54%)	31,47,47	2.22	7 (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
3	GDP	C	301	4	-	0/12/32/32	0/3/3/3
3	GDP	A	301	4	-	1/12/32/32	0/3/3/3

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	301	GDP	C2'-C1'	-17.12	1.27	1.53
3	C	301	GDP	C2'-C1'	-17.11	1.27	1.53
3	C	301	GDP	C3'-C4'	-10.28	1.26	1.53
3	A	301	GDP	C3'-C4'	-10.27	1.26	1.53
3	A	301	GDP	O4'-C1'	9.97	1.55	1.41
3	C	301	GDP	O4'-C1'	9.53	1.54	1.41
3	C	301	GDP	C4-N3	7.28	1.47	1.35
3	A	301	GDP	C4-N3	7.27	1.47	1.35
3	C	301	GDP	C5-C6	7.16	1.53	1.41
3	A	301	GDP	C5-C6	6.76	1.53	1.41
3	C	301	GDP	C2-N2	6.32	1.46	1.33
3	A	301	GDP	C2-N2	6.13	1.46	1.33
3	C	301	GDP	C6-N1	6.01	1.43	1.33
3	A	301	GDP	C6-N1	5.88	1.43	1.33
3	C	301	GDP	O4'-C4'	5.30	1.56	1.45
3	A	301	GDP	O4'-C4'	5.07	1.56	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	301	GDP	C2-N1	4.97	1.44	1.35
3	C	301	GDP	C2'-C3'	4.97	1.66	1.53
3	A	301	GDP	C2'-C3'	4.77	1.66	1.53
3	A	301	GDP	C2-N1	4.75	1.43	1.35
3	C	301	GDP	O3'-C3'	2.67	1.49	1.43
3	A	301	GDP	O3'-C3'	2.63	1.49	1.43
3	C	301	GDP	C2-N3	2.41	1.46	1.34
3	C	301	GDP	C5'-C4'	2.39	1.59	1.51
3	A	301	GDP	C2-N3	2.35	1.45	1.34
3	A	301	GDP	C5'-C4'	2.27	1.58	1.51
3	C	301	GDP	PA-O5'	2.03	1.67	1.59

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	301	GDP	C1'-N9-C4	8.80	142.11	126.64
3	A	301	GDP	C1'-N9-C4	8.03	140.75	126.64
3	C	301	GDP	N3-C2-N1	-5.12	120.40	127.22
3	A	301	GDP	N3-C2-N1	-4.91	120.67	127.22
3	C	301	GDP	C2-N3-C4	3.68	119.56	115.36
3	A	301	GDP	PA-O3A-PB	-3.61	120.44	132.83
3	C	301	GDP	N2-C2-N1	3.61	122.87	117.25
3	C	301	GDP	C3'-C2'-C1'	3.52	106.27	100.98
3	C	301	GDP	PA-O3A-PB	-3.48	120.87	132.83
3	A	301	GDP	C2-N3-C4	3.29	119.12	115.36
3	A	301	GDP	N2-C2-N1	3.09	122.05	117.25
3	A	301	GDP	C5-C6-N1	-2.78	119.63	123.43
3	C	301	GDP	C5-C6-N1	-2.78	119.63	123.43
3	C	301	GDP	C2'-C3'-C4'	2.77	108.03	102.64
3	C	301	GDP	C2-N1-C6	2.22	119.45	115.93
3	A	301	GDP	C2-N1-C6	2.18	119.40	115.93
3	C	301	GDP	C4-C5-N7	-2.09	107.22	109.40

There are no chirality outliers.

