



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 10:56 pm BST

PDB ID : 6B9D
Title : Human ATL1 mutant - R77A bound to GDP
Authors : O'Donnell, J.P.; Sondermann, H.
Deposited on : 2017-10-10
Resolution : 1.95 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 : 2.11
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

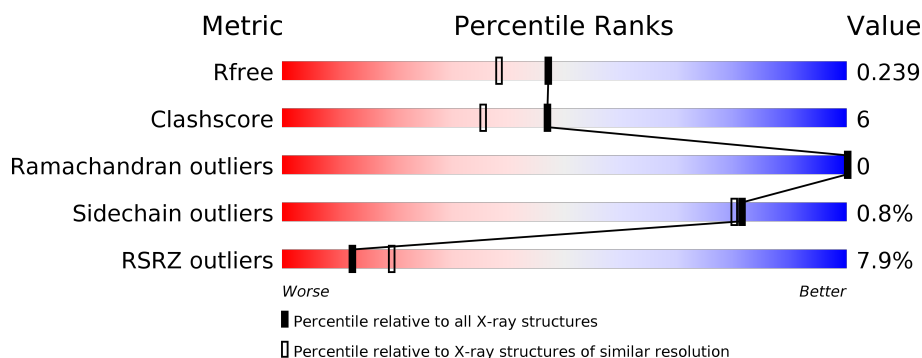
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	458	<div> <div>7%</div> <div> <div></div> <div>71%</div> <div>12%</div> <div>17%</div> </div> </div>
1	B	458	<div> <div>6%</div> <div> <div></div> <div>70%</div> <div>13%</div> <div>17%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6247 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 Atlastin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	0	0
			2929	1879	487	550	13			
1	B	380	Total	C	N	O	S	0	0	0
			2943	1893	491	546	13			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q8WXF7
A	77	ALA	ARG	engineered mutation	UNP Q8WXF7
A	447	ALA	-	expression tag	UNP Q8WXF7
A	448	ALA	-	expression tag	UNP Q8WXF7
A	449	ALA	-	expression tag	UNP Q8WXF7
A	450	LEU	-	expression tag	UNP Q8WXF7
A	451	GLU	-	expression tag	UNP Q8WXF7
A	452	HIS	-	expression tag	UNP Q8WXF7
A	453	HIS	-	expression tag	UNP Q8WXF7
A	454	HIS	-	expression tag	UNP Q8WXF7
A	455	HIS	-	expression tag	UNP Q8WXF7
A	456	HIS	-	expression tag	UNP Q8WXF7
A	457	HIS	-	expression tag	UNP Q8WXF7
B	0	SER	-	expression tag	UNP Q8WXF7
B	77	ALA	ARG	engineered mutation	UNP Q8WXF7
B	447	ALA	-	expression tag	UNP Q8WXF7
B	448	ALA	-	expression tag	UNP Q8WXF7
B	449	ALA	-	expression tag	UNP Q8WXF7
B	450	LEU	-	expression tag	UNP Q8WXF7
B	451	GLU	-	expression tag	UNP Q8WXF7
B	452	HIS	-	expression tag	UNP Q8WXF7
B	453	HIS	-	expression tag	UNP Q8WXF7
B	454	HIS	-	expression tag	UNP Q8WXF7
B	455	HIS	-	expression tag	UNP Q8WXF7
B	456	HIS	-	expression tag	UNP Q8WXF7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	457	HIS	-	expression tag	UNP Q8WXF7

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- The image displays the chemical structure of Guanosine Diphosphate (GDP). It consists of a guanine base (a purine derivative) linked to a ribose sugar, which is in turn linked to two phosphate groups. The guanine base is shown with atoms labeled N1, N3, N7, C2, C4, C6, and C8. The ribose sugar is shown with atoms labeled C1', C2', C3', C4', and C5'. The two phosphate groups are shown with atoms labeled P A(5), P B, and various oxygen atoms (O1A, O1B, O2A, O2B, O3A, O3B, O4', O5'). The structure is rendered in a 3D perspective, with the guanine base in blue and the ribose sugar and phosphate groups in red and green.

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

- | Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|-----------------|---------|---------|
| 3 | B | 1 | Total Mg
1 1 | 0 | 0 |
| 3 | A | 1 | Total Mg
1 1 | 0 | 0 |

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

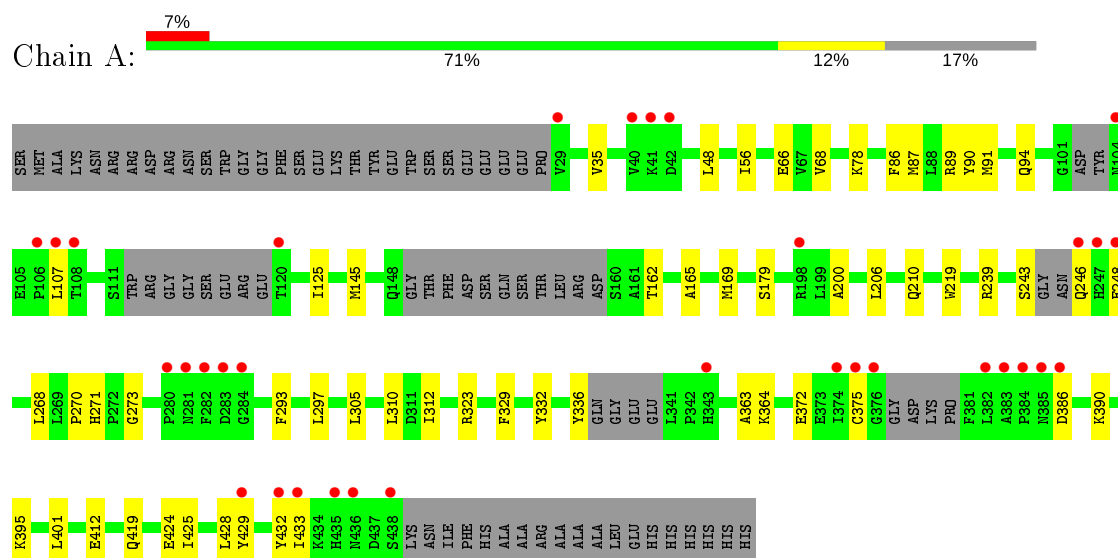
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	147	Total	O	0	0
			147	147		
5	B	134	Total	O	0	0
			134	134		

