Full wwPDB X-ray Structure Validation Report

Nov 15, 2023 – 01:42 PM JST

PDB ID	:	6J9V
Title	:	Crystal structure of Trypanosoma brucei gambiense glycerol kinase complex with ADP.
Authors	:	Balogun, E.O.; Chishima, T.; Ichinose, M.; Inaoka, D.K.; Kido, Y.; Ibrahim, B.; de Koning, H.; McKerrow, J.H.; Watanabe, Y.; Nozaki, T.; Michels, P.A.M.; Harada, S.; Kita, K.; Shiba, T.
Deposited on	:	2019-01-24
Resolution	:	2.40 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 (1)) were used in the production of this report:

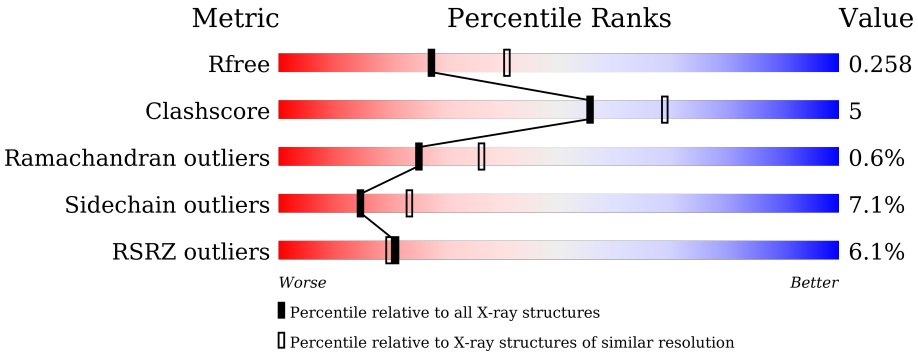
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 <=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	518	3% 84% 14% ··
1	B	518	9% 81% 16% ··

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 7993 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.

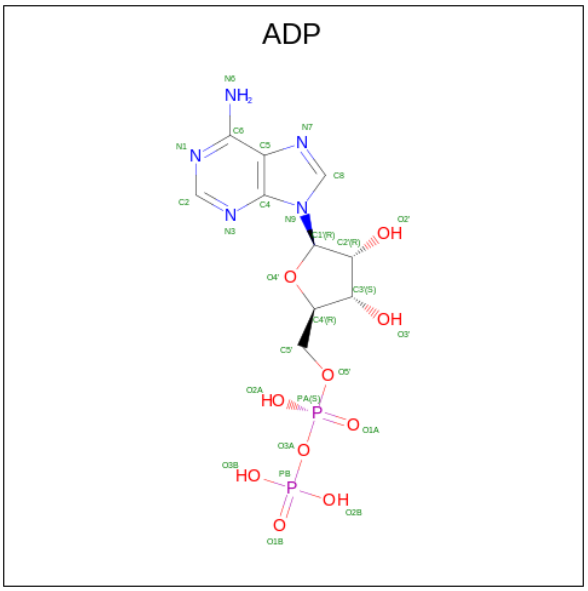
- Molecule 1 is a protein called Glycerol kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	513	Total	C	N	O	S	0	0	0
			3957	2499	694	731	33			
1	B	513	Total	C	N	O	S	0	0	0
			3957	2499	694	731	33			

There are 12 discrepancies between the modelled and reference sequences:

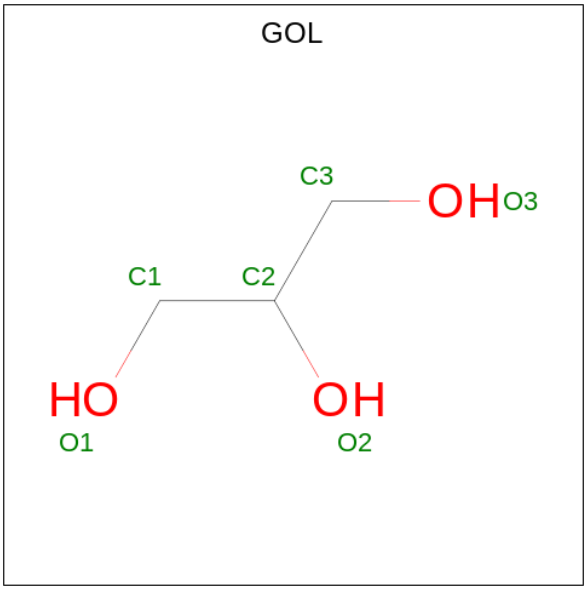
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP D3KVM3
A	-4	ILE	-	expression tag	UNP D3KVM3
A	-3	ASP	-	expression tag	UNP D3KVM3
A	-2	PRO	-	expression tag	UNP D3KVM3
A	-1	PHE	-	expression tag	UNP D3KVM3
A	0	THR	-	expression tag	UNP D3KVM3
B	-5	GLY	-	expression tag	UNP D3KVM3
B	-4	ILE	-	expression tag	UNP D3KVM3
B	-3	ASP	-	expression tag	UNP D3KVM3
B	-2	PRO	-	expression tag	UNP D3KVM3
B	-1	PHE	-	expression tag	UNP D3KVM3
B	0	THR	-	expression tag	UNP D3KVM3

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

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

- Molecule 4 is water.