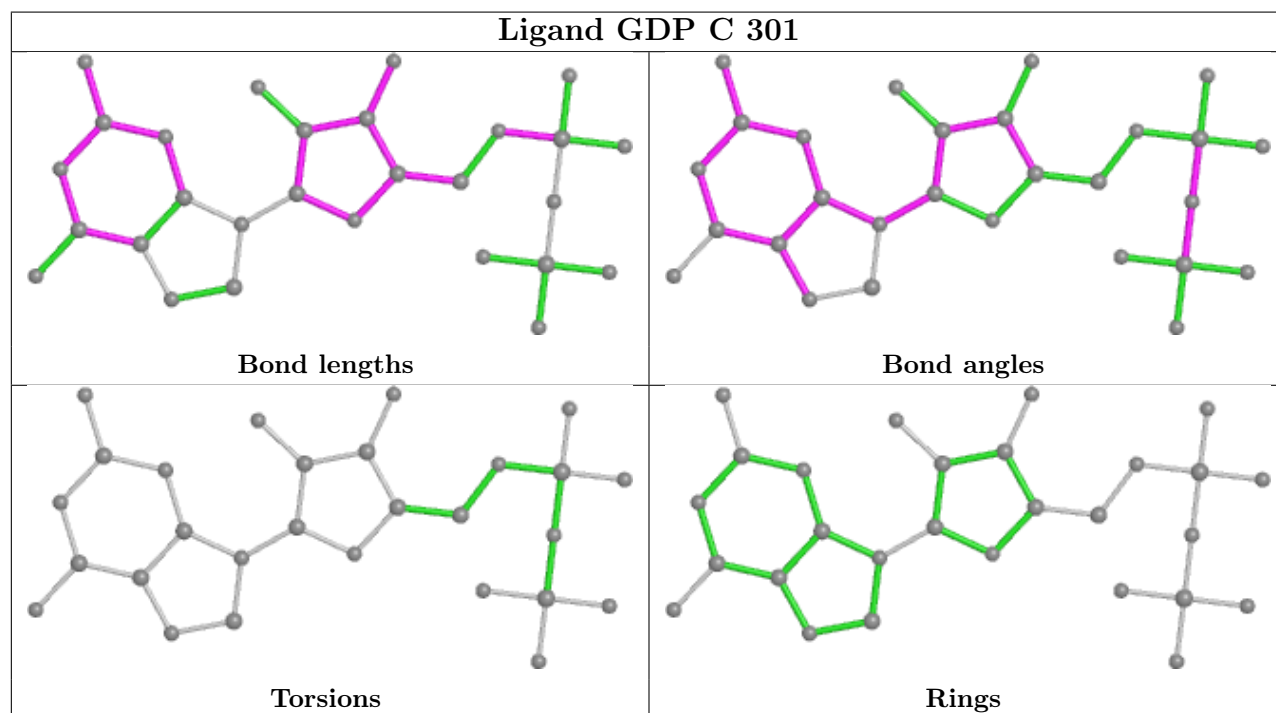
All (1) torsion outliers are listed below:

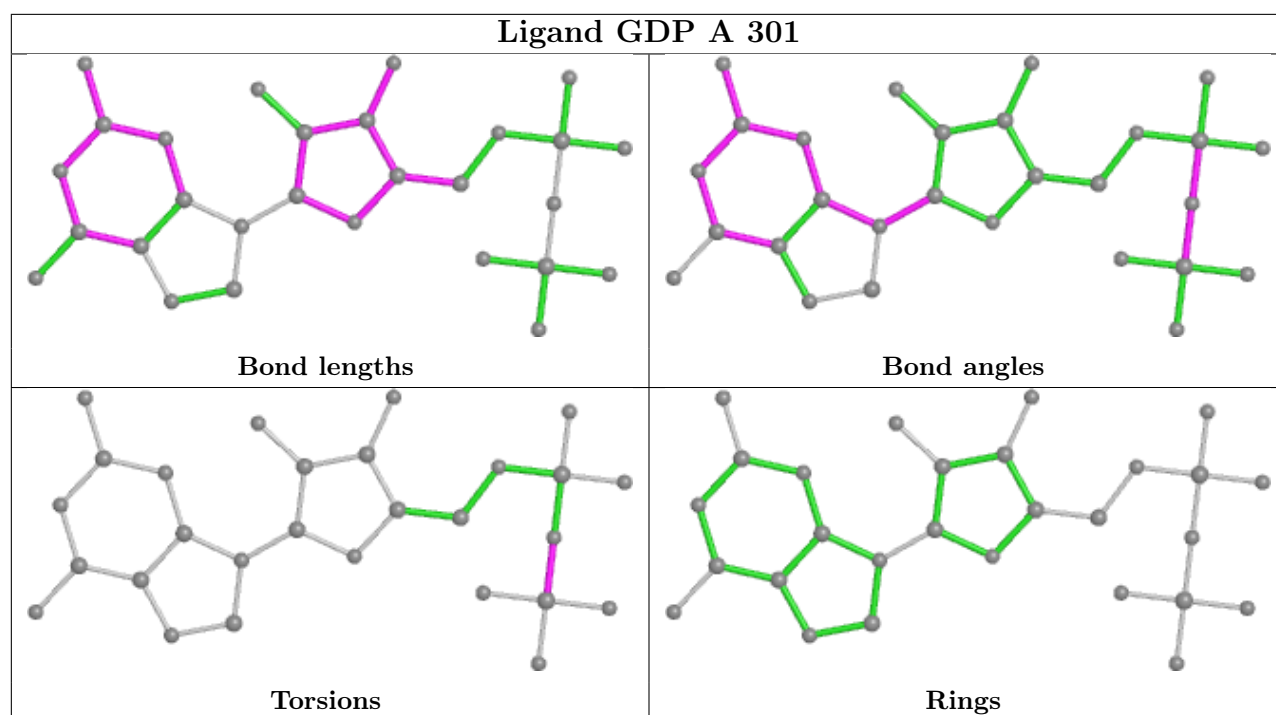
Mol	Chain	Res	Type	Atoms
3	A	301	GDP	PA-O3A-PB-O1B

