



wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 15, 2020 – 10:10 am BST

PDB ID : 5HCI
Title : GPN-loop GTPase Npa3 in complex with GDP
Authors : Niesser, J.; Wagner, F.R.; Kostrewa, D.; Muehlbacher, W.; Cramer, P.
Deposited on : 2016-01-04
Resolution : 2.30 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 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	:	2.11
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)	:	2.11

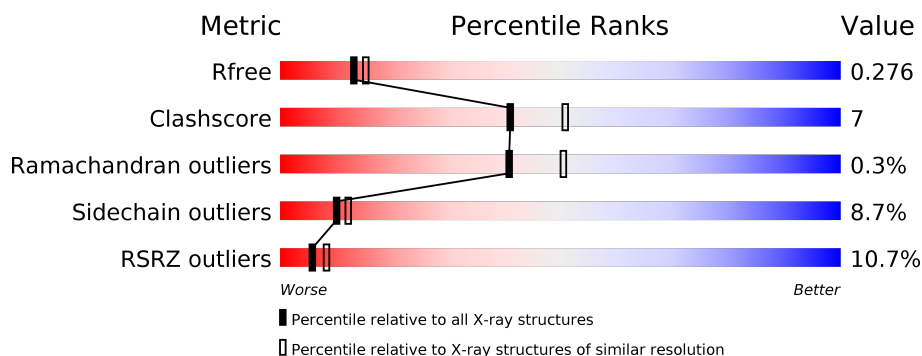
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 on 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	261	<div> <div>4%</div> <div>81% 15% ..</div> </div>
1	B	261	<div> <div>14%</div> <div>75% 17% 5% ..</div> </div>
1	C	261	<div> <div>11%</div> <div>72% 19% . 6%</div> </div>
1	D	261	<div> <div>8%</div> <div>84% 12% ...</div> </div>
1	E	261	<div> <div>7%</div> <div>81% 15% ..</div> </div>
1	F	261	<div> <div>18%</div> <div>82% 12% .. 5%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GOL	A	303	-	-	-	X
4	GOL	E	303	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12383 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 GPN-loop GTPase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	257	Total	C	N	O	S	0	0	0
			2032	1300	324	391	17			
1	B	258	Total	C	N	O	S	0	0	0
			2042	1306	327	392	17			
1	C	246	Total	C	N	O	S	0	0	0
			1937	1238	309	374	16			
1	D	258	Total	C	N	O	S	0	0	0
			2042	1306	327	392	17			
1	E	259	Total	C	N	O	S	0	0	0
			2052	1312	330	393	17			
1	F	249	Total	C	N	O	S	0	0	0
			1959	1252	314	377	16			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	LEU	deletion	UNP P47122
A	?	-	ASN	deletion	UNP P47122
A	?	-	GLY	deletion	UNP P47122
A	?	-	ASP	deletion	UNP P47122
A	?	-	ASN	deletion	UNP P47122
A	?	-	GLY	deletion	UNP P47122
A	?	-	LEU	deletion	UNP P47122
A	?	-	GLY	deletion	UNP P47122
A	?	-	SER	deletion	UNP P47122
A	265	LYS	-	expression tag	UNP P47122
A	266	HIS	-	expression tag	UNP P47122
A	267	HIS	-	expression tag	UNP P47122
A	268	HIS	-	expression tag	UNP P47122
A	269	HIS	-	expression tag	UNP P47122
A	270	HIS	-	expression tag	UNP P47122
A	271	HIS	-	expression tag	UNP P47122
B	?	-	LEU	deletion	UNP P47122

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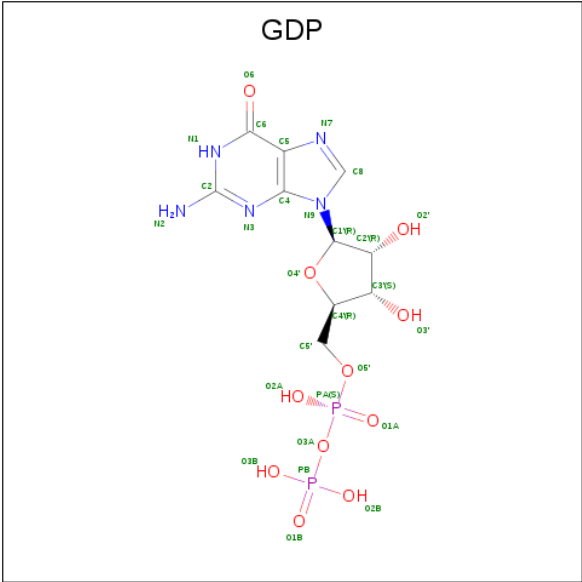
Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	ASN	deletion	UNP P47122
B	?	-	GLY	deletion	UNP P47122
B	?	-	ASP	deletion	UNP P47122
B	?	-	ASN	deletion	UNP P47122
B	?	-	GLY	deletion	UNP P47122
B	?	-	LEU	deletion	UNP P47122
B	?	-	GLY	deletion	UNP P47122
B	?	-	SER	deletion	UNP P47122
B	265	LYS	-	expression tag	UNP P47122
B	266	HIS	-	expression tag	UNP P47122
B	267	HIS	-	expression tag	UNP P47122
B	268	HIS	-	expression tag	UNP P47122
B	269	HIS	-	expression tag	UNP P47122
B	270	HIS	-	expression tag	UNP P47122
B	271	HIS	-	expression tag	UNP P47122
C	?	-	LEU	deletion	UNP P47122
C	?	-	ASN	deletion	UNP P47122
C	?	-	GLY	deletion	UNP P47122
C	?	-	ASP	deletion	UNP P47122
C	?	-	ASN	deletion	UNP P47122
C	?	-	GLY	deletion	UNP P47122
C	?	-	LEU	deletion	UNP P47122
C	?	-	GLY	deletion	UNP P47122
C	?	-	SER	deletion	UNP P47122
C	265	LYS	-	expression tag	UNP P47122
C	266	HIS	-	expression tag	UNP P47122
C	267	HIS	-	expression tag	UNP P47122
C	268	HIS	-	expression tag	UNP P47122
C	269	HIS	-	expression tag	UNP P47122
C	270	HIS	-	expression tag	UNP P47122
C	271	HIS	-	expression tag	UNP P47122
D	?	-	LEU	deletion	UNP P47122
D	?	-	ASN	deletion	UNP P47122
D	?	-	GLY	deletion	UNP P47122
D	?	-	ASP	deletion	UNP P47122
D	?	-	ASN	deletion	UNP P47122
D	?	-	GLY	deletion	UNP P47122
D	?	-	LEU	deletion	UNP P47122
D	?	-	GLY	deletion	UNP P47122
D	?	-	SER	deletion	UNP P47122
D	265	LYS	-	expression tag	UNP P47122
D	266	HIS	-	expression tag	UNP P47122