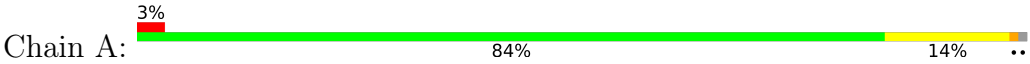
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	6	Total O 6 6	0	0
4	B	1	Total O 1 1	0	0

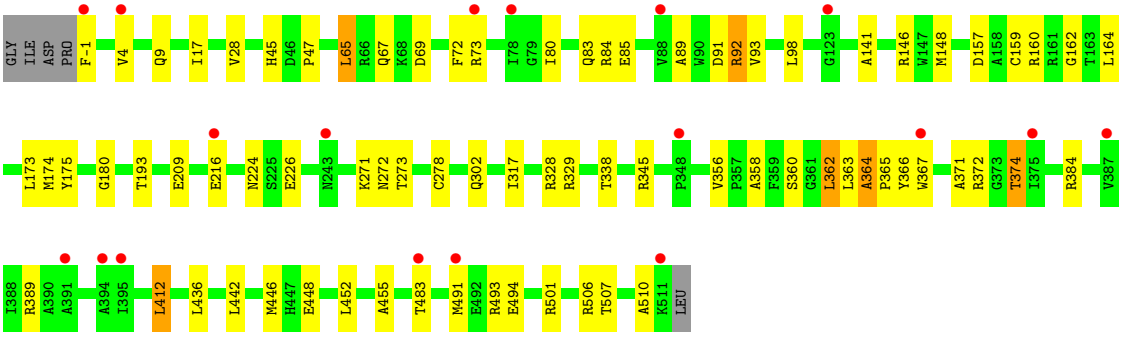
3 Residue-property plots (i)

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: Glycerol kinase

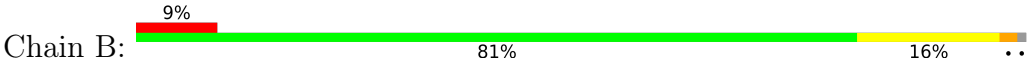
Chain A:

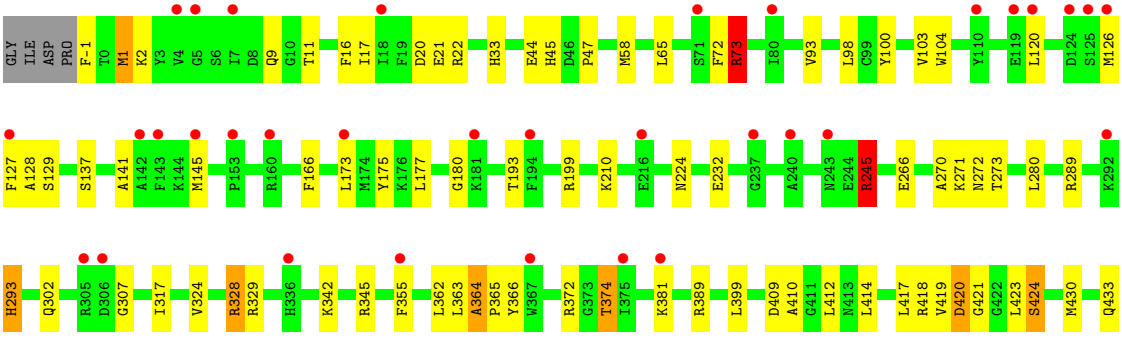




- Molecule 1: Glycerol kinase

Chain B:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	62.42Å 120.95Å 154.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 19.99 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.6 (20.00-2.40) 99.9 (19.99-2.40)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.46 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.193 , 0.254 0.201 , 0.258	Depositor DCC
R_{free} test set	2340 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	70.4	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 58.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7993	wwPDB-VP
Average B, all atoms (Å^2)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: ADP, GOL