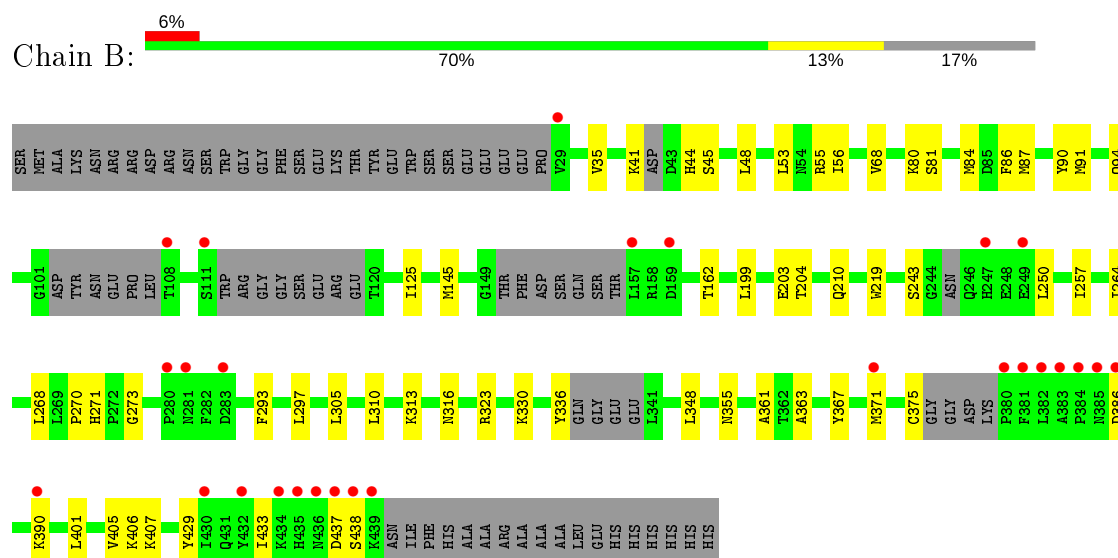
3 Residue-property plots [i](#)

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: Atlastin-1



• Molecule 1: Atlastin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	51.06 Å 68.43 Å 75.88 Å 117.00° 89.81° 99.04°	Depositor
Resolution (Å)	25.18 – 1.95 42.36 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.3 (25.18-1.95) 84.6 (42.36-1.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.42 (at 1.79 Å)	Xtriage
Refinement program	PHENIX (1.12_2829)	Depositor
R, R_{free}	0.204 , 0.239 0.204 , 0.239	Depositor DCC
R_{free} test set	1980 reflections (2.48%)	wwPDB-VP
Wilson B-factor (Å ²)	24.3	Xtriage
Anisotropy	0.467	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 56.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6247	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.84% 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

5.1 Standard geometry

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	0.38	0/2983	0.54	0/4033
1	B	0.39	0/2998	0.55	0/4049
All	All	0.39	0/5981	0.55	0/8082

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

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	2929	0	2819	35	0
1	B	2943	0	2851	37	0
2	A	28	0	12	0	0
2	B	28	0	12	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	18	0	24	2	0
4	B	18	0	24	3	0
5	A	147	0	0	3	0
5	B	134	0	0	1	0
All	All	6247	0	5742	72	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 6.