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Chain	Residue	Modelled	Actual	Comment	Reference
D	267	HIS	-	expression tag	UNP P47122
D	268	HIS	-	expression tag	UNP P47122
D	269	HIS	-	expression tag	UNP P47122
D	270	HIS	-	expression tag	UNP P47122
D	271	HIS	-	expression tag	UNP P47122
E	?	-	LEU	deletion	UNP P47122
E	?	-	ASN	deletion	UNP P47122
E	?	-	GLY	deletion	UNP P47122
E	?	-	ASP	deletion	UNP P47122
E	?	-	ASN	deletion	UNP P47122
E	?	-	GLY	deletion	UNP P47122
E	?	-	LEU	deletion	UNP P47122
E	?	-	GLY	deletion	UNP P47122
E	?	-	SER	deletion	UNP P47122
E	265	LYS	-	expression tag	UNP P47122
E	266	HIS	-	expression tag	UNP P47122
E	267	HIS	-	expression tag	UNP P47122
E	268	HIS	-	expression tag	UNP P47122
E	269	HIS	-	expression tag	UNP P47122
E	270	HIS	-	expression tag	UNP P47122
E	271	HIS	-	expression tag	UNP P47122
F	?	-	LEU	deletion	UNP P47122
F	?	-	ASN	deletion	UNP P47122
F	?	-	GLY	deletion	UNP P47122
F	?	-	ASP	deletion	UNP P47122
F	?	-	ASN	deletion	UNP P47122
F	?	-	GLY	deletion	UNP P47122
F	?	-	LEU	deletion	UNP P47122
F	?	-	GLY	deletion	UNP P47122
F	?	-	SER	deletion	UNP P47122
F	265	LYS	-	expression tag	UNP P47122
F	266	HIS	-	expression tag	UNP P47122
F	267	HIS	-	expression tag	UNP P47122
F	268	HIS	-	expression tag	UNP P47122
F	269	HIS	-	expression tag	UNP P47122
F	270	HIS	-	expression tag	UNP P47122
F	271	HIS	-	expression tag	UNP P47122

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).

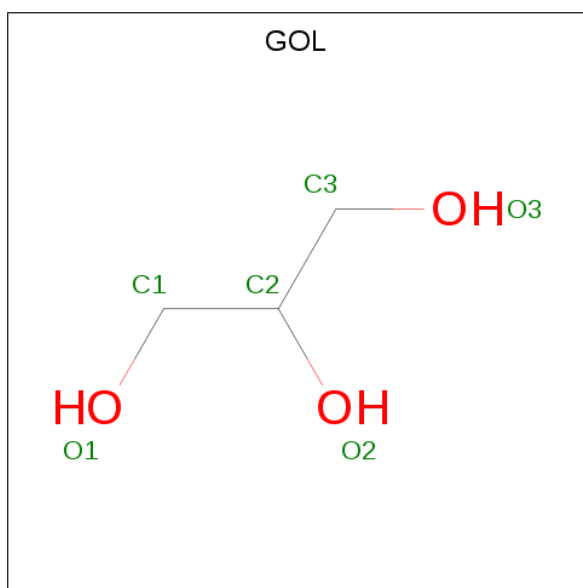


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	C	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	D	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	E	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
2	F	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		
3	F	1	Total	Mg	0	0
			1	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		

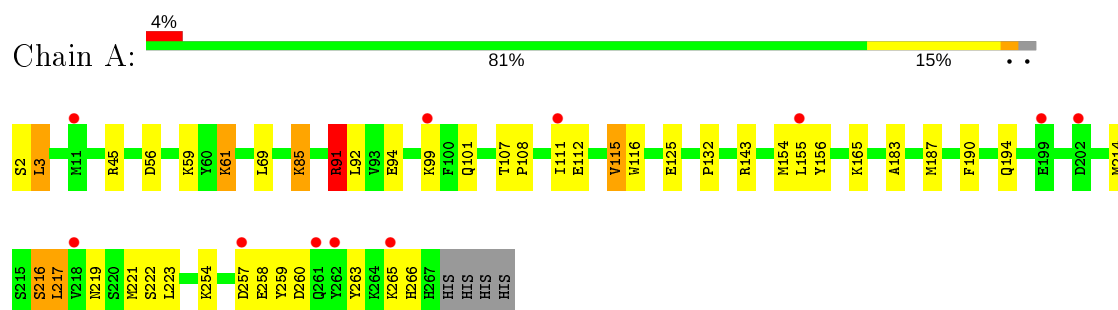
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	33	Total	O	0	0
			33	33		
5	B	30	Total	O	0	0
			30	30		
5	C	16	Total	O	0	0
			16	16		
5	D	16	Total	O	0	0
			16	16		
5	E	21	Total	O	0	0
			21	21		
5	F	5	Total	O	0	0
			5	5		

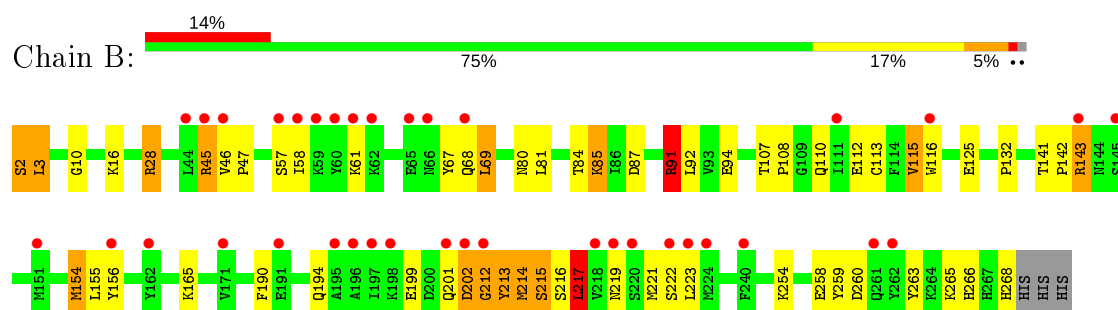
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