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	0.71	0/4039	0.88	5/5465 (0.1%)
1	B	0.67	0/4039	0.88	5/5465 (0.1%)
All	All	0.69	0/8078	0.88	10/10930 (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	B	0	1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	245	ARG	NE-CZ-NH1	7.51	124.06	120.30
1	A	389	ARG	NE-CZ-NH1	6.73	123.67	120.30
1	A	506	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	B	389	ARG	NE-CZ-NH2	-5.53	117.54	120.30
1	A	146	ARG	NE-CZ-NH2	-5.45	117.58	120.30
1	A	412	LEU	CA-CB-CG	5.36	127.63	115.30
1	B	345	ARG	NE-CZ-NH2	5.26	122.93	120.30
1	A	389	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	B	307	GLY	N-CA-C	-5.08	100.40	113.10
1	B	389	ARG	NE-CZ-NH1	5.04	122.82	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	363	LEU	Peptide

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	3957	0	3970	37	0
1	B	3957	0	3970	44	0
2	A	27	0	12	0	0
2	B	27	0	12	0	0
3	A	6	0	8	2	0
3	B	12	0	16	0	0
4	A	6	0	0	0	0
4	B	1	0	0	0	0
All	All	7993	0	7988	81	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 5.

All (81) 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:22:ARG:HB3	1:B:478:GLU:HG2	1.73	0.71
1:A:364:ALA:O	1:A:367:TRP:NE1	2.30	0.65
1:A:92:ARG:NH2	1:A:162:GLY:O	2.33	0.62
1:A:364:ALA:HB1	1:A:365:PRO:CD	2.30	0.62
1:B:364:ALA:HB1	1:B:365:PRO:CD	2.30	0.61
1:B:1:MET:HE1	1:B:21:GLU:H	1.66	0.60
1:B:1:MET:CE	1:B:2:LYS:O	2.50	0.59
1:A:224:ASN:HD22	1:A:302:GLN:H	1.50	0.59
1:B:374:THR:HG21	1:B:507:THR:HA	1.84	0.57
1:A:272:ASN:HD21	1:A:278:CYS:HB3	1.70	0.57
1:A:374:THR:HG21	1:A:507:THR:HA	1.87	0.56
1:A:446:MET:CE	1:A:452:LEU:HD22	2.36	0.56
1:B:1:MET:HE1	1:B:21:GLU:N	2.21	0.56
1:A:89:ALA:HB2	1:A:148:MET:HE1	1.88	0.55
1:B:328:ARG:HG2	1:B:329:ARG:N	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:421:GLY:O	1:B:424:SER:HB2	2.06	0.54
1:A:17:ILE:HD13	1:A:448:GLU:HG2	1.89	0.54
1:A:358:ALA:HB2	1:A:362:LEU:HD13	1.89	0.53
1:A:65:LEU:HD13	1:A:72:PHE:CG	2.44	0.52
1:B:364:ALA:CB	1:B:365:PRO:CD	2.87	0.52
1:B:1:MET:HE3	1:B:2:LYS:N	2.24	0.52
1:B:293:HIS:CD2	1:B:409:ASP:OD1	2.62	0.51
1:A:374:THR:HG23	1:A:507:THR:HG22	1.93	0.51
1:B:224:ASN:HD22	1:B:302:GLN:H	1.58	0.51
1:A:141:ALA:HB3	1:A:193:THR:HA	1.91	0.50
1:B:419:VAL:HG21	1:B:430:MET:SD	2.50	0.50
1:A:83:GLN:O	1:A:83:GLN:HG3	2.10	0.50
