



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:59 PM UTC

PDB ID : 2G77 / pdb_00002g77
Title : Crystal Structure of Gyp1 TBC domain in complex with Rab33 GTPase bound to GDP and AlF3
Authors : Pan, X.; Eathiraj, S.; Munson, M.; Lambright, D.G.
Deposited on : 2006-02-27
Resolution : 2.26 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 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

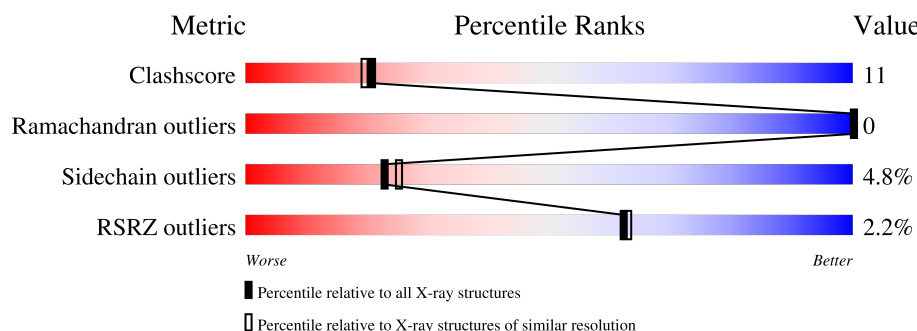
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.26 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))
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	410	
2	B	198	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4590 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 GTPase-activating protein GYP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	0	0
			2767	1786	466	504	11			

There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	228	MET	-	cloning artifact	UNP Q08484
A	229	GLY	-	cloning artifact	UNP Q08484
A	230	HIS	-	cloning artifact	UNP Q08484
A	231	HIS	-	cloning artifact	UNP Q08484
A	232	HIS	-	cloning artifact	UNP Q08484
A	233	HIS	-	cloning artifact	UNP Q08484
A	234	HIS	-	cloning artifact	UNP Q08484
A	235	HIS	-	cloning artifact	UNP Q08484
A	236	GLY	-	cloning artifact	UNP Q08484
A	237	SER	-	cloning artifact	UNP Q08484
A	238	LEU	-	cloning artifact	UNP Q08484
A	239	VAL	-	cloning artifact	UNP Q08484
A	240	PRO	-	cloning artifact	UNP Q08484
A	241	ARG	-	cloning artifact	UNP Q08484
A	242	GLY	-	cloning artifact	UNP Q08484
A	243	SER	-	cloning artifact	UNP Q08484
A	405	LYS	GLU	engineered mutation	UNP Q08484

- Molecule 2 is a protein called Ras-related protein Rab-33B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	173	Total	C	N	O	S	Se	0	0
			1374	869	248	249	4	4		

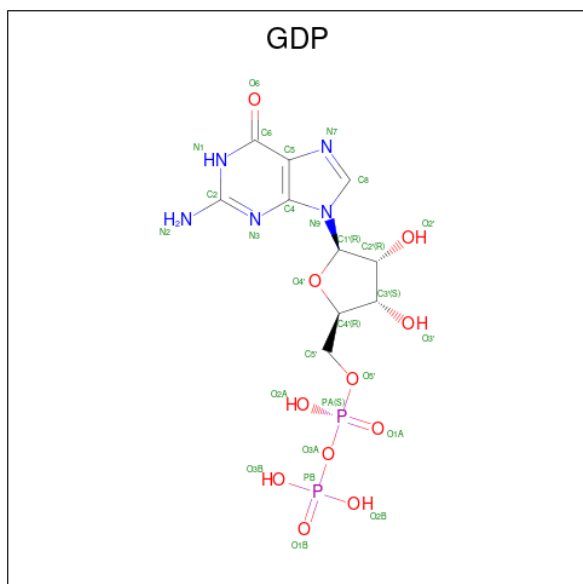
There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	5	MET	-	cloning artifact	UNP O35963
B	6	GLY	-	cloning artifact	UNP O35963
B	7	HIS	-	cloning artifact	UNP O35963
B	8	HIS	-	cloning artifact	UNP O35963
B	9	HIS	-	cloning artifact	UNP O35963
B	10	HIS	-	cloning artifact	UNP O35963
B	11	HIS	-	cloning artifact	UNP O35963
B	12	HIS	-	cloning artifact	UNP O35963
B	13	GLY	-	cloning artifact	UNP O35963
B	116	MSE	MET	modified residue	UNP O35963
B	119	MSE	MET	modified residue	UNP O35963
B	172	MSE	MET	modified residue	UNP O35963
B	193	MSE	MET	modified residue	UNP O35963

