



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:17 AM UTC

PDB ID : 7RCM / pdb_00007rcm
Title : Crystal Structure of ADP-bound Galactokinase
Authors : Whitby, F.G.
Deposited on : 2021-07-07
Resolution : 2.10 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

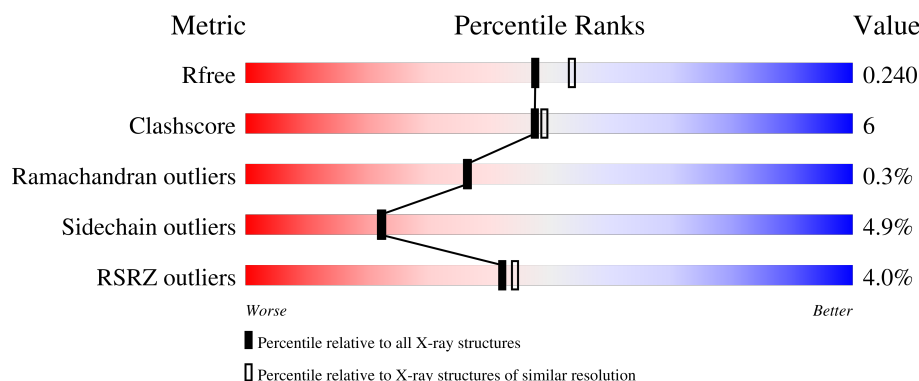
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	392	
1	B	392	

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 6287 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.

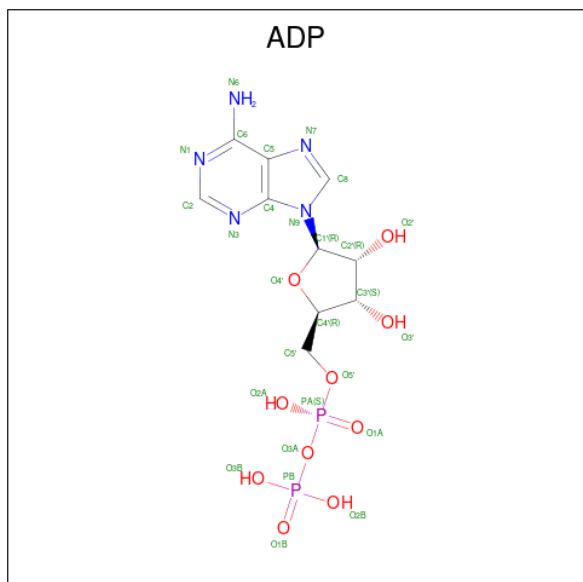
- Molecule 1 is a protein called Galactokinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	S	0	2	0
			2963	1849	536	562	16			
1	B	391	Total	C	N	O	S	0	1	0
			2957	1846	535	560	16			

There are 4 discrepancies between the modelled and reference sequences:

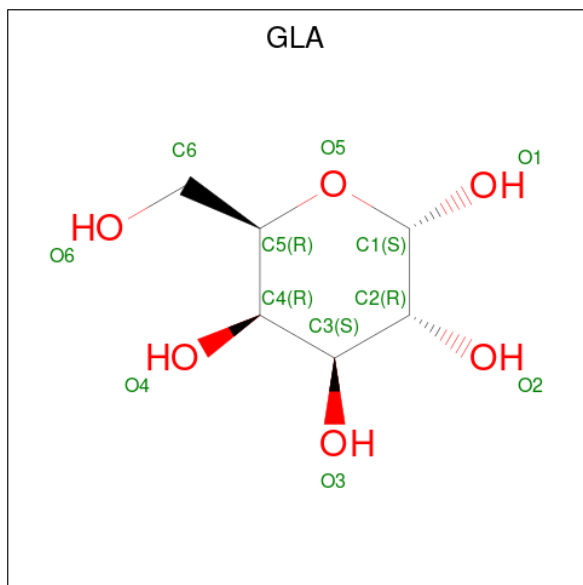
Chain	Residue	Modelled	Actual	Comment	Reference
A	252	ALA	LYS	engineered mutation	UNP P51570
A	253	ALA	GLU	engineered mutation	UNP P51570
B	252	ALA	LYS	engineered mutation	UNP P51570
B	253	ALA	GLU	engineered mutation	UNP P51570

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: 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 alpha-D-galactopyranose (CCD ID: GLA) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		
5	B	1	Total	Mg	0	0
			1	1		

- Molecule 6 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Na	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Na	0	0
			1	1		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