1:A:364:ALA:CB	1:A:365:PRO:CD	2.90	0.50
1:B:72:PHE:O	1:B:73:ARG:HB2	2.11	0.50
1:B:418:ARG:HG3	1:B:442:LEU:HB2	1.94	0.50
1:A:374:THR:CG2	1:A:507:THR:HG22	2.42	0.49
1:A:436:LEU:O	1:A:493:ARG:NE	2.39	0.49
1:A:84:ARG:HH21	3:A:702:GOL:H31	1.78	0.48
1:A:4:VAL:HG11	1:A:455:ALA:O	2.14	0.48
1:B:1:MET:HE3	1:B:2:LYS:H	1.79	0.48
1:A:273:THR:O	1:A:278:CYS:HA	2.14	0.47
1:B:45:HIS:O	1:B:47:PRO:HD3	2.15	0.46
1:B:272:ASN:HB2	1:B:280:LEU:HD12	1.97	0.46
1:B:16:PHE:CD2	1:B:58:MET:HA	2.51	0.46
1:B:145:MET:HG2	1:B:166:PHE:CG	2.51	0.46
1:B:175:TYR:CE1	1:B:180:GLY:HA2	2.50	0.46
1:B:72:PHE:O	1:B:73:ARG:CB	2.64	0.46
1:B:374:THR:HG23	1:B:507:THR:HG22	1.98	0.46
1:B:374:THR:CG2	1:B:507:THR:HG22	2.46	0.45
1:B:141:ALA:HB3	1:B:193:THR:HA	1.97	0.45
1:B:372:ARG:O	1:B:374:THR:HG22	2.17	0.45
1:A:17:ILE:HD13	1:A:448:GLU:CG	2.47	0.44
1:A:17:ILE:HG12	1:A:28:VAL:HG23	1.99	0.44
1:B:399:LEU:HD12	1:B:433:GLN:OE1	2.18	0.44
1:A:89:ALA:HB2	1:A:148:MET:CE	2.47	0.44
1:A:364:ALA:HB1	1:A:365:PRO:HD3	2.00	0.43
1:A:345:ARG:HH21	1:A:345:ARG:HG3	1.83	0.43
1:A:363:LEU:HD22	1:A:367:TRP:CZ3	2.53	0.43
1:B:232:GLU:OE1	1:B:245:ARG:NH2	2.51	0.43
1:A:80:ILE:HD12	1:A:174:MET:HG3	2.00	0.43
1:A:356:VAL:HG11	1:A:507:THR:HG21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:PHE:CG	1:B:128:ALA:N	2.87	0.42
1:B:410:ALA:HB1	1:B:412:LEU:HD12	2.01	0.42
1:B:270:ALA:HB3	1:B:414:LEU:HD11	2.02	0.42
1:B:273:THR:HG23	1:B:420:ASP:OD1	2.19	0.42
1:A:175:TYR:CE1	1:A:180:GLY:HA2	2.55	0.42
1:B:417:LEU:O	1:B:441:ILE:HA	2.19	0.42
1:A:45:HIS:O	1:A:47:PRO:HD3	2.18	0.42
1:B:293:HIS:CD2	1:B:293:HIS:N	2.88	0.42
1:B:364:ALA:HB1	1:B:365:PRO:HD2	2.01	0.42
1:A:91:ASP:OD1	1:A:93:VAL:HG22	2.19	0.42
1:A:84:ARG:HH21	3:A:702:GOL:C3	2.32	0.41
1:B:11:THR:O	1:B:33:HIS:NE2	2.52	0.41
1:A:84:ARG:O	1:A:85:GLU:HB2	2.20	0.41
1:B:103:VAL:O	1:B:104:TRP:C	2.58	0.41
1:A:159:CYS:HB2	1:A:164:LEU:HD22	2.02	0.41
1:B:120:LEU:HD13	1:B:210:LYS:HG2	2.03	0.41
1:B:324:VAL:HG21	1:B:423:LEU:HD21	2.03	0.41
1:A:65:LEU:HD22	1:A:69:ASP:HB3	2.02	0.41
1:A:89:ALA:CB	1:A:148:MET:HE1	2.51	0.40
1:B:17:ILE:HD13	1:B:451:ALA:CB	2.51	0.40
1:B:289:ARG:HD3	1:B:410:ALA:HA	2.03	0.40
1:B:1:MET:CE	1:B:20:ASP:HB2	2.51	0.40
1:A:371:ALA:O	1:A:372:ARG:NH1	2.50	0.40
1:B:44:GLU:CG	1:B:100:TYR:HB3	2.51	0.40
1:B:364:ALA:HB1	1:B:365:PRO:HD3	2.02	0.40