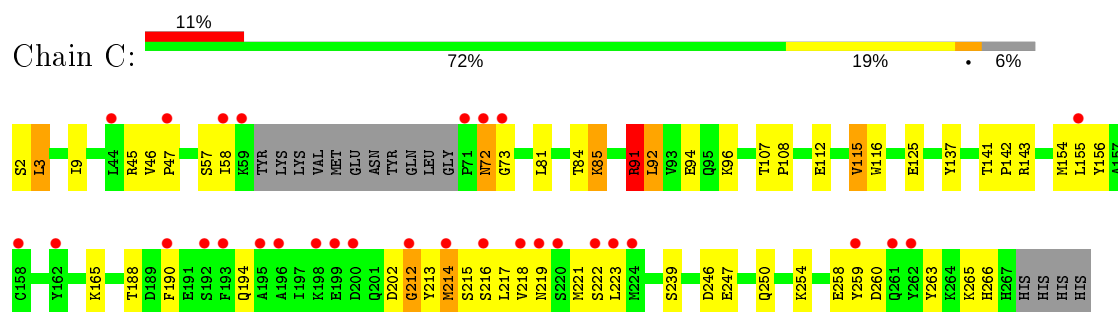
• Molecule 1: GPN-loop GTPase 1



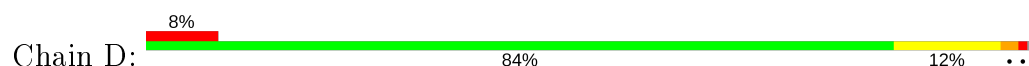
• Molecule 1: GPN-loop GTPase 1

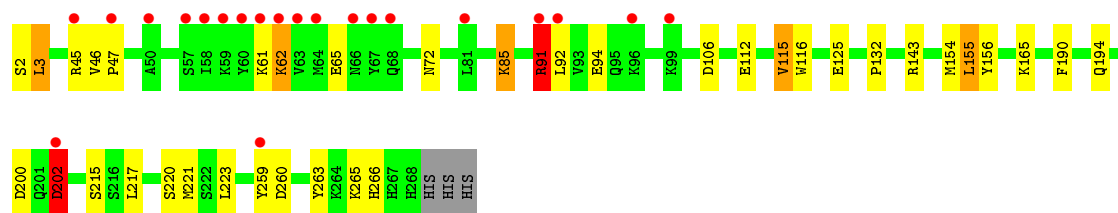


• Molecule 1: GPN-loop GTPase 1

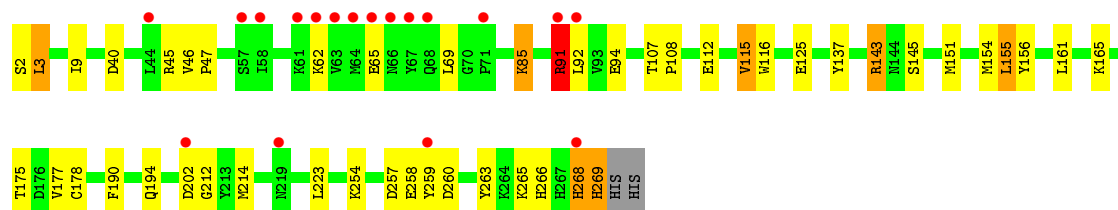
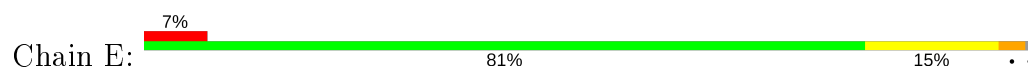


• Molecule 1: GPN-loop GTPase 1

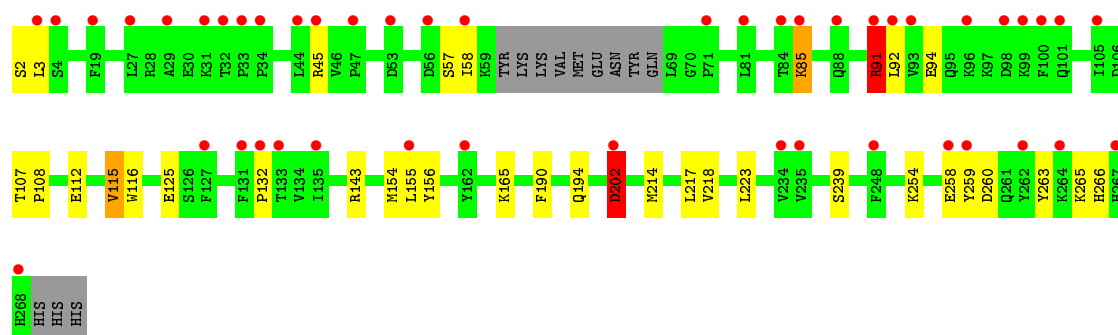
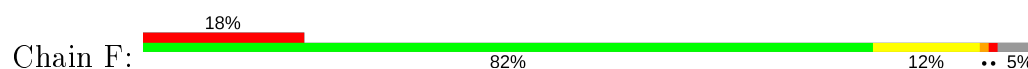




● Molecule 1: GPN-loop GTPase 1



● Molecule 1: GPN-loop GTPase 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	107.99Å 119.18Å 347.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.25 – 2.30 49.15 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.25-2.30) 99.8 (49.15-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.29Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.238 , 0.276 0.241 , 0.276	Depositor DCC
R_{free} test set	1988 reflections (2.00%)	wwPDB-VP
Wilson B-factor (Å ²)	54.6	Xtriage
Anisotropy	0.533	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 54.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12383	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.59% 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: GDP, GOL, MG

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	0.55	0/2075	0.65	2/2806 (0.1%)
1	B	0.58	0/2086	0.90	7/2821 (0.2%)
1	C	0.62	1/1977 (0.1%)	0.70	2/2673 (0.1%)
1	D	0.40	0/2086	0.60	3/2821 (0.1%)
1	E	0.57	0/2097	0.68	3/2836 (0.1%)
1	F	0.38	0/2000	0.58	1/2705 (0.0%)
All	All	0.52	1/12321 (0.0%)	0.69	18/16662 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	D	0	1
1	F	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	73	GLY	N-CA	-5.39	1.38	1.46

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	45	ARG	NE-CZ-NH1	-20.10	110.25	120.30
1	B	45	ARG	NE-CZ-NH2	19.68	130.14	120.30
1	E	91	ARG	NE-CZ-NH1	10.69	125.64	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	45	ARG	CD-NE-CZ	9.86	137.40	123.60
1	D	91	ARG	NE-CZ-NH1	9.20	124.90	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	72	ASN	Peptide
1	D	202	ASP	Peptide
1	F	202	ASP	Peptide

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	2032	0	2009	22	1
1	B	2042	0	2016	58	0
1	C	1937	0	1914	52	0
1	D	2042	0	2016	20	0
1	E	2052	0	2023	26	1
1	F	1959	0	1933	22	0
2	A	28	0	12	0	0
2	B	28	0	12	0	0
2	C	28	0	12	1	0
2	D	28	0	12	0	0
2	E	28	0	12	0	0
2	F	28	0	12	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	6	0	8	1	0
4	C	6	0	8	0	0
4	E	12	0	16	1	0
5	A	33	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	30	0	0	2	0
5	C	16	0	0	0	0
5	D	16	0	0	1	0
5	E	21	0	0	3	0
5	F	5	0	0	0	0
All	All	12383	0	12015	173	1

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 7.