All (72) 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:125:ILE:HG22	1:B:145:MET:HG2	1.68	0.74
1:A:68:VAL:HG21	1:A:310:LEU:HB3	1.72	0.71
1:B:243:SER:HB3	1:B:250:LEU:HD23	1.72	0.70
1:B:68:VAL:HG21	1:B:310:LEU:HB3	1.77	0.66
1:B:330:LYS:HB2	4:B:505:GOL:H31	1.79	0.64
1:A:271:HIS:CD2	1:A:273:GLY:H	2.17	0.63
1:B:271:HIS:CD2	1:B:273:GLY:H	2.17	0.62
1:B:91:MET:HE1	1:B:305:LEU:HB3	1.82	0.61
1:A:125:ILE:HG22	1:A:145:MET:HG2	1.84	0.60
1:B:323:ARG:NH1	5:B:603:HOH:O	2.32	0.59
1:B:355:ASN:HD21	1:B:407:LYS:HA	1.68	0.58
1:A:87:MET:HE3	1:A:305:LEU:HD11	1.85	0.58
1:A:48:LEU:HD11	1:A:329:PHE:HB3	1.87	0.56
1:B:35:VAL:HG21	1:B:56:ILE:HD11	1.88	0.56
1:B:386:ASP:OD2	1:B:390:LYS:NZ	2.40	0.55
1:B:94:GLN:HB2	4:B:504:GOL:H32	1.87	0.55
1:A:372:GLU:HG2	1:A:432:TYR:OH	2.08	0.54
1:B:355:ASN:ND2	1:B:407:LYS:HA	2.23	0.54
1:A:270:PRO:HG2	1:A:293:PHE:HA	1.90	0.54
1:A:91:MET:HE1	1:A:305:LEU:HB3	1.91	0.53
1:B:204:THR:HB	1:B:316:ASN:HD22	1.73	0.52
1:B:271:HIS:HD2	1:B:273:GLY:H	1.57	0.52
1:A:200:ALA:HB1	4:A:504:GOL:H31	1.91	0.52
1:A:91:MET:CE	1:A:94:GLN:HE21	2.23	0.51
1:A:271:HIS:HD2	1:A:273:GLY:H	1.56	0.51
1:B:437:ASP:OD1	1:B:438:SER:N	2.44	0.51
1:A:419:GLN:NE2	5:A:607:HOH:O	2.38	0.51
1:B:48:LEU:HD11	1:B:53:LEU:HD22	1.93	0.51
1:A:243:SER:HB3	1:A:246:GLN:N	2.27	0.50
1:A:395:LYS:HG3	1:A:425:ILE:HG21	1.92	0.50
1:A:206:LEU:O	4:A:504:GOL:O3	2.31	0.49
1:A:364:LYS:HE3	1:A:424:GLU:HB3	1.94	0.49
1:A:219:TRP:CE2	1:A:268:LEU:HD13	2.48	0.49
1:A:364:LYS:HE2	1:A:428:LEU:HD11	1.95	0.48
1:A:323:ARG:NH1	5:A:609:HOH:O	2.39	0.48
1:B:41:LYS:N	1:B:45:SER:O	2.43	0.48
1:A:35:VAL:HG21	1:A:56:ILE:HD11	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:199:LEU:HB3	1:B:348:LEU:HD22	1.95	0.48
1:B:257:ILE:HG22	1:B:264:ILE:CD1	2.43	0.48
1:A:363:ALA:HA	1:A:401:LEU:HD23	1.98	0.46
1:B:367:TYR:CZ	1:B:371:MET:HG3	2.51	0.46
1:A:78:LYS:HD2	1:A:179:SER:HB2	1.98	0.45
1:A:412:GLU:O	5:A:601:HOH:O	2.21	0.45
1:B:204:THR:HB	1:B:316:ASN:ND2	2.31	0.45
1:B:363:ALA:HA	1:B:401:LEU:HD23	1.99	0.45
1:A:248:GLU:H	1:A:248:GLU:CD	2.21	0.45
1:A:210:GLN:NE2	1:A:312:ILE:O	2.44	0.44
1:B:162:THR:HG23	1:B:336:TYR:HD1	1.82	0.44
1:A:332:TYR:O	1:A:336:TYR:HD2	1.99	0.44
1:A:89:ARG:HH21	1:A:107:LEU:HA	1.82	0.44
1:B:90:TYR:CD2	1:B:91:MET:CE	3.00	0.44
1:B:86:PHE:CG	1:B:297:LEU:HD21	2.53	0.43
1:B:41:LYS:CB	1:B:44:HIS:H	2.31	0.43
1:B:55:ARG:HA	4:B:503:GOL:H11	2.01	0.43
1:B:80:LYS:O	1:B:84:MET:HG3	2.19	0.43
1:B:84:MET:HA	1:B:87:MET:HE3	2.01	0.43
1:A:165:ALA:O	1:A:169:MET:HG2	2.19	0.43
1:A:90:TYR:CD2	1:A:91:MET:CE	3.02	0.43
1:A:386:ASP:OD2	1:A:390:LYS:NZ	2.52	0.42
1:A:429:TYR:CZ	1:A:433:ILE:HD11	2.54	0.42
1:B:219:TRP:CE2	1:B:268:LEU:HD13	2.54	0.42
1:A:86:PHE:CG	1:A:297:LEU:HD21	2.55	0.42
1:B:270:PRO:HG2	1:B:293:PHE:HA	2.02	0.42
1:A:66:GLU:OE2	1:A:312:ILE:HD13	2.19	0.41
1:B:84:MET:HA	1:B:87:MET:CE	2.50	0.41
1:B:429:TYR:CZ	1:B:433:ILE:HD11	2.55	0.41
1:B:429:TYR:CE2	1:B:433:ILE:HD11	2.56	0.41
1:B:401:LEU:O	1:B:405:VAL:HG23	2.20	0.41
1:A:162:THR:HG23	1:A:336:TYR:CD1	2.55	0.41
1:A:239:ARG:HH11	1:A:239:ARG:HD3	1.75	0.40
1:B:323:ARG:CZ	1:B:361:ALA:HB1	2.51	0.40
1:B:210:GLN:HE22	1:B:313:LYS:HD2	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	365/458 (80%)	356 (98%)	9 (2%)	0	100	100
1	B	364/458 (80%)	356 (98%)	8 (2%)	0	100	100
All	All	729/916 (80%)	712 (98%)	17 (2%)	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	A	303/402 (75%)	302 (100%)	1 (0%)	92	92
1	B	306/402 (76%)	302 (99%)	4 (1%)	69	65
All	All	609/804 (76%)	604 (99%)	5 (1%)	81	80