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	511/518 (99%)	488 (96%)	20 (4%)	3 (1%)	25 36

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	511/518 (99%)	475 (93%)	33 (6%)	3 (1%)	25 36
All	All	1022/1036 (99%)	963 (94%)	53 (5%)	6 (1%)	25 36

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	364	ALA
1	B	364	ALA
1	B	448	GLU
1	B	73	ARG
1	A	226	GLU
1	A	510	ALA

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	421/425 (99%)	393 (93%)	28 (7%)	16 26
1	B	421/425 (99%)	389 (92%)	32 (8%)	13 20
All	All	842/850 (99%)	782 (93%)	60 (7%)	14 23

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-1	PHE
1	A	9	GLN
1	A	65	LEU
1	A	67	GLN
1	A	73	ARG
1	A	92	ARG
1	A	98	LEU
1	A	157	ASP
1	A	160	ARG
1	A	173	LEU

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Mol	Chain	Res	Type
1	A	209	GLU
1	A	216	GLU
1	A	271	LYS
1	A	317	ILE
1	A	328	ARG
1	A	329	ARG
1	A	338	THR
1	A	360	SER
1	A	362	LEU
1	A	366	TYR
1	A	374	THR
1	A	384	ARG
1	A	412	LEU
1	A	442	LEU
1	A	483	THR
1	A	491	MET
1	A	494	GLU
1	A	501	ARG
1	B	-1	PHE
1	B	1	MET
1	B	9	GLN
1	B	65	LEU
1	B	73	ARG
1	B	93	VAL
1	B	98	LEU
1	B	126	MET
1	B	129	SER
1	B	137	SER
1	B	173	LEU
1	B	177	LEU
1	B	199	ARG
1	B	245	ARG
1	B	266	GLU
1	B	271	LYS
1	B	293	HIS
1	B	317	ILE
1	B	328	ARG
1	B	342	LYS
1	B	355	PHE
1	B	362	LEU
1	B	366	TYR
1	B	374	THR

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Mol	Chain	Res	Type
1	B	381	LYS
1	B	420	ASP
1	B	424	SER
1	B	442	LEU
1	B	480	SER
1	B	485	SER
1	B	491	MET
1	B	511	LYS

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

Mol	Chain	Res	Type
1	A	23	GLN
1	A	30	GLN
1	A	151	ASN
1	A	224	ASN
1	A	272	ASN
1	A	302	GLN
1	A	351	GLN
1	B	23	GLN
1	B	224	ASN
1	B	243	ASN
1	B	272	ASN
1	B	302	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	GOL	B	703	-	5,5,5	0.51	0	5,5,5	1.03	0
3	GOL	A	702	-	5,5,5	0.73	0	5,5,5	1.56	1 (20%)
2	ADP	B	701	-	24,29,29	1.09	1 (4%)	29,45,45	1.44	5 (17%)
2	ADP	A	701	-	24,29,29	1.00	1 (4%)	29,45,45	1.63	6 (20%)
3	GOL	B	702	-	5,5,5	0.68	0	5,5,5	0.86	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
3	GOL	B	703	-	-	2/4/4/4	-
3	GOL	A	702	-	-	4/4/4/4	-
2	ADP	B	701	-	-	0/12/32/32	0/3/3/3
2	ADP	A	701	-	-	6/12/32/32	0/3/3/3
3	GOL	B	702	-	-	2/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	ADP	C5-C4	2.69	1.48	1.40
2	B	701	ADP	C5-C4	2.29	1.47	1.40

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	701	ADP	N3-C2-N1	-4.80	121.17	128.68
2	A	701	ADP	C3'-C2'-C1'	3.47	106.20	100.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	701	ADP	N3-C2-N1	-3.46	123.27	128.68
2	A	701	ADP	C2-N1-C6	2.98	123.86	118.75
2	B	701	ADP	PA-O3A-PB	-2.90	122.86	132.83
2	A	701	ADP	C1'-N9-C4	-2.58	122.11	126.64
3	A	702	GOL	O3-C3-C2	2.55	122.43	110.20
2	B	701	ADP	N6-C6-N1	2.37	123.49	118.57
2	B	701	ADP	C2-N1-C6	2.25	122.60	118.75
2	B	701	ADP	C1'-N9-C4	-2.14	122.87	126.64
2	A	701	ADP	O3B-PB-O2B	2.14	115.83	107.64
2	A	701	ADP	C4-C5-N7	-2.12	107.19	109.40

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	ADP	PA-O3A-PB-O2B
2	A	701	ADP	PA-O3A-PB-O3B
3	A	702	GOL	O1-C1-C2-C3
3	B	702	GOL	C1-C2-C3-O3
3	A	702	GOL	C1-C2-C3-O3
3	B	703	GOL	O1-C1-C2-C3
3	A	702	GOL	O1-C1-C2-O2
3	B	702	GOL	O2-C2-C3-O3
3	B	703	GOL	O1-C1-C2-O2
3	A	702	GOL	O2-C2-C3-O3
2	A	701	ADP	O4'-C4'-C5'-O5'
2	A	701	ADP	PB-O3A-PA-O1A
2	A	701	ADP	PA-O3A-PB-O1B
2	A	701	ADP	C3'-C4'-C5'-O5'

There are no ring outliers.