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 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	A	201/236 (85%)	0.88	31 (15%) 2 1	48, 72, 141, 195	0
1	C	202/236 (85%)	1.31	49 (24%) 0 0	63, 105, 175, 213	0
2	B	38/43 (88%)	1.08	9 (23%) 0 0	58, 78, 149, 174	0
2	D	38/43 (88%)	1.38	10 (26%) 0 0	70, 96, 166, 185	0
All	All	479/558 (85%)	1.11	99 (20%) 1 0	48, 89, 170, 213	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	73	GLY	10.6
2	B	1604	THR	7.7
1	A	72	PHE	7.7
1	C	71	LYS	7.6
2	D	1597	MET	7.1
1	C	203	VAL	7.1
1	C	139	HIS	7.0
1	C	73	GLY	6.8
1	C	72	PHE	6.2
2	D	1602	MET	6.2
1	C	131	VAL	6.1
1	A	143	ASN	5.8
1	C	126	ILE	5.6
1	A	141	LYS	5.3
2	B	1597	MET	5.2
2	D	1634	LYS	4.9
2	B	1606	LYS	4.9
1	C	7	PRO	4.6
2	D	1600	GLU	4.6
1	C	95	ARG	4.5
1	C	193	LEU	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	113	GLU	4.4
1	A	120	CYS	4.4
2	B	1600	GLU	4.3
2	D	1599	PHE	4.2
1	C	199	HIS	3.6
1	C	99	LYS	3.6
1	A	87	ILE	3.5
2	B	1607	GLU	3.5
1	A	140	ARG	3.4
1	C	191	PRO	3.3
1	C	39	TYR	3.3
1	C	134	LYS	3.2
1	C	188	VAL	3.2
1	A	206	THR	3.2
1	C	41	ALA	3.2
1	C	136	ILE	3.2
1	C	42	THR	3.2
1	C	197	TYR	3.1
1	C	110	ARG	3.1
1	C	87	ILE	3.1
2	B	1602	MET	3.1
1	C	119	LEU	3.1
1	C	117	ILE	3.1
1	C	88	ILE	3.0
1	C	187	VAL	3.0
2	D	1614	VAL	3.0
1	A	119	LEU	2.9
2	B	1605	LYS	2.9
1	A	207	THR	2.9
2	D	1598	GLY	2.9
1	C	92	VAL	2.9
1	C	128	ASP	2.9
1	C	132	LYS	2.9
1	A	89	MET	2.8
1	C	138	PHE	2.8
1	A	88	ILE	2.8
1	A	86	ALA	2.8
1	A	147	TYR	2.7
1	C	120	CYS	2.7
1	C	135	SER	2.7
1	C	86	ALA	2.6
1	A	71	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	118	VAL	2.6
2	D	1604	THR	2.6
2	B	1598	GLY	2.6
1	A	110	ARG	2.5
1	A	203	VAL	2.5
1	A	142	LYS	2.5
1	C	129	ARG	2.4
1	A	208	ALA	2.4
1	A	63	VAL	2.4
1	A	118	VAL	2.4
1	C	190	ASP	2.4
1	C	200	ASP	2.4
1	A	161	PHE	2.4
1	C	106	ARG	2.4
1	C	196	GLN	2.4
2	D	1610	TRP	2.4
1	A	76	ARG	2.3
1	A	90	PHE	2.3
1	C	148	ASP	2.3
1	C	94	SER	2.3
1	A	74	GLY	2.3
1	C	69	GLN	2.3
1	C	14	VAL	2.3
1	C	186	GLU	2.3
2	D	1633	GLY	2.2
1	A	61	PHE	2.2
1	A	117	ILE	2.2
1	A	16	VAL	2.1
2	B	1599	PHE	2.1
1	A	13	LEU	2.1
1	A	121	GLY	2.1
1	C	146	TYR	2.1
1	C	15	LEU	2.1
1	C	147	TYR	2.1
1	C	121	GLY	2.1
1	C	163	TRP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates [i](#)