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	375	CYS
1	B	81	SER
1	B	203	GLU
1	B	375	CYS
1	B	406	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	GLN
1	A	191	GLN
1	A	271	HIS
1	A	318	ASN
1	A	419	GLN
1	B	54	ASN
1	B	174	GLN
1	B	210	GLN
1	B	271	HIS
1	B	316	ASN
1	B	349	GLN
1	B	355	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 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	B	505	-	5,5,5	1.12	0	5,5,5	0.86	0
4	GOL	A	505	-	5,5,5	0.91	0	5,5,5	1.10	0
2	GDP	B	501	-	24,30,30	1.25	3 (12%)	31,47,47	2.00	10 (32%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	B	503	-	5,5,5	0.99	0	5,5,5	0.86	0
4	GOL	B	504	-	5,5,5	1.07	0	5,5,5	1.00	0
4	GOL	A	503	-	5,5,5	0.83	0	5,5,5	1.05	0
2	GDP	A	501	-	24,30,30	1.34	2 (8%)	31,47,47	1.99	9 (29%)
4	GOL	A	504	-	5,5,5	0.77	0	5,5,5	1.03	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
4	GOL	B	505	-	-	2/4/4/4	-
4	GOL	A	505	-	-	0/4/4/4	-
2	GDP	B	501	-	-	1/12/32/32	0/3/3/3
4	GOL	B	503	-	-	4/4/4/4	-
4	GOL	B	504	-	-	2/4/4/4	-
4	GOL	A	503	-	-	2/4/4/4	-
2	GDP	A	501	-	-	2/12/32/32	0/3/3/3
4	GOL	A	504	-	-	3/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	GDP	C6-C5	4.56	1.49	1.41
2	B	501	GDP	C6-C5	4.03	1.48	1.41
2	B	501	GDP	C5-C4	2.24	1.46	1.40
2	A	501	GDP	C5-C4	2.18	1.46	1.40
2	B	501	GDP	C6-N1	2.04	1.36	1.33

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	GDP	C2-N3-C4	5.18	121.27	115.36
2	A	501	GDP	C2-N3-C4	4.77	120.81	115.36
2	A	501	GDP	C6-C5-C4	-4.64	116.37	120.80
2	B	501	GDP	C6-C5-C4	-4.53	116.47	120.80
2	A	501	GDP	C5-C6-N1	-3.84	118.19	123.43
2	A	501	GDP	C6-N1-C2	3.68	121.77	115.93
2	B	501	GDP	C5-C6-N1	-3.65	118.44	123.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	GDP	C6-N1-C2	3.60	121.64	115.93
2	B	501	GDP	N3-C2-N1	-3.38	122.72	127.22
2	A	501	GDP	N3-C2-N1	-3.08	123.12	127.22
2	A	501	GDP	C4-C5-N7	-2.80	106.48	109.40
2	B	501	GDP	O4'-C1'-C2'	-2.77	102.88	106.93
2	A	501	GDP	C1'-N9-C4	-2.71	121.88	126.64
2	A	501	GDP	O4'-C1'-C2'	-2.38	103.45	106.93
2	B	501	GDP	C4-C5-N7	-2.29	107.02	109.40
2	B	501	GDP	N2-C2-N1	2.25	120.75	117.25
2	B	501	GDP	PA-O3A-PB	-2.20	125.29	132.83
2	A	501	GDP	N2-C2-N1	2.15	120.59	117.25
2	B	501	GDP	O3B-PB-O3A	2.00	111.36	104.64

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	505	GOL	C1-C2-C3-O3
4	B	503	GOL	O1-C1-C2-C3
4	B	503	GOL	C1-C2-C3-O3
4	B	503	GOL	O2-C2-C3-O3
4	B	504	GOL	O1-C1-C2-C3
4	A	503	GOL	O1-C1-C2-C3
4	A	504	GOL	O1-C1-C2-C3
4	A	504	GOL	C1-C2-C3-O3
4	B	505	GOL	O2-C2-C3-O3
4	B	504	GOL	O1-C1-C2-O2
4	A	504	GOL	O1-C1-C2-O2
2	B	501	GDP	PA-O3A-PB-O1B
2	A	501	GDP	PA-O3A-PB-O3B
4	B	503	GOL	O1-C1-C2-O2
4	A	503	GOL	O1-C1-C2-O2
2	A	501	GDP	PA-O3A-PB-O2B

There are no ring outliers.