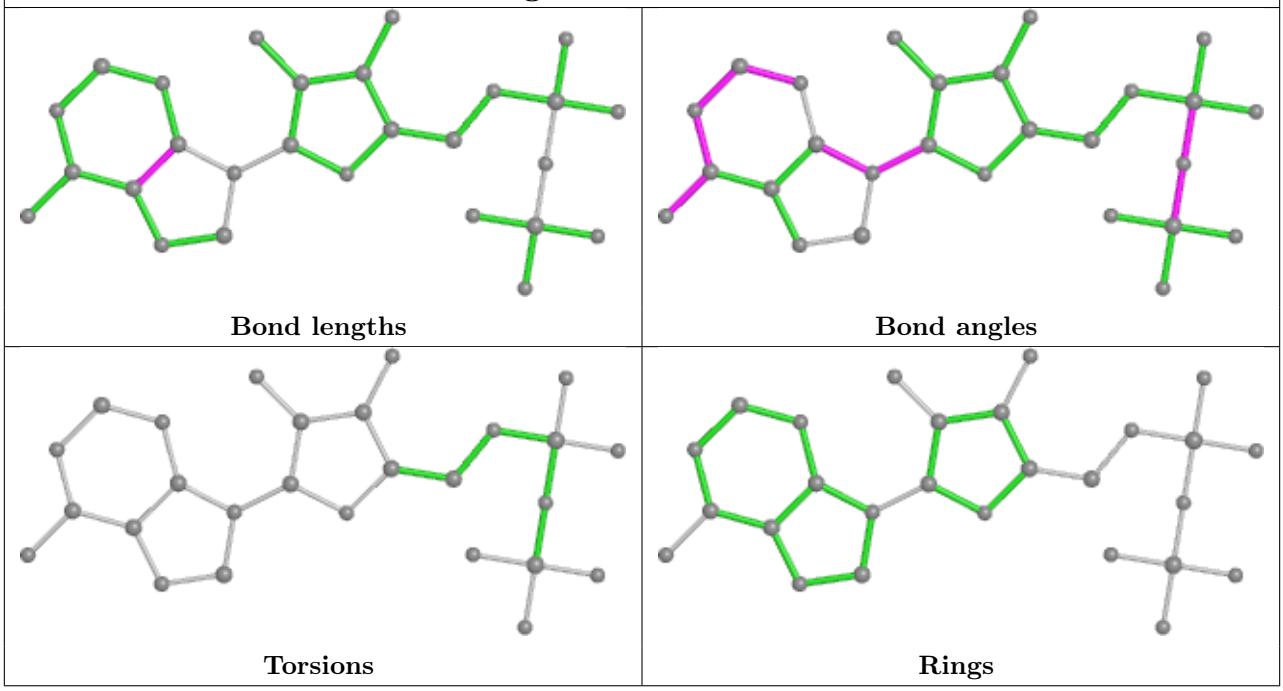
1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	702	GOL	2	0

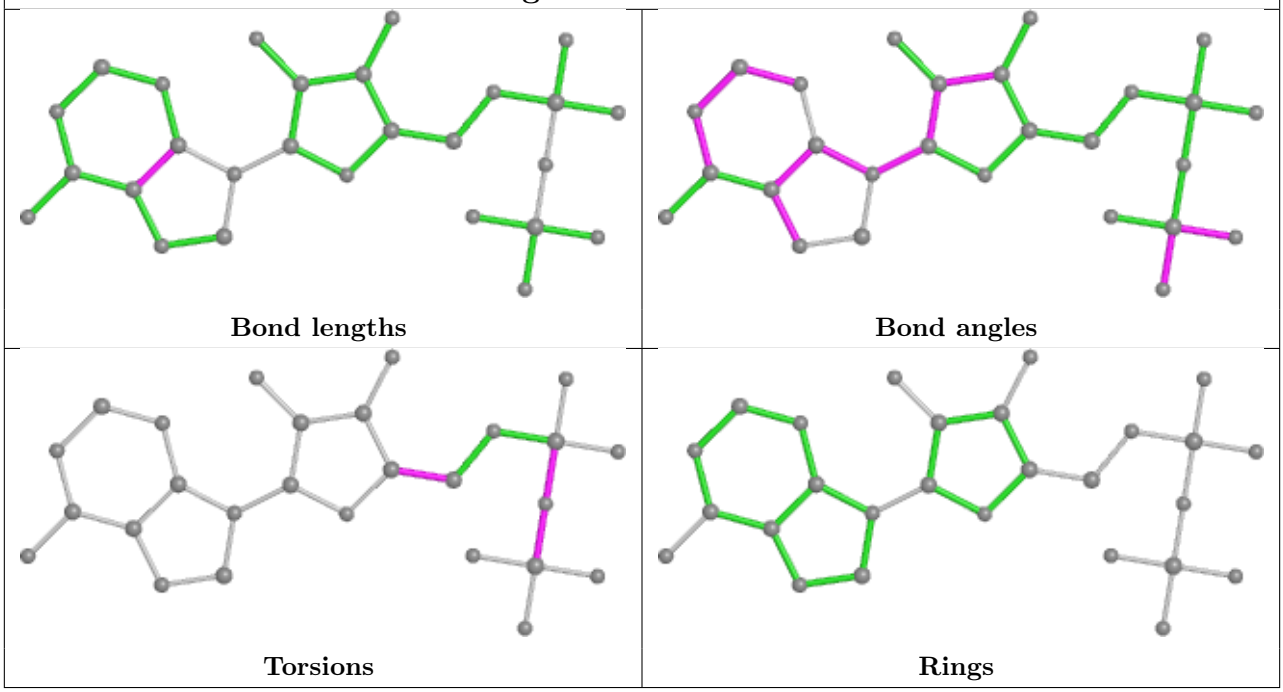
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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.