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

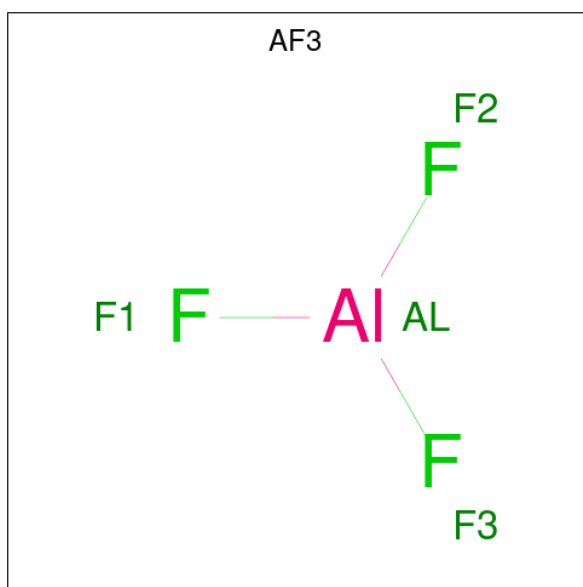
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mg 1 1	0	0

- Molecule 4 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



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

- Molecule 5 is ALUMINUM FLUORIDE (CCD ID: AF3) (formula: AlF₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total 4	Al 1	F 3	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	284	Total 284	O 284	0	0
6	B	132	Total 132	O 132	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.20Å 94.53Å 102.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.26 8.00 – 2.26	Depositor EDS
% Data completeness (in resolution range)	98.7 (8.00-2.26) 96.2 (8.00-2.26)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.10 (at 2.27Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.205 , 0.244 0.181 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	32.9	Xtriage
Anisotropy	0.957	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 66.0	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	0.96	EDS
Total number of atoms	4590	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.66% 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: MG, GDP, AF3

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	1.25	1/2840 (0.0%)	1.14	8/3857 (0.2%)
2	B	1.11	1/1400 (0.1%)	1.11	3/1887 (0.2%)
All	All	1.21	2/4240 (0.0%)	1.13	11/5744 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	150	CYS	C-N	-9.46	1.21	1.33
1	A	605	PRO	CA-C	5.36	1.57	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	588	ASP	N-CA-C	6.56	118.98	111.11
1	A	604	ASN	CA-C-N	-6.11	114.09	120.38
1	A	604	ASN	C-N-CA	-6.11	114.09	120.38
2	B	153	ARG	NE-CZ-NH1	-5.87	115.63	121.50
1	A	281	ILE	CA-C-N	-5.48	114.12	119.76
1	A	281	ILE	C-N-CA	-5.48	114.12	119.76
1	A	286	ARG	CA-C-N	-5.45	113.30	118.97
1	A	286	ARG	C-N-CA	-5.45	113.30	118.97
1	A	266	THR	N-CA-C	-5.40	105.50	112.68
2	B	151	ASP	CA-C-N	-5.20	114.22	122.23
2	B	151	ASP	C-N-CA	-5.20	114.22	122.23

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	2767	0	2704	38	0
2	B	1374	0	1341	54	0
3	B	1	0	0	0	0
4	B	28	0	12	0	0
5	B	4	0	0	0	0
6	A	284	0	0	12	0
6	B	132	0	0	8	0
All	All	4590	0	4057	91	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 11.