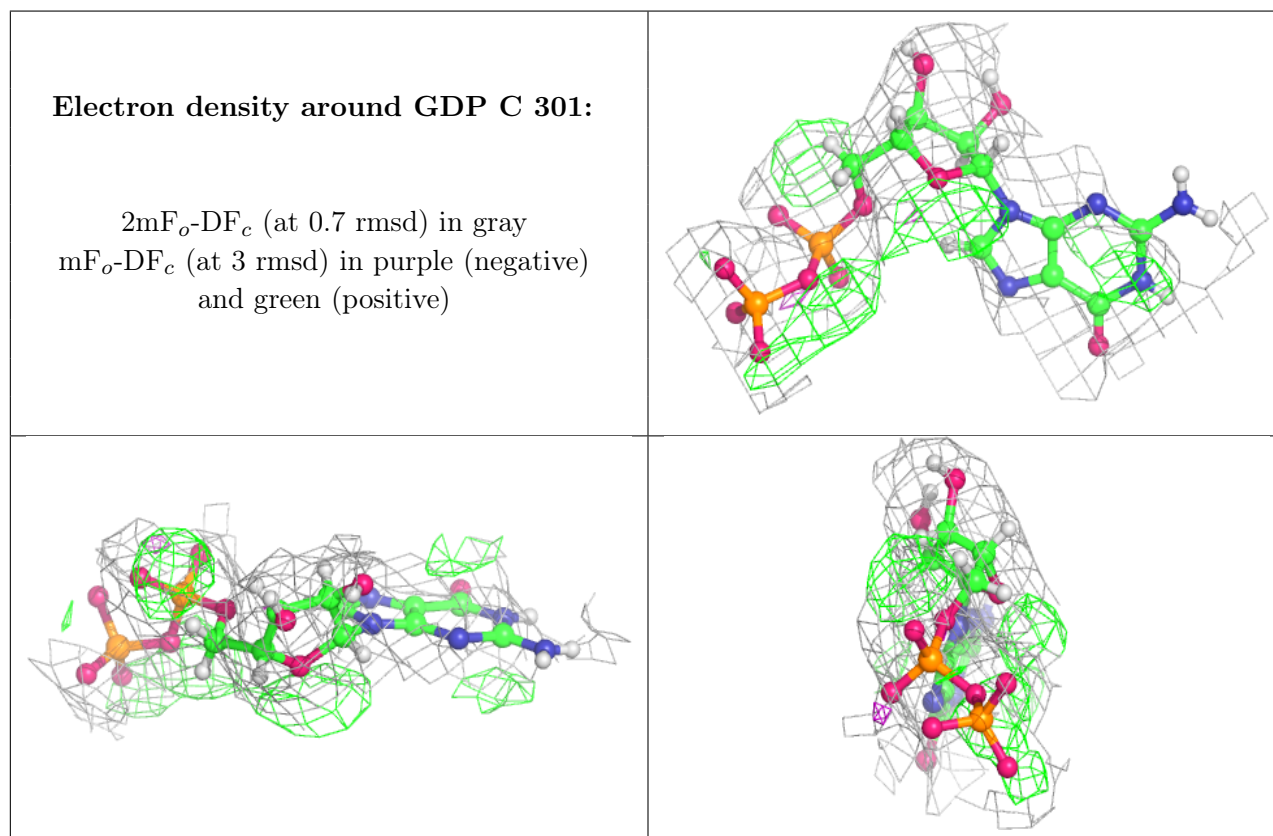
There are no monosaccharides in this entry.