The worst 5 of 173 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:B:91:ARG:HG3	1:D:202:ASP:HB3	1.23	1.13
1:C:202:ASP:O	1:C:213:TYR:N	1.85	1.07
1:C:91:ARG:HG3	1:F:202:ASP:HB3	1.48	0.94
1:B:143:ARG:NH2	5:B:401:HOH:O	2.01	0.92
1:C:91:ARG:CG	1:F:202:ASP:HB3	2.08	0.82

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:ASP:O	1:E:91:ARG:NH2[5_455]	2.13	0.07

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	255/261 (98%)	251 (98%)	4 (2%)	0	100	100
1	B	256/261 (98%)	250 (98%)	4 (2%)	2 (1%)	19	23

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	242/261 (93%)	236 (98%)	5 (2%)	1 (0%)	34	42
1	D	256/261 (98%)	252 (98%)	4 (2%)	0	100	100
1	E	257/261 (98%)	252 (98%)	4 (2%)	1 (0%)	34	42
1	F	245/261 (94%)	241 (98%)	4 (2%)	0	100	100
All	All	1511/1566 (96%)	1482 (98%)	25 (2%)	4 (0%)	41	50

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	212	GLY
1	C	212	GLY
1	B	213	TYR
1	E	212	GLY

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	A	231/235 (98%)	211 (91%)	20 (9%)	10	12
1	B	232/235 (99%)	209 (90%)	23 (10%)	8	9
1	C	221/235 (94%)	203 (92%)	18 (8%)	11	15
1	D	232/235 (99%)	211 (91%)	21 (9%)	9	11
1	E	233/235 (99%)	212 (91%)	21 (9%)	9	11
1	F	223/235 (95%)	206 (92%)	17 (8%)	13	16
All	All	1372/1410 (97%)	1252 (91%)	120 (9%)	10	12

5 of 120 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	143	ARG
1	D	72	ASN
1	F	112	GLU

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Mol	Chain	Res	Type
1	C	154	MET
1	C	260	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	194	GLN
1	F	194	GLN
1	D	194	GLN
1	B	201	GLN
1	C	250	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 6 are monoatomic - leaving 10 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	$\# Z > 2$
4	GOL	E	304	-	5,5,5	0.37	0	5,5,5	0.54	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GDP	F	301	3	24,30,30	1.19	2 (8%)	31,47,47	1.96	7 (22%)
2	GDP	D	301	3	24,30,30	1.29	2 (8%)	31,47,47	1.99	8 (25%)
4	GOL	A	303	-	5,5,5	0.36	0	5,5,5	0.63	0
4	GOL	C	303	-	5,5,5	0.36	0	5,5,5	0.29	0
4	GOL	E	303	-	5,5,5	0.35	0	5,5,5	0.21	0
2	GDP	C	301	-	24,30,30	1.39	3 (12%)	31,47,47	2.12	7 (22%)
2	GDP	A	301	3	24,30,30	1.18	2 (8%)	31,47,47	1.99	8 (25%)
2	GDP	E	301	3	24,30,30	1.32	2 (8%)	31,47,47	2.42	10 (32%)
2	GDP	B	301	3	24,30,30	1.23	2 (8%)	31,47,47	1.98	9 (29%)

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
4	GOL	E	304	-	-	1/4/4/4	-
2	GDP	F	301	3	-	5/12/32/32	0/3/3/3
2	GDP	D	301	3	-	4/12/32/32	0/3/3/3
4	GOL	A	303	-	-	2/4/4/4	-
4	GOL	C	303	-	-	2/4/4/4	-
4	GOL	E	303	-	-	2/4/4/4	-
2	GDP	C	301	-	-	3/12/32/32	0/3/3/3
2	GDP	A	301	3	-	0/12/32/32	0/3/3/3
2	GDP	E	301	3	-	1/12/32/32	0/3/3/3
2	GDP	B	301	3	-	0/12/32/32	0/3/3/3

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	GDP	C6-C5	5.28	1.50	1.41
2	E	301	GDP	C6-C5	4.49	1.49	1.41
2	D	301	GDP	C6-C5	4.42	1.49	1.41
2	B	301	GDP	C6-C5	4.23	1.48	1.41
2	A	301	GDP	C6-C5	4.16	1.48	1.41

The worst 5 of 49 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	301	GDP	C2-N3-C4	7.25	123.64	115.36
2	C	301	GDP	C2-N3-C4	6.38	122.65	115.36
2	A	301	GDP	C5-C6-N1	-5.13	116.42	123.43
2	B	301	GDP	C5-C6-N1	-4.97	116.63	123.43
2	D	301	GDP	C2-N3-C4	4.94	121.00	115.36

There are no chirality outliers.

5 of 20 torsion outliers are listed below:

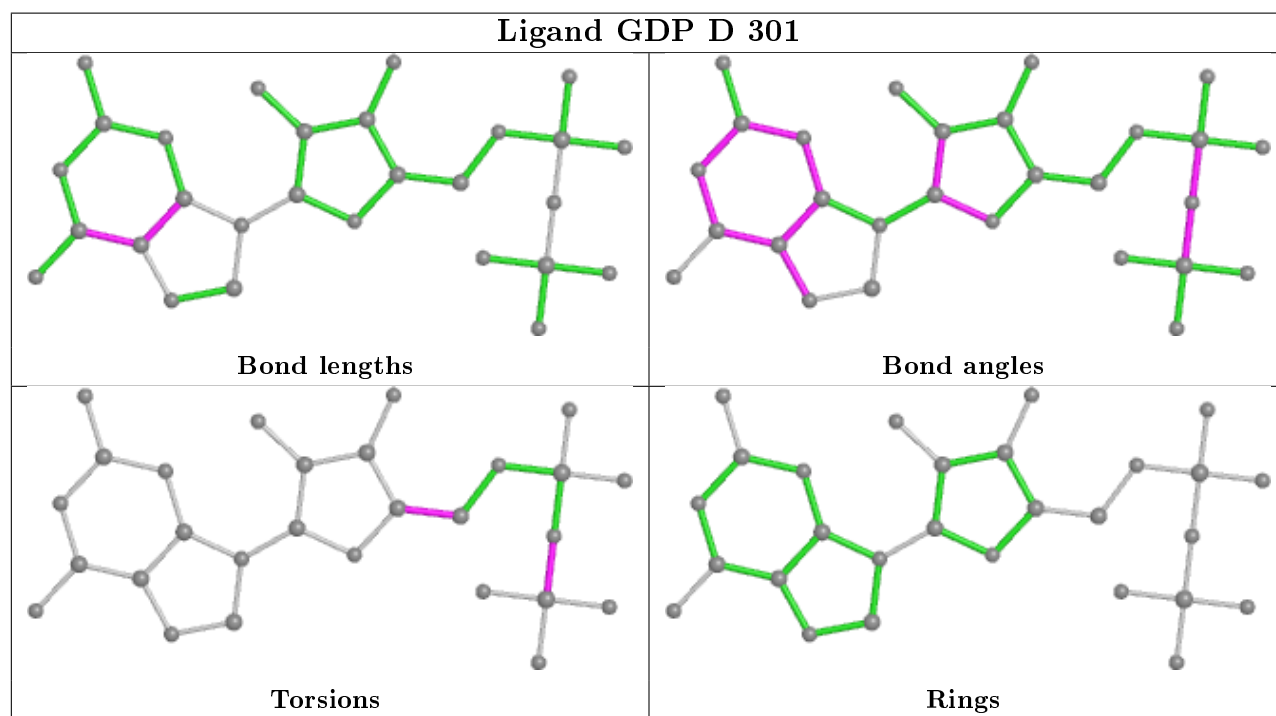
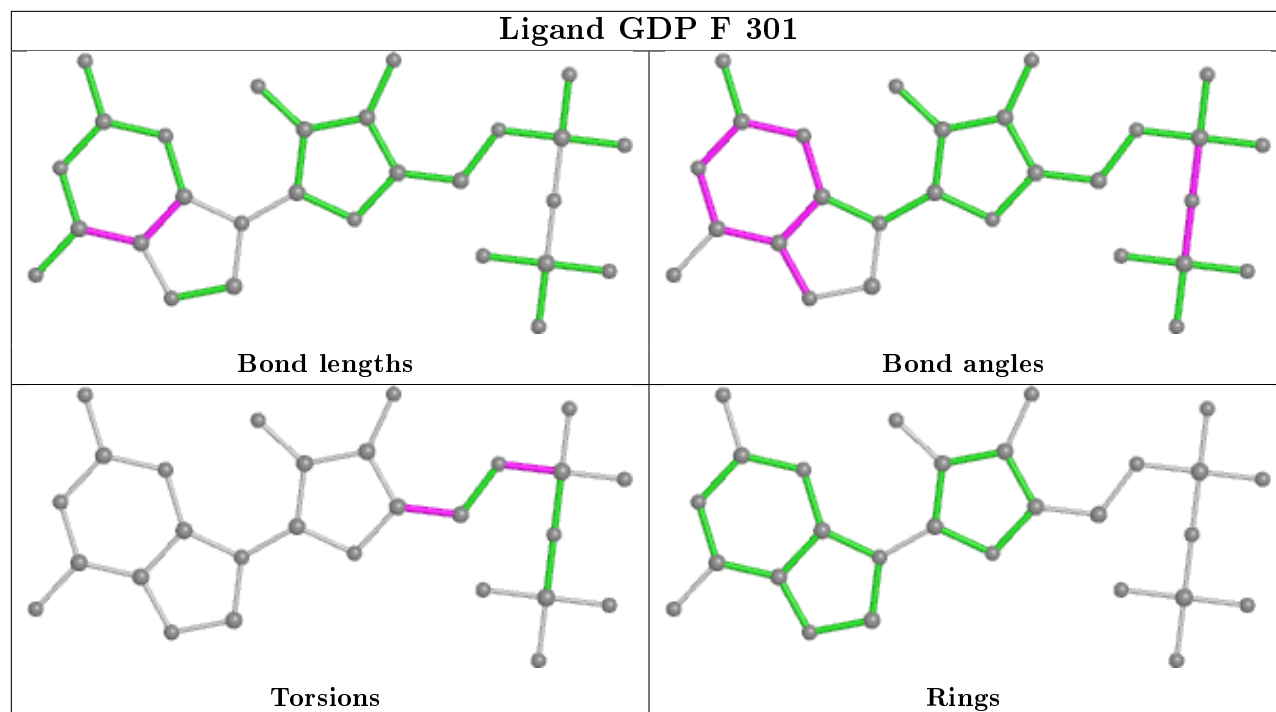
Mol	Chain	Res	Type	Atoms
2	F	301	GDP	C5'-O5'-PA-O1A
2	F	301	GDP	O4'-C4'-C5'-O5'
4	A	303	GOL	O1-C1-C2-C3
4	C	303	GOL	O1-C1-C2-C3
4	E	303	GOL	C1-C2-C3-O3