4 monomers are involved in 5 short contacts:

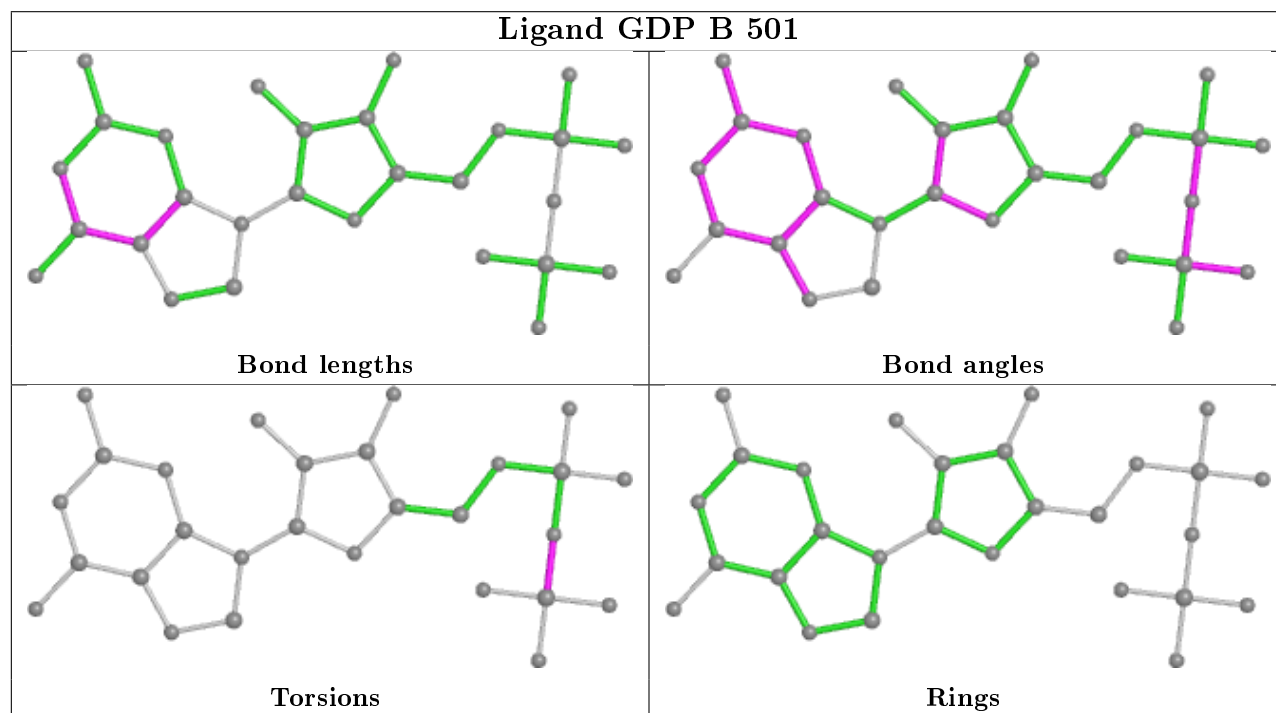
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	505	GOL	1	0
4	B	503	GOL	1	0
4	B	504	GOL	1	0

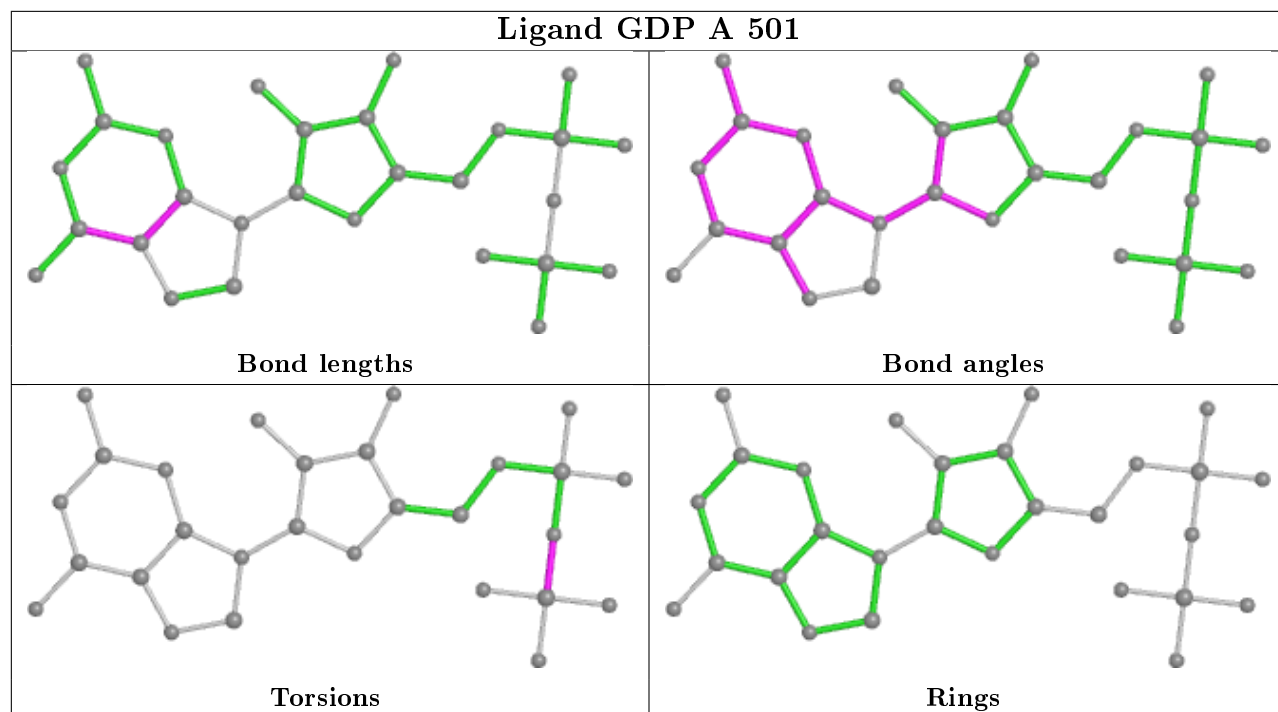
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	504	GOL	2	0