Ligand ADP B 701



Ligand ADP A 701



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, 95^{th} 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	513/518 (99%)	0.14	18 (3%) 44 43	48, 66, 90, 111	0
1	B	513/518 (99%)	0.35	45 (8%) 10 9	53, 75, 106, 128	0
All	All	1026/1036 (99%)	0.24	63 (6%) 21 20	48, 70, 102, 128	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	511	LYS	4.6
1	B	447	HIS	4.5
1	B	488	GLY	4.3
1	B	306	ASP	4.2
1	B	491	MET	4.2
1	B	490	ALA	3.8
1	B	120	LEU	3.7
1	A	348	PRO	3.6
1	B	119	GLU	3.5
1	B	292	LYS	3.5
1	A	88	VAL	3.4
1	B	367	TRP	3.2
1	A	491	MET	3.2
1	A	483	THR	3.0
1	B	160	ARG	2.9
1	B	145	MET	2.9
1	B	110	TYR	2.9
1	B	143	PHE	2.9
1	A	367	TRP	2.8
1	B	7	ILE	2.8
1	A	391	ALA	2.7
1	B	381	LYS	2.7
1	A	216	GLU	2.7
1	B	489	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	216	GLU	2.6
1	B	18	ILE	2.6
1	B	375	ILE	2.6
1	B	495	ALA	2.6
1	B	240	ALA	2.6
1	B	173	LEU	2.6
1	B	124	ASP	2.5
1	B	153	PRO	2.5
1	B	127	PHE	2.5
1	B	4	VAL	2.5
1	B	305	ARG	2.4
1	B	487	SER	2.4
1	A	375	ILE	2.4
1	B	445	SER	2.4
1	B	142	ALA	2.4
1	A	-1	PHE	2.4
1	A	243	ASN	2.4
1	B	126	MET	2.4
1	A	73	ARG	2.4
1	B	5	GLY	2.3
1	B	510	ALA	2.3
1	A	78	ILE	2.3
1	B	243	ASN	2.2
1	A	387	VAL	2.2
1	A	395	ILE	2.2
1	A	511	LYS	2.2
1	B	125	SER	2.2
1	A	394	ALA	2.2
1	B	477	ARG	2.2
1	B	336	HIS	2.2
1	B	80	ILE	2.2
1	B	237	GLY	2.2
1	A	4	VAL	2.1
1	B	469	GLU	2.1
1	B	194	PHE	2.1
1	B	355	PHE	2.1
1	B	181	LYS	2.0
1	A	123	GLY	2.0
1	B	71	SER	2.0

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 monosaccharides in this entry.

6.4 Ligands (i)

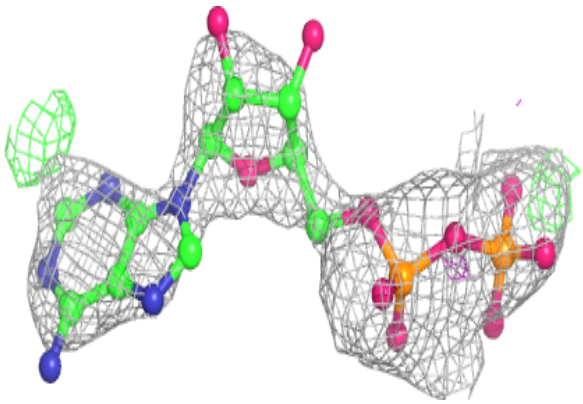
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

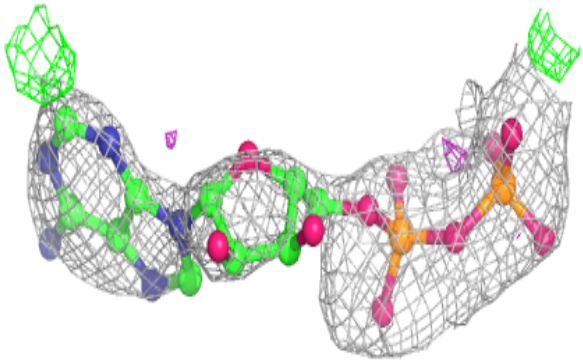
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GOL	A	702	6/6	0.78	0.24	65,69,74,81	0
2	ADP	A	701	27/27	0.80	0.33	79,97,106,111	27
3	GOL	B	703	6/6	0.81	0.14	65,76,90,93	0
3	GOL	B	702	6/6	0.85	0.28	63,67,71,75	0
2	ADP	B	701	27/27	0.93	0.13	55,66,115,132	0