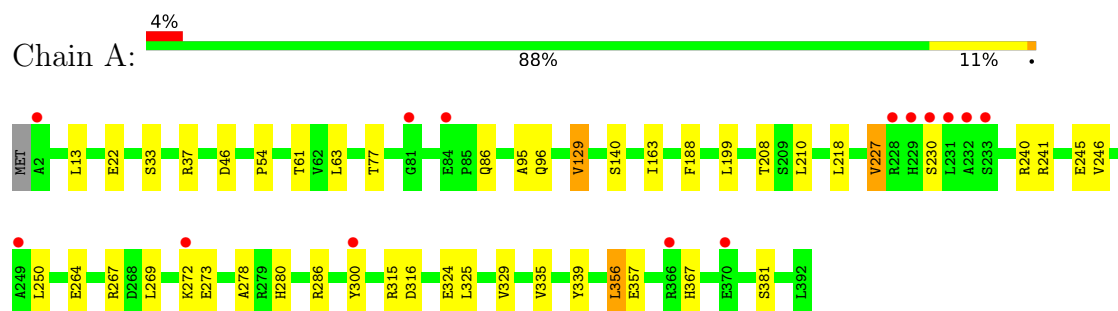
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	111	Total	O	0	0
			111	111		
8	B	138	Total	O	0	0
			138	138		

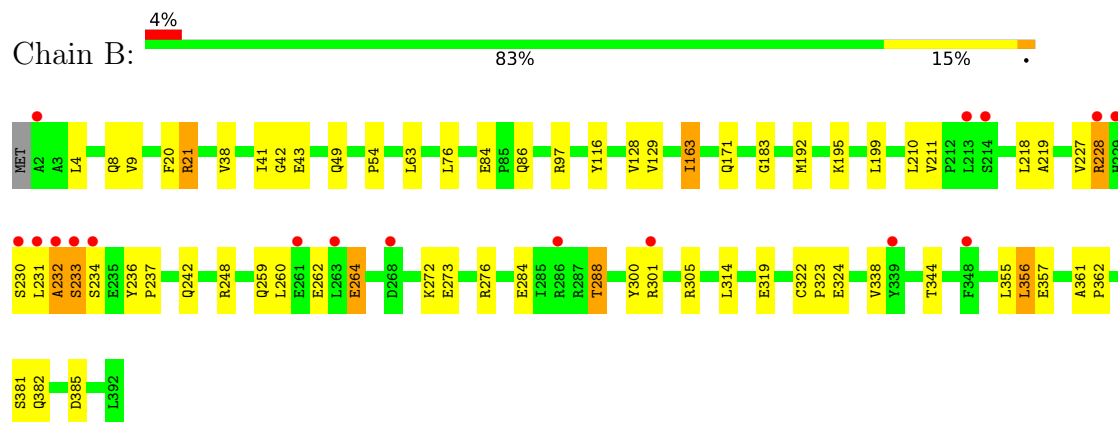
3 Residue-property plots [i](#)

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: Galactokinase



• Molecule 1: Galactokinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	128.79Å 128.79Å 240.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.82 – 2.10 39.82 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.1 (39.82-2.10) 99.1 (39.82-2.10)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.196 , 0.236 0.203 , 0.240	Depositor DCC
R_{free} test set	3072 reflections (4.44%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.654	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6287	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.93% 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: GOL, MG, ADP, GLA, NA, PO4

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	1.09	0/3017	1.46	3/4092 (0.1%)
1	B	1.13	0/3011	1.49	7/4084 (0.2%)
All	All	1.11	0/6028	1.48	10/8176 (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	A	0	1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	21	ARG	CA-C-N	5.99	128.56	120.65
1	B	21	ARG	C-N-CA	5.99	128.56	120.65
1	A	280	HIS	CA-CB-CG	-5.91	107.89	113.80
1	A	46	ASP	CA-CB-CG	5.70	118.30	112.60
1	B	20	PHE	CA-C-N	5.64	127.84	120.28
1	B	20	PHE	C-N-CA	5.64	127.84	120.28
1	B	262	GLU	CB-CA-C	-5.35	102.48	110.88
1	B	183	GLY	CA-C-O	-5.28	116.77	122.47
1	B	9	VAL	CA-C-O	-5.03	116.04	121.27
1	A	240	ARG	CG-CD-NE	-5.02	100.95	112.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	230	SER	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	2963	0	2957	28	0
1	B	2957	0	2952	40	0
2	A	27	0	12	0	0
2	B	27	0	12	0	0
3	A	12	0	11	0	0
3	B	12	0	11	0	0
4	A	15	0	0	1	0
4	B	15	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	B	6	0	8	2	0
8	A	111	0	0	2	1
8	B	138	0	0	7	1
All	All	6287	0	5963	67	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 6.