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	379/458 (82%)	0.37	33 (8%) 10 16	22, 40, 74, 91	0
1	B	380/458 (82%)	0.39	27 (7%) 16 24	22, 40, 76, 100	0
All	All	759/916 (82%)	0.38	60 (7%) 12 19	22, 40, 76, 100	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	383	ALA	6.8
1	B	437	ASP	6.6
1	B	382	LEU	6.4
1	B	384	PRO	5.6
1	A	382	LEU	5.3
1	B	435	HIS	5.2
1	B	29	VAL	5.2
1	A	383	ALA	4.9
1	B	438	SER	4.7
1	A	120	THR	4.6
1	A	384	PRO	4.6
1	A	42	ASP	4.6
1	A	107	LEU	4.6
1	A	436	ASN	4.5
1	B	385	ASN	3.8
1	A	280	PRO	3.8
1	A	343	HIS	3.7
1	A	40	VAL	3.7
1	B	157	LEU	3.6
1	A	106	PRO	3.6
1	B	386	ASP	3.6
1	A	435	HIS	3.5
1	B	381	PHE	3.5
1	B	111	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	29	VAL	3.4
1	A	281	ASN	3.3
1	A	284	GLY	3.3
1	B	436	ASN	3.3
1	A	433	ILE	3.3
1	B	439	LYS	3.1
1	B	371	MET	3.1
1	A	385	ASN	3.0
1	A	374	ILE	3.0
1	A	247	HIS	3.0
1	A	283	ASP	3.0
1	A	376	GLY	3.0
1	B	432	TYR	3.0
1	A	282	PHE	2.9
1	A	429	TYR	2.8
1	A	246	GLN	2.8
1	A	108	THR	2.8
1	B	281	ASN	2.7
1	A	386	ASP	2.6
1	A	438	SER	2.6
1	B	247	HIS	2.5
1	A	248	GLU	2.4
1	B	390	LYS	2.3
1	B	108	THR	2.3
1	B	280	PRO	2.3
1	A	432	TYR	2.3
1	A	104	ASN	2.2
1	A	41	LYS	2.2
1	B	434	LYS	2.2
1	B	430	ILE	2.2
1	B	159	ASP	2.2
1	A	375	CYS	2.2
1	B	283	ASP	2.1
1	B	249	GLU	2.1
1	A	198	ARG	2.1
1	B	380	PRO	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

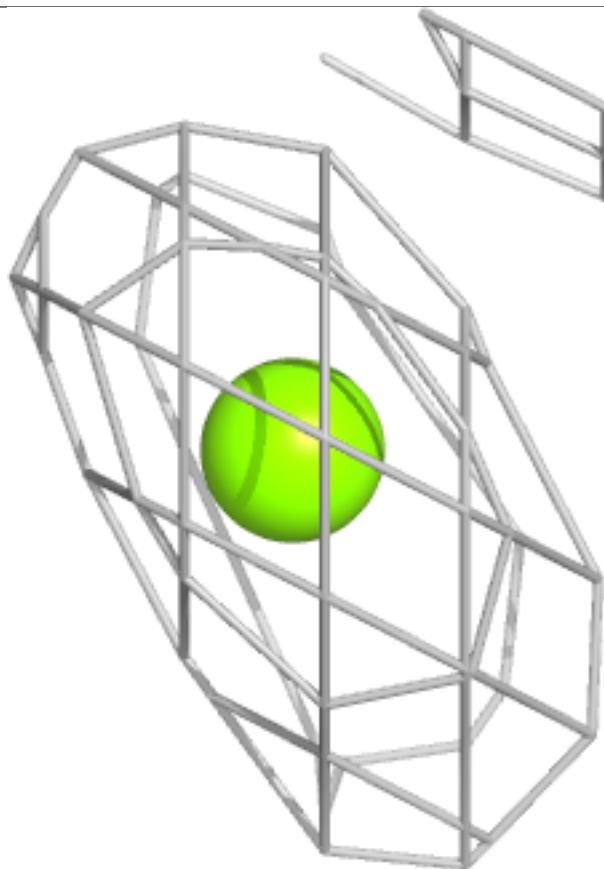
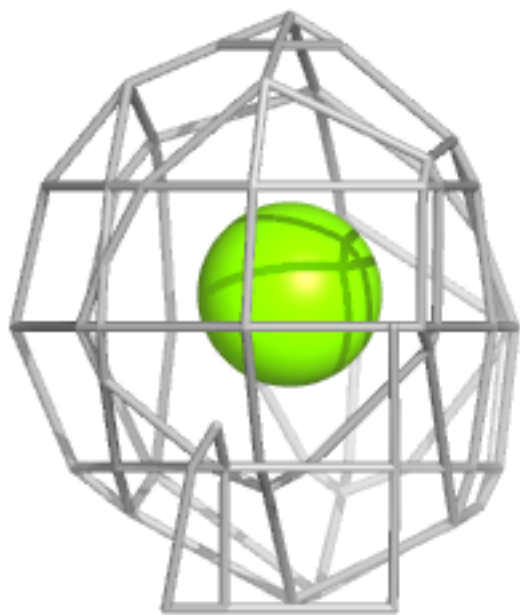
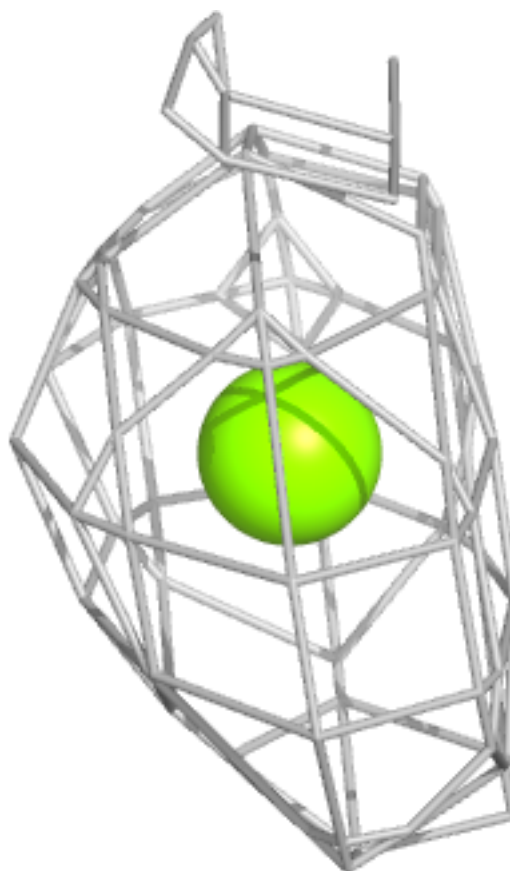
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MG	B	502	1/1	0.71	0.22	59,59,59,59	0
4	GOL	A	504	6/6	0.73	0.19	50,58,59,61	0
4	GOL	B	503	6/6	0.83	0.16	55,63,63,69	0
4	GOL	B	504	6/6	0.83	0.20	46,54,58,66	0
4	GOL	B	505	6/6	0.89	0.21	34,53,58,60	0
4	GOL	A	503	6/6	0.90	0.16	26,40,49,49	0
4	GOL	A	505	6/6	0.91	0.17	36,41,51,53	0
3	MG	A	502	1/1	0.93	0.10	55,55,55,55	0
2	GDP	A	501	28/28	0.97	0.10	22,31,39,41	0
2	GDP	B	501	28/28	0.97	0.10	26,34,39,48	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.