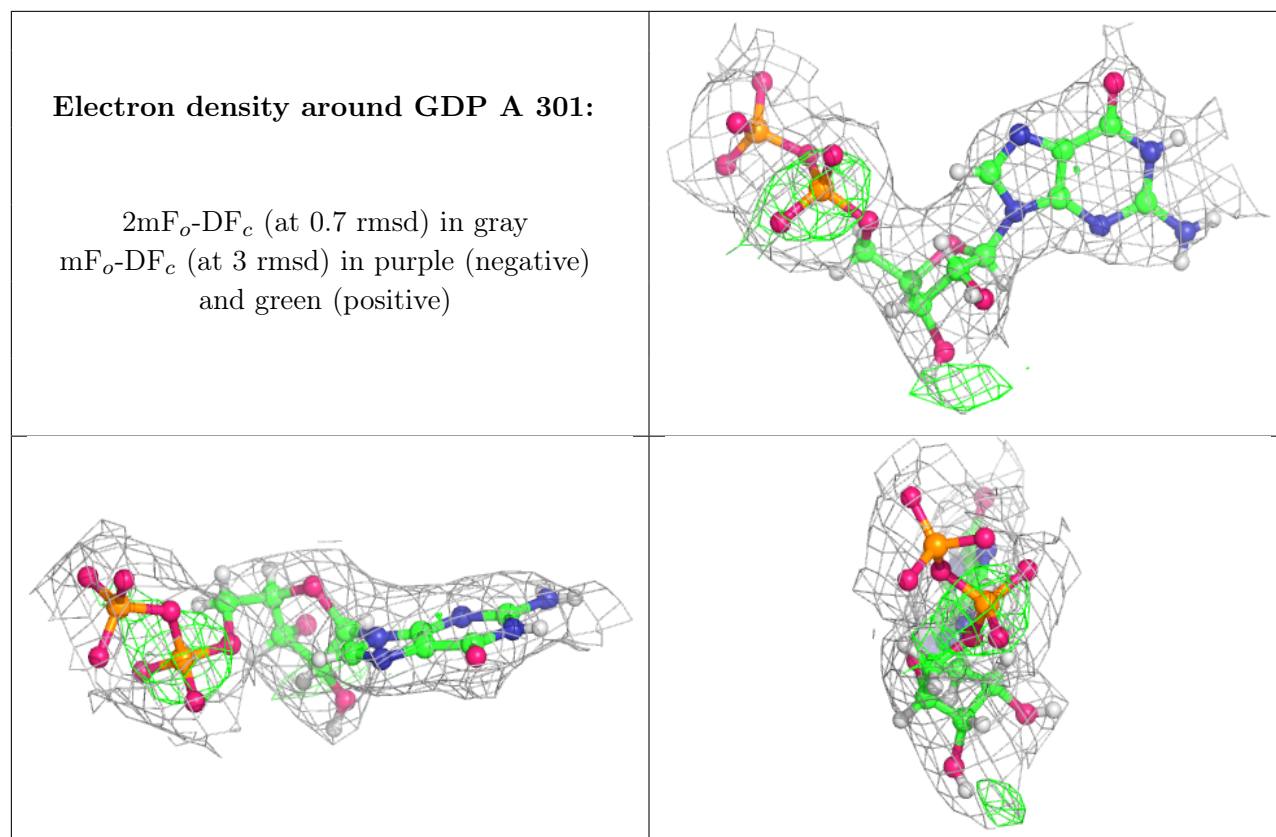
6.4 Ligands [i](#)

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, 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(\AA^2)	Q<0.9
4	MG	A	302	1/1	0.75	0.22	75,75,75,75	0
3	GDP	C	301	28/28	0.88	0.26	64,113,159,161	0
4	MG	C	302	1/1	0.92	0.19	102,102,102,102	0
3	GDP	A	301	28/28	0.95	0.22	50,66,85,89	0
5	ZN	B	1701	1/1	0.97	0.04	61,61,61,61	0
5	ZN	D	1701	1/1	0.99	0.05	88,88,88,88	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.