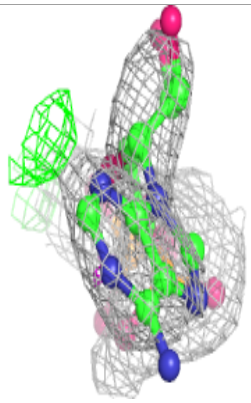
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP A 701:

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

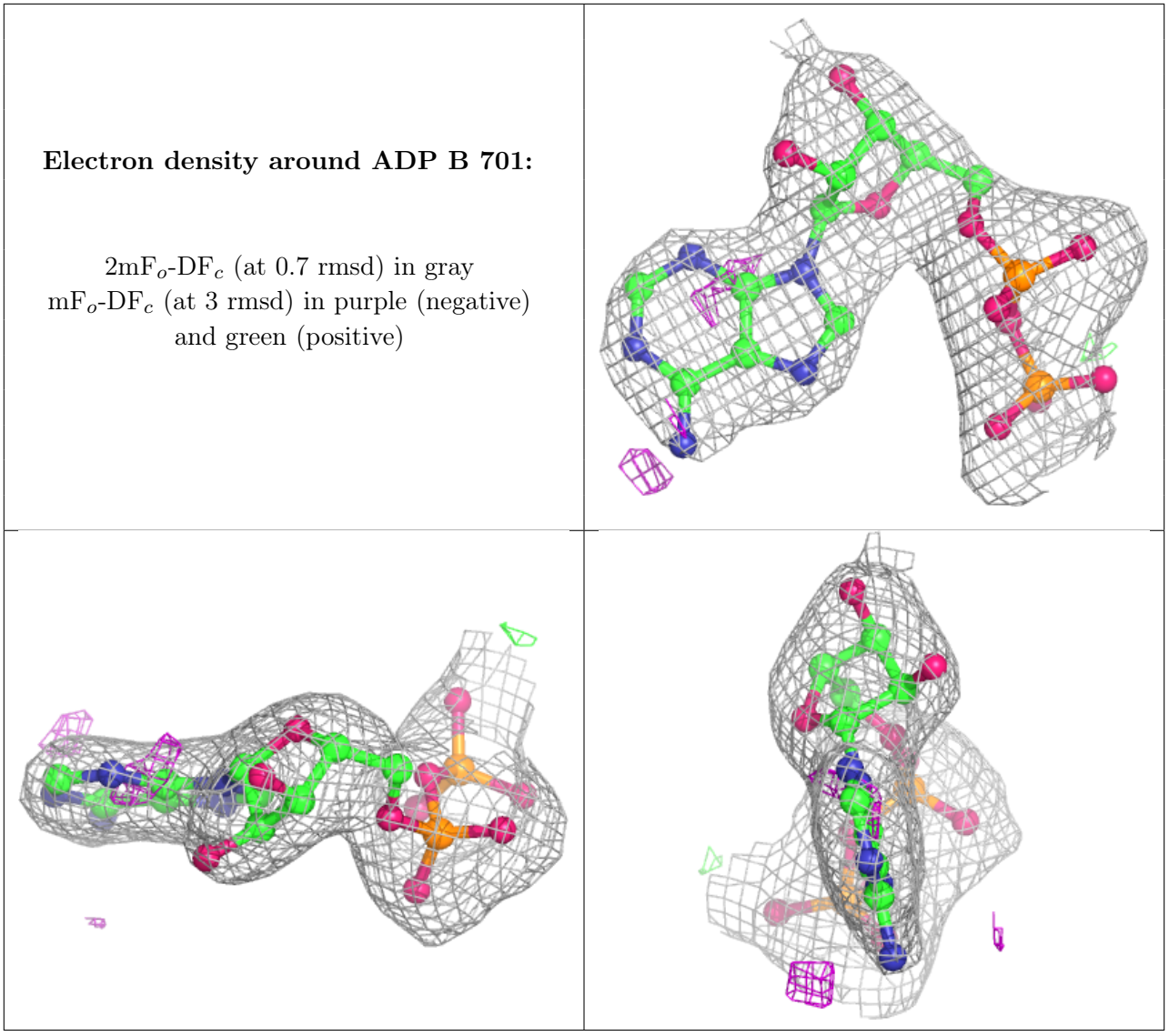






Electron density around ADP B 701:

$2mF_o$-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.