All (91) 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)
2:B:136:LEU:HD21	2:B:138:ALA:O	1.45	1.17
2:B:150:CYS:HB2	6:B:614:HOH:O	1.50	1.11
2:B:77:ILE:HB	2:B:200:LYS:HD2	1.43	0.99
1:A:366:LEU:HB2	6:A:920:HOH:O	1.64	0.97
2:B:35:LYS:H	2:B:108:HIS:HD2	1.12	0.96
2:B:77:ILE:O	2:B:200:LYS:HE3	1.64	0.96
1:A:284:ILE:HD11	1:A:398:PRO:HB3	1.47	0.93
1:A:325:ASP:HB2	6:A:705:HOH:O	1.68	0.92
1:A:616:GLU:HG3	6:A:827:HOH:O	1.69	0.91
2:B:119:MSE:CE	2:B:122:PHE:HD2	1.85	0.89
2:B:119:MSE:HE3	2:B:122:PHE:HD2	1.36	0.88
2:B:46:LYS:O	2:B:50:THR:HG23	1.75	0.84
2:B:136:LEU:CD2	2:B:138:ALA:O	2.28	0.80
2:B:119:MSE:CE	2:B:122:PHE:CD2	2.65	0.79
2:B:119:MSE:HE2	2:B:123:HIS:NE2	1.99	0.78
2:B:35:LYS:H	2:B:108:HIS:CD2	2.01	0.78
2:B:136:LEU:HG	2:B:137:LEU:H	1.50	0.75
1:A:327:HIS:HE1	1:A:365:TYR:OH	1.69	0.75
1:A:482:ARG:HD3	1:A:490:ARG:HB2	1.70	0.74
1:A:275:GLN:HG3	6:A:918:HOH:O	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:93:GLU:OE2	2:B:97:LYS:NZ	2.24	0.70
2:B:119:MSE:HE3	2:B:122:PHE:CD2	2.22	0.70
2:B:119:MSE:HE1	2:B:122:PHE:CD2	2.27	0.70
1:A:285:HIS:HD2	6:A:856:HOH:O	1.75	0.70
2:B:193:MSE:HA	2:B:193:MSE:HE2	1.73	0.70
2:B:136:LEU:HG	2:B:137:LEU:N	2.08	0.69
2:B:136:LEU:CD2	2:B:138:ALA:H	2.07	0.67
1:A:482:ARG:HD3	1:A:490:ARG:CB	2.25	0.67
1:A:464:ASP:C	1:A:591:MET:HE1	2.21	0.65
2:B:77:ILE:CB	2:B:200:LYS:HD2	2.22	0.65
2:B:101:GLN:OE1	2:B:102:HIS:HD2	1.79	0.65
2:B:131:GLU:HG2	6:B:635:HOH:O	1.98	0.63
2:B:53:PHE:CD1	2:B:192:PHE:HB2	2.35	0.62
2:B:194:THR:HG23	6:B:572:HOH:O	2.00	0.61
1:A:271:GLN:HE22	1:A:274:ARG:HH11	1.51	0.59
2:B:191:ILE:O	2:B:194:THR:HG22	2.03	0.59
2:B:96:ARG:HG3	2:B:128:TRP:CH2	2.38	0.58
1:A:572:ASN:ND2	6:A:888:HOH:O	2.36	0.58
2:B:116:MSE:SE	2:B:150:CYS:HB3	2.56	0.56
1:A:284:ILE:HD11	1:A:398:PRO:CB	2.29	0.56
2:B:150:CYS:CB	6:B:614:HOH:O	2.28	0.55
2:B:36:ILE:HD13	2:B:195:LEU:HD13	1.88	0.55
2:B:117:THR:HG21	2:B:152:LEU:HB2	1.89	0.55
2:B:36:ILE:HB	2:B:86:LEU:HD23	1.89	0.55
1:A:464:ASP:HB3	1:A:591:MET:HE1	1.89	0.53
2:B:96:ARG:O	2:B:100:VAL:HG12	2.08	0.53
2:B:192:PHE:CD2	2:B:193:MSE:HE3	2.43	0.53
2:B:131:GLU:HB3	6:B:635:HOH:O	2.09	0.53
1:A:263:LYS:HE2	6:A:882:HOH:O	2.09	0.52