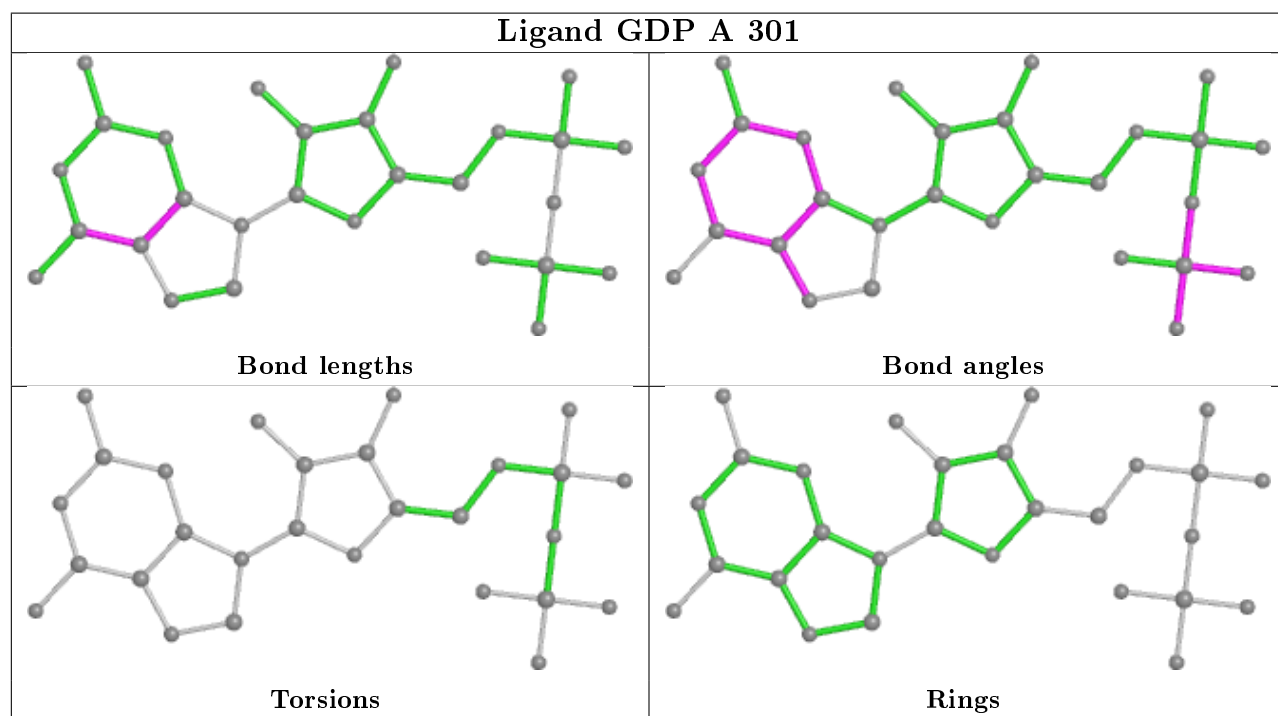
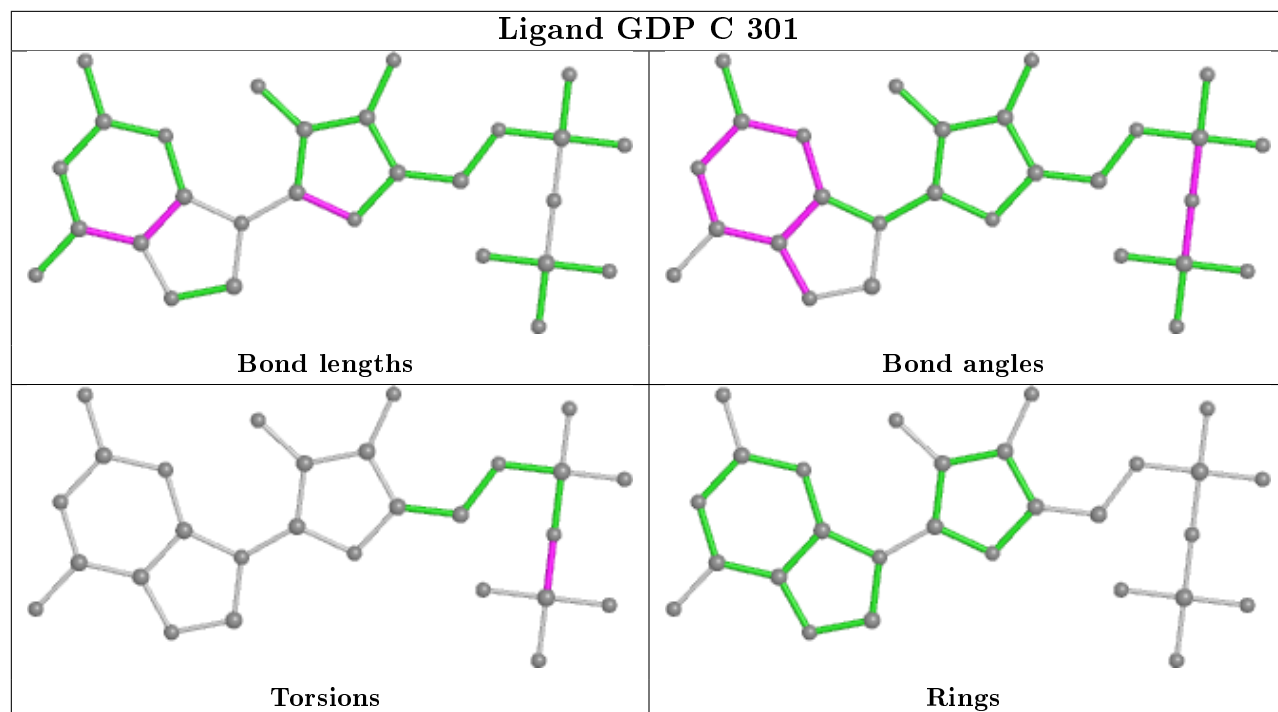
There are no ring outliers.

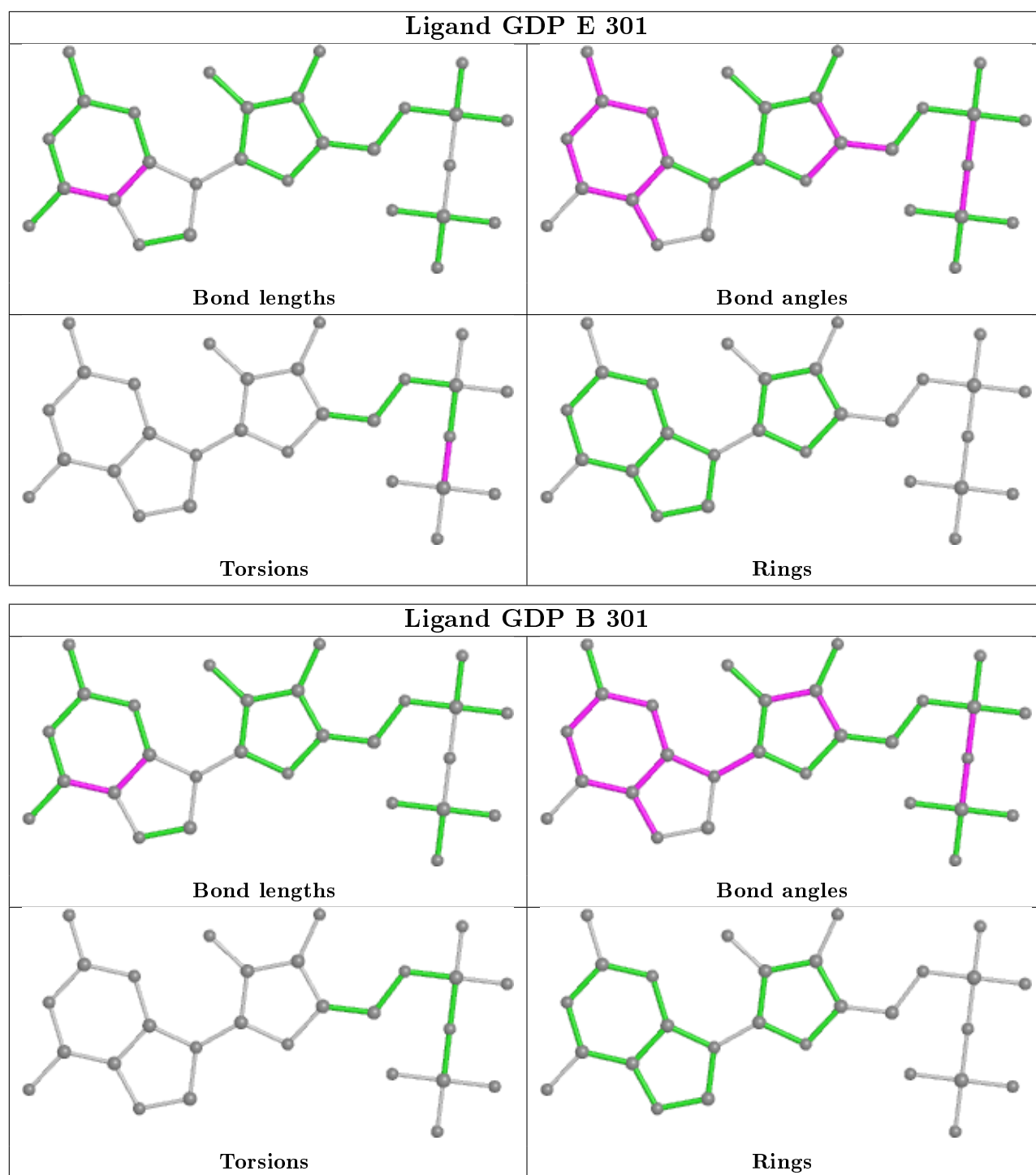
4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	301	GDP	1	0
4	A	303	GOL	1	0
4	E	303	GOL	1	0
2	C	301	GDP	1	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	A	257/261 (98%)	0.39	11 (4%) 35 42	42, 60, 106, 198	0
1	B	258/261 (98%)	1.01	37 (14%) 2 3	44, 69, 146, 237	0
1	C	246/261 (94%)	0.86	30 (12%) 4 6	48, 72, 158, 280	0
1	D	258/261 (98%)	0.62	21 (8%) 12 16	56, 83, 135, 186	0
1	E	259/261 (99%)	0.50	18 (6%) 16 22	44, 63, 127, 211	0
1	F	249/261 (95%)	1.19	46 (18%) 1 1	55, 104, 146, 195	0
All	All	1527/1566 (97%)	0.76	163 (10%) 6 8	42, 74, 141, 280	0