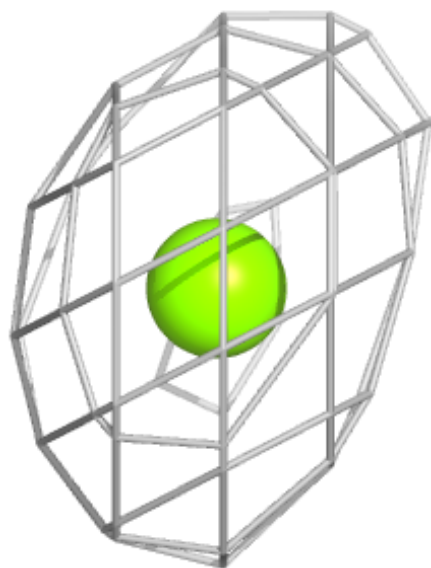
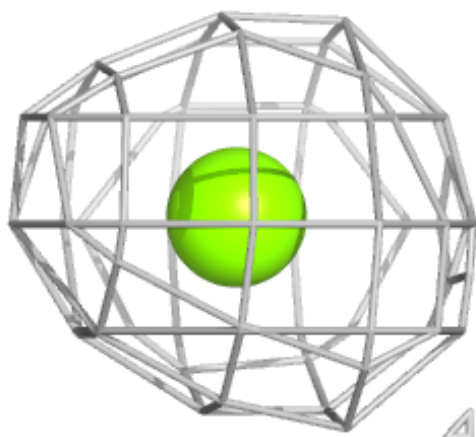
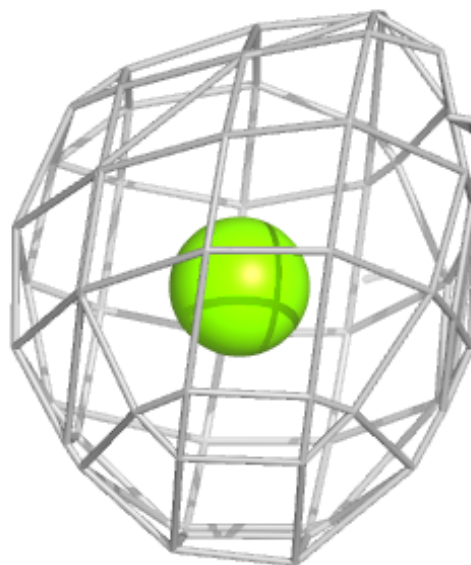
Electron density around MG B 502:

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



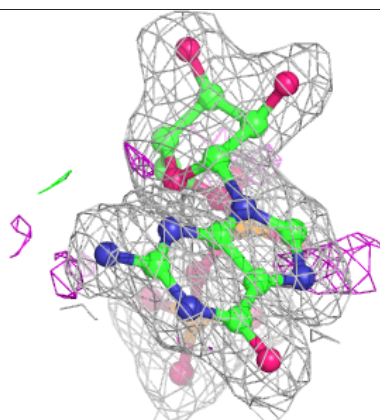
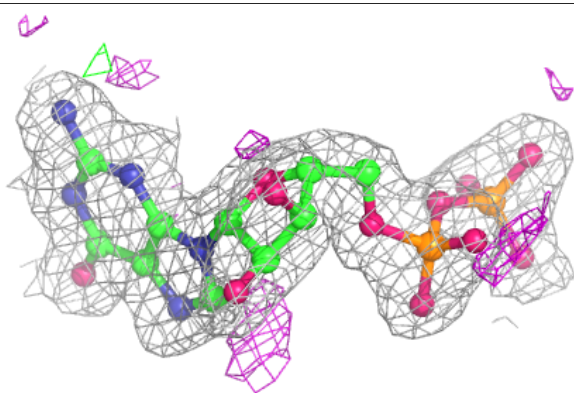
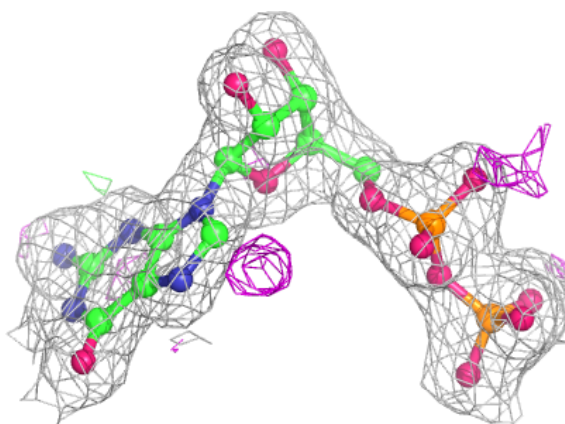
Electron density around MG A 502:

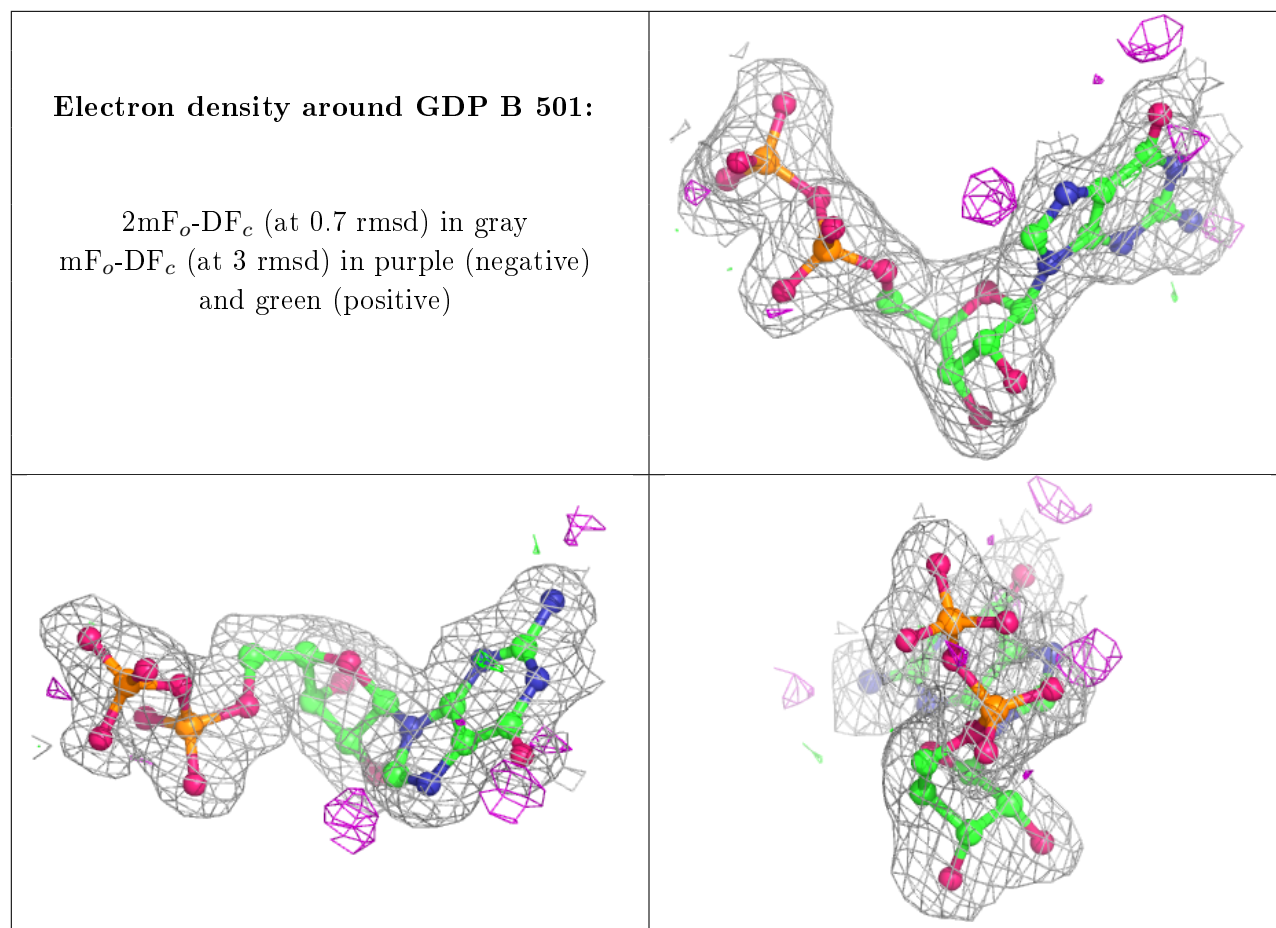
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around GDP A 501:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.