2:B:104:TYR:HE2	2:B:132:CYS:SG	2.33	0.51
1:A:439:ASN:HD22	1:A:446:GLY:HA3	1.76	0.51
1:A:246:GLN:CB	1:A:249:ASN:HB3	2.40	0.50
1:A:339:ILE:HD12	2:B:43:ASN:HB2	1.94	0.50
1:A:284:ILE:CD1	1:A:398:PRO:HB3	2.31	0.50
2:B:61:ARG:HD2	6:B:531:HOH:O	2.12	0.49
2:B:136:LEU:HD21	2:B:138:ALA:H	1.77	0.49
1:A:327:HIS:CE1	1:A:365:TYR:OH	2.58	0.49
1:A:275:GLN:CG	6:A:918:HOH:O	2.56	0.47
2:B:136:LEU:HD21	2:B:138:ALA:C	2.31	0.47
2:B:119:MSE:HE2	2:B:123:HIS:CD2	2.51	0.46
2:B:102:HIS:HB2	6:B:622:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:39:ILE:HD11	2:B:112:PHE:CE2	2.51	0.46
1:A:449:ARG:NH1	6:A:898:HOH:O	2.48	0.45
1:A:582:PHE:HA	1:A:618:LEU:HD11	1.98	0.45
1:A:287:PRO:O	1:A:291:LYS:HG3	2.16	0.45
1:A:321:HIS:HD2	6:A:666:HOH:O	1.98	0.45
2:B:50:THR:HG21	2:B:88:ASP:HB2	1.98	0.45
1:A:474:VAL:HG21	1:A:598:THR:HG21	1.99	0.45
2:B:192:PHE:CD2	2:B:193:MSE:CE	3.00	0.45
1:A:492:PHE:CE2	1:A:500:MET:HE1	2.51	0.44
1:A:345:ASN:N	1:A:346:PRO:HD3	2.33	0.44
1:A:464:ASP:HB3	1:A:591:MET:CE	2.48	0.43
2:B:99:MET:O	2:B:100:VAL:C	2.61	0.43
2:B:129:ILE:HD11	2:B:172:MSE:HE1	2.01	0.43
1:A:482:ARG:NH2	6:A:853:HOH:O	2.40	0.43
2:B:136:LEU:CG	2:B:137:LEU:N	2.79	0.42
1:A:330:ASP:OD1	1:A:332:PRO:HD2	2.20	0.42
2:B:151:ASP:OD1	2:B:178:SER:OG	2.33	0.41
1:A:381:ASN:O	1:A:384:VAL:HG12	2.20	0.41
2:B:77:ILE:CG2	2:B:200:LYS:HD2	2.50	0.41
1:A:582:PHE:CA	1:A:618:LEU:HD11	2.51	0.41
2:B:96:ARG:HD2	2:B:131:GLU:OE1	2.21	0.41
1:A:507:GLU:OE2	1:A:569:SER:HB3	2.21	0.41
2:B:54:CYS:SG	2:B:86:LEU:HD12	2.60	0.41
2:B:117:THR:CG2	2:B:152:LEU:HB2	2.50	0.41
2:B:153:ARG:HB2	6:B:608:HOH:O	2.20	0.41
1:A:281:ILE:O	1:A:282:PRO:C	2.62	0.41
2:B:37:ILE:HG22	2:B:87:TRP:HB2	2.02	0.41
1:A:271:GLN:HE22	1:A:274:ARG:NH1	2.17	0.40
1:A:283:LYS:HE3	6:A:813:HOH:O	2.21	0.40
1:A:489:MET:HE2	1:A:489:MET:HB3	1.96	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	327/410 (80%)	319 (98%)	8 (2%)	0	100	100
2	B	171/198 (86%)	162 (95%)	9 (5%)	0	100	100
All	All	498/608 (82%)	481 (97%)	17 (3%)	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	308/386 (80%)	297 (96%)	11 (4%)	31	39
2	B	146/166 (88%)	135 (92%)	11 (8%)	12	11
All	All	454/552 (82%)	432 (95%)	22 (5%)	23	25