The worst 5 of 163 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	220	SER	14.7
1	D	67	TYR	11.1
1	B	223	LEU	10.0
1	C	223	LEU	9.7
1	D	63	VAL	8.7

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 carbohydrates in this entry.

6.4 Ligands ⓘ

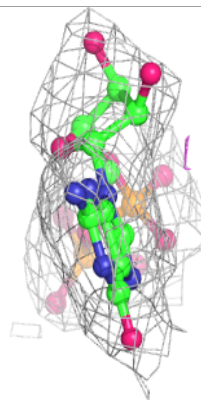
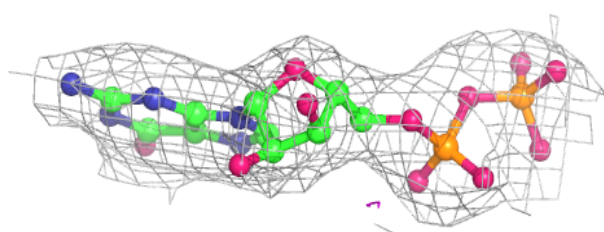
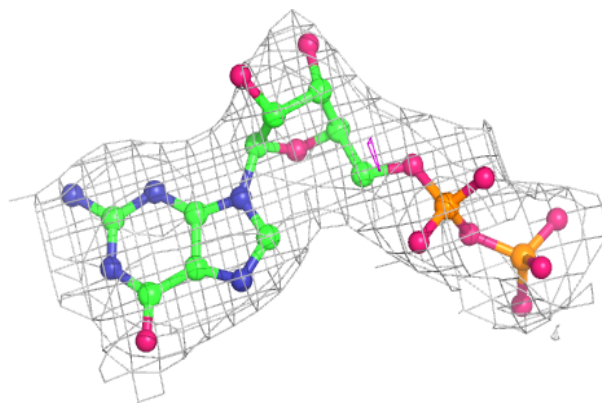
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	E	303	6/6	0.60	0.46	91,102,106,108	0
4	GOL	A	303	6/6	0.64	0.47	104,119,129,129	0
4	GOL	C	303	6/6	0.72	0.23	90,102,119,125	0
4	GOL	E	304	6/6	0.86	0.26	69,98,110,124	0
2	GDP	F	301	28/28	0.93	0.13	71,93,103,114	0
2	GDP	C	301	28/28	0.93	0.16	45,66,77,84	0
3	MG	D	302	1/1	0.93	0.05	65,65,65,65	0
3	MG	A	302	1/1	0.94	0.24	55,55,55,55	0
3	MG	B	302	1/1	0.94	0.12	52,52,52,52	0
2	GDP	E	301	28/28	0.95	0.20	54,61,86,110	0
3	MG	C	302	1/1	0.96	0.19	63,63,63,63	0
2	GDP	D	301	28/28	0.96	0.11	55,68,91,136	0
3	MG	F	302	1/1	0.97	0.21	85,85,85,85	0
3	MG	E	302	1/1	0.97	0.07	61,61,61,61	0
2	GDP	B	301	28/28	0.97	0.14	51,53,63,65	0
2	GDP	A	301	28/28	0.98	0.18	57,59,77,79	0

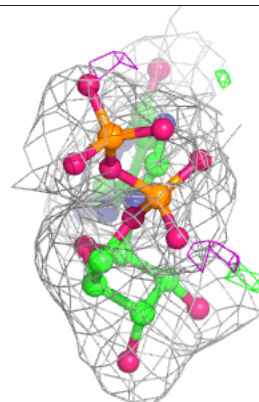
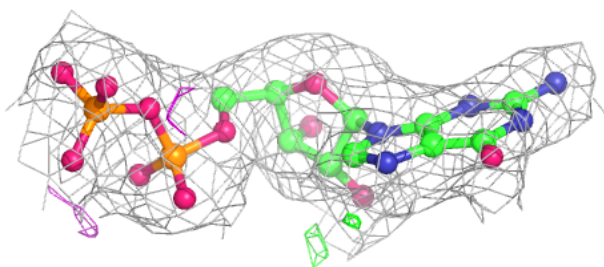
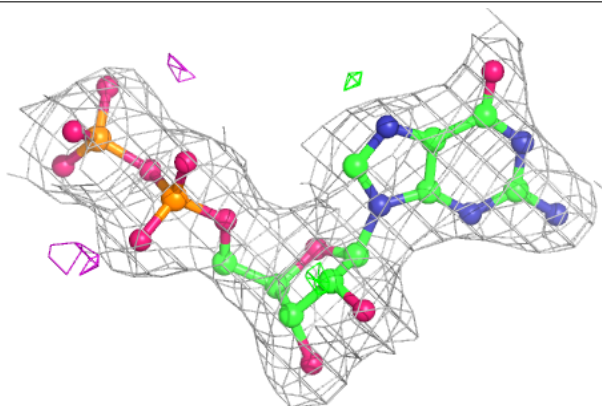
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GDP F 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