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	251	ILE
1	A	263	LYS
1	A	284	ILE
1	A	302	LYS
1	A	319	LEU
1	A	409	PRO
1	A	421	LEU
1	A	457	LEU
1	A	482	ARG
1	A	500	MET
1	A	588	ASP
2	B	49	LEU
2	B	50	THR
2	B	71	ARG
2	B	75	VAL

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Mol	Chain	Res	Type
2	B	96	ARG
2	B	107	VAL
2	B	134	GLN
2	B	136	LEU
2	B	143	ARG
2	B	194	THR
2	B	195	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	253	GLN
1	A	270	GLN
1	A	271	GLN
1	A	275	GLN
1	A	321	HIS
1	A	327	HIS
1	A	400	GLN
1	A	439	ASN
1	A	450	GLN
1	A	453	ASN
1	A	467	ASN
1	A	471	ASN
1	A	626	GLN
2	B	102	HIS
2	B	108	HIS
2	B	118	ASN
2	B	135	HIS
2	B	187	HIS
2	B	197	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is 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	$\# Z > 2$
5	AF3	B	502	3,4,6	0,3,3	-	-	-		
4	GDP	B	501	3,5	29,30,30	1.34	4 (13%)	45,47,47	1.95	7 (15%)

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	GDP	B	501	3,5	-	3/16/32/32	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	501	GDP	C6-N1	-3.20	1.32	1.38
4	B	501	GDP	C8-N9	-2.47	1.32	1.37
4	B	501	GDP	C5-C4	2.33	1.45	1.38
4	B	501	GDP	PB-O3B	-2.03	1.47	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	501	GDP	C5-C4-N3	-6.35	118.29	128.39
4	B	501	GDP	C2-N3-C4	4.90	120.75	112.30
4	B	501	GDP	N9-C4-N3	4.25	134.45	125.95
4	B	501	GDP	C6-C5-N7	3.91	137.40	130.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	501	GDP	O2A-PA-O3A	3.86	117.71	107.27
4	B	501	GDP	C4-C5-N7	-3.71	104.78	110.67
4	B	501	GDP	C8-N7-C5	2.69	109.05	104.26

There are no chirality outliers.

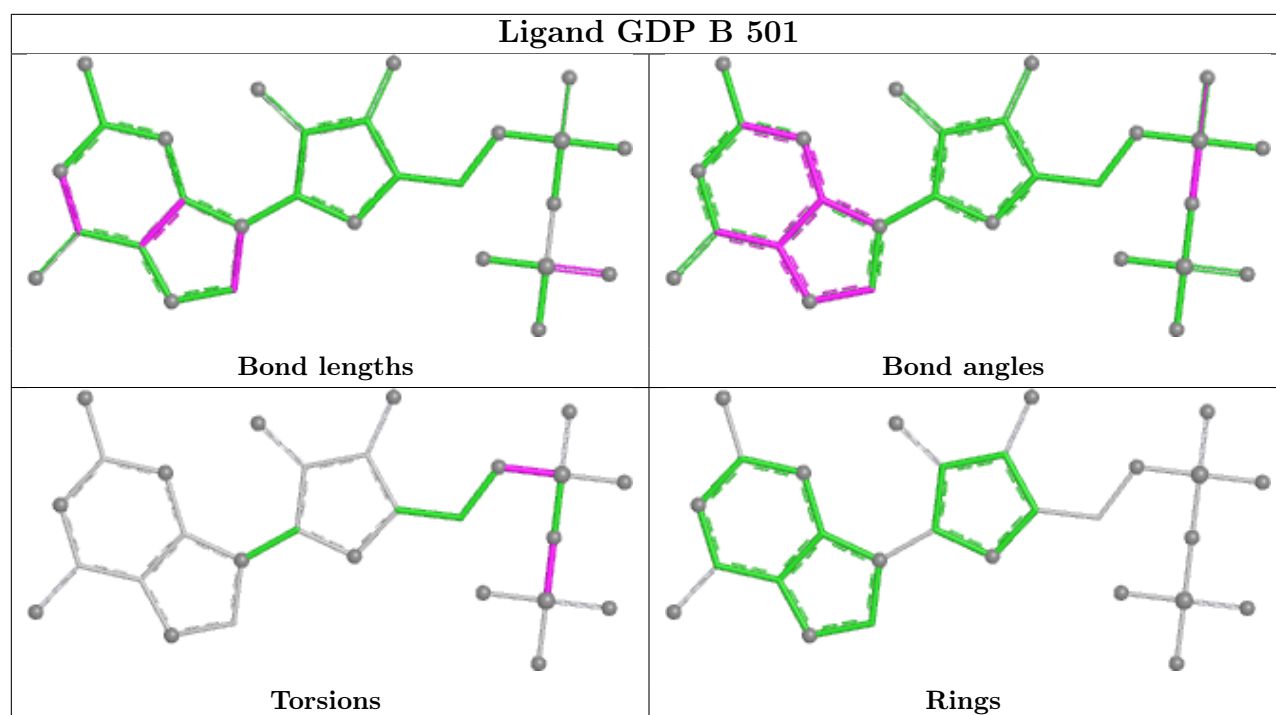
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	501	GDP	PA-O3A-PB-O3B
4	B	501	GDP	C5'-O5'-PA-O1A
4	B	501	GDP	PA-O3A-PB-O2B