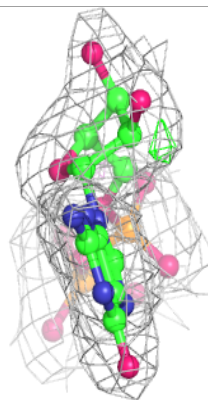
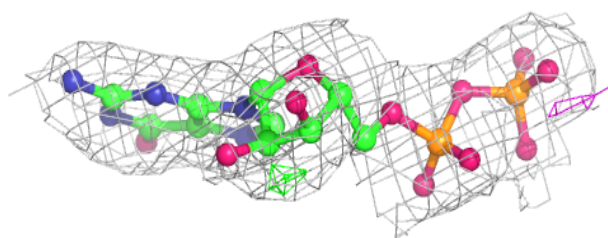
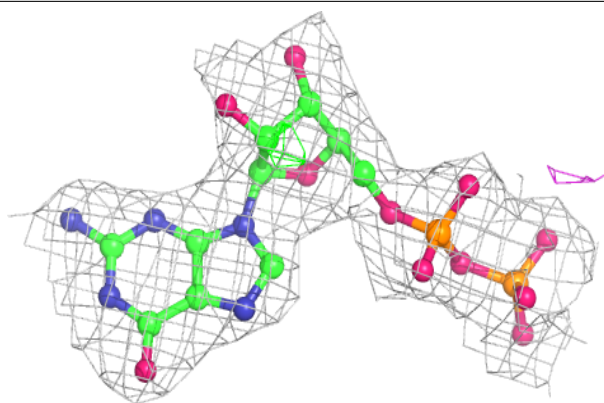
**Electron density around GDP C 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

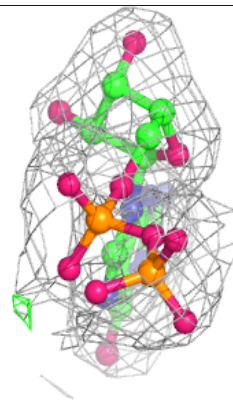
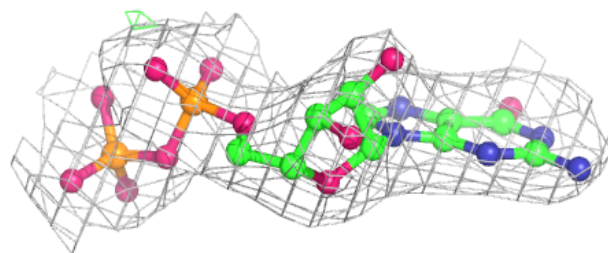
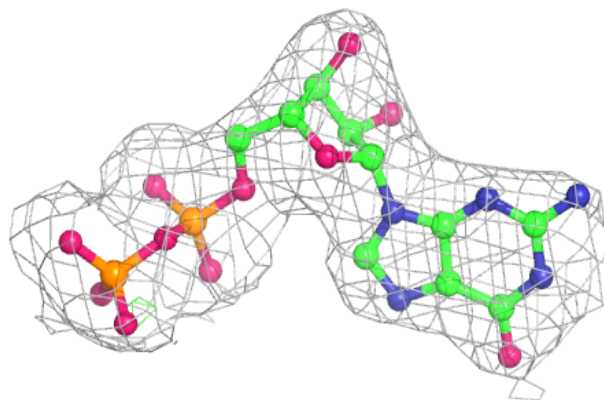


Electron density around GDP E 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

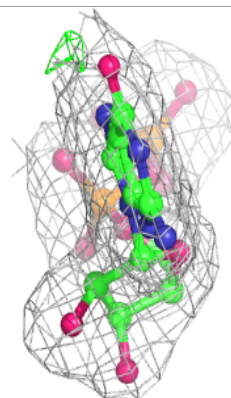
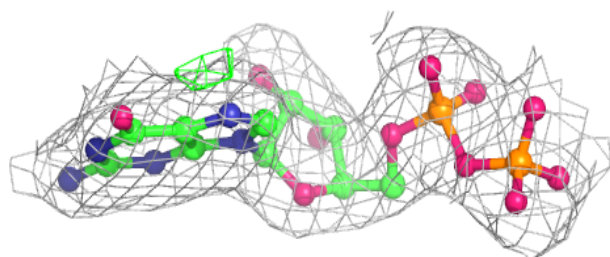
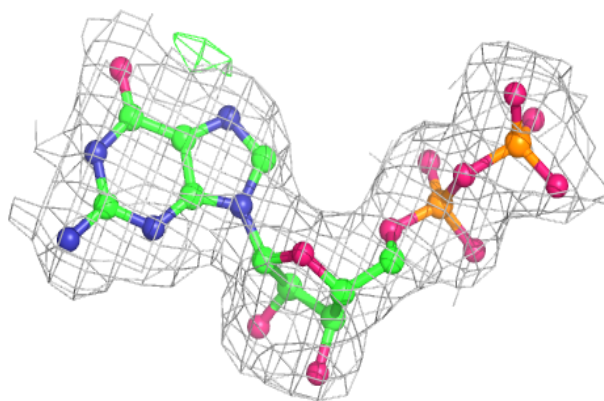
**Electron density around GDP D 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

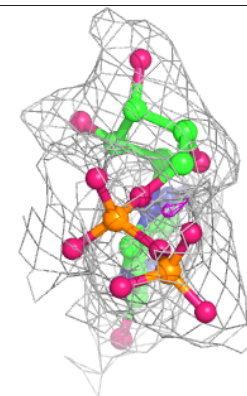
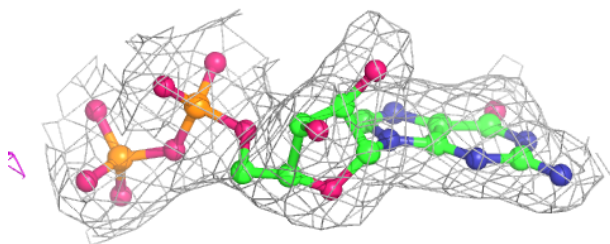
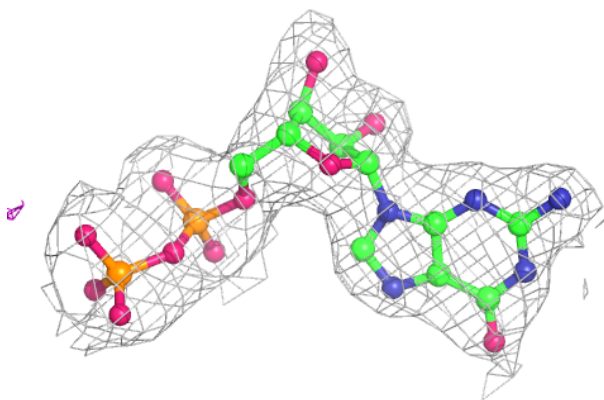


Electron density around GDP B 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around GDP A 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.