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	331/410 (80%)	-0.62	7 (2%) 63 64	7, 34, 53, 79	5 (1%)
2	B	169/198 (85%)	-0.24	4 (2%) 59 61	26, 45, 81, 92	0
All	All	500/608 (82%)	-0.49	11 (2%) 62 63	7, 37, 73, 92	5 (1%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	633	THR	6.0
1	A	627	SER	3.3
2	B	137	LEU	3.1
1	A	615	ILE	3.0
1	A	620	SER	3.0
1	A	473	HIS	2.9
1	A	632	ALA	2.8
2	B	138	ALA	2.5
2	B	140	ASP	2.5
1	A	569	SER	2.3
2	B	100	VAL	2.1

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](#)

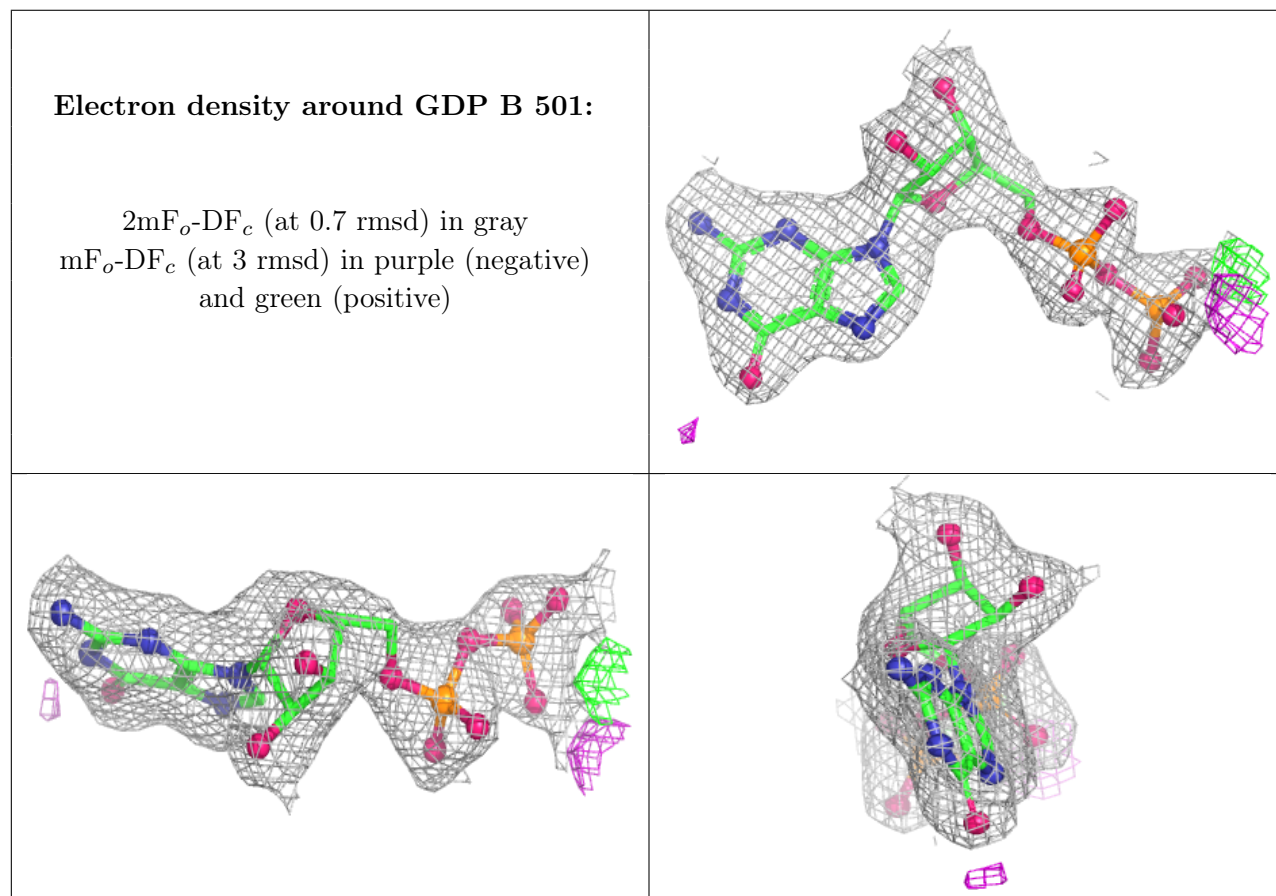
There are no oligosaccharides in this entry.

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
3	MG	B	503	1/1	0.80	0.18	13,13,13,13	0
5	AF3	B	502	4/4	0.86	0.10	33,33,41,44	0
4	GDP	B	501	28/28	0.98	0.04	22,30,33,34	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.