All (67) 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:63:LEU:HD11	1:B:129:VAL:HG22	1.38	1.03
1:B:171:GLN:HE22	7:B:407:GOL:H12	1.32	0.92
1:B:63:LEU:CD1	1:B:129:VAL:HG22	2.10	0.81
1:A:163:ILE:H	1:A:163:ILE:HD12	1.57	0.69
1:B:116:TYR:O	8:B:501:HOH:O	2.12	0.67
1:A:54:PRO:HD2	1:A:199:LEU:O	1.94	0.67
1:B:284:GLU:O	1:B:288:THR:HG23	1.95	0.66
1:B:171:GLN:NE2	7:B:407:GOL:H12	2.10	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:284:GLU:O	1:B:288:THR:CG2	2.48	0.62
1:B:218:LEU:HD22	1:B:300:TYR:CE1	2.35	0.61
1:B:361:ALA:HB3	1:B:362:PRO:HD3	1.83	0.60
1:B:276:ARG:NH2	1:B:319:GLU:OE2	2.33	0.60
1:B:195:LYS:HB2	1:B:385:ASP:OD1	2.02	0.59
1:A:218:LEU:HD22	1:A:300:TYR:CE1	2.39	0.58
1:B:41:ILE:HG23	1:B:288:THR:HB	1.86	0.57
1:B:356:LEU:HD13	1:B:357:GLU:O	2.04	0.57
1:B:38:VAL:HG23	1:B:344:THR:HG21	1.86	0.56
1:B:231:LEU:HD23	8:B:557:HOH:O	2.07	0.54
1:B:43:GLU:HB3	1:B:314:LEU:HD21	1.90	0.53
1:B:236:TYR:HB3	1:B:237:PRO:HD3	1.90	0.53
1:A:335:VAL:HG23	1:A:367:HIS:HD2	1.75	0.52
1:B:8:GLN:HB3	8:B:593:HOH:O	2.10	0.51
1:A:77:THR:HA	1:A:129[A]:VAL:HG22	1.92	0.50
1:B:227:VAL:HG21	1:B:324:GLU:HG2	1.93	0.50
1:B:54:PRO:HD2	1:B:199:LEU:O	2.12	0.49
1:B:356:LEU:C	1:B:356:LEU:HD12	2.37	0.49
1:A:241:ARG:HB3	1:A:241:ARG:NH1	2.29	0.47
1:A:199:LEU:HB2	8:A:589:HOH:O	2.14	0.47
1:B:232:ALA:C	1:B:234:SER:H	2.22	0.47
1:A:315:ARG:HD2	1:A:316:ASP:OD1	2.15	0.47
1:B:219:ALA:HB3	1:B:356:LEU:CD1	2.45	0.47
1:B:301:ARG:HH12	1:B:305:ARG:HD3	1.80	0.46
1:A:356:LEU:HD12	1:A:356:LEU:H	1.80	0.46
1:A:77:THR:HA	1:A:129[B]:VAL:HG13	1.98	0.45
1:A:163:ILE:HD12	1:A:163:ILE:N	2.29	0.45
1:A:356:LEU:HD12	1:A:356:LEU:N	2.32	0.44
1:A:246:VAL:HG11	1:A:278:ALA:HB2	1.98	0.44
1:B:356:LEU:C	1:B:356:LEU:CD1	2.91	0.44
1:B:233:SER:O	1:B:233:SER:OG	2.32	0.44
1:A:188:PHE:CE1	1:A:208:THR:HG21	2.53	0.44
1:A:250:LEU:HD23	1:A:269:LEU:HD22	2.00	0.43
1:B:219:ALA:HB3	1:B:356:LEU:HD12	2.00	0.43
1:A:325:LEU:O	1:A:329:VAL:HG23	2.18	0.43
1:B:228:ARG:HD3	8:B:506:HOH:O	2.18	0.43
1:B:381:SER:HB3	8:B:588:HOH:O	2.18	0.43
1:A:227:VAL:HG21	1:A:324:GLU:HG2	2.01	0.43
1:A:267:ARG:HH22	4:A:403:PO4:P	2.41	0.43
1:A:339:TYR:OH	1:A:357:GLU:OE2	2.34	0.43
1:B:322:CYS:O	1:B:323:PRO:C	2.62	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:LEU:HD13	1:B:192:MET:HE3	2.01	0.42
1:A:286:ARG:NH2	8:A:518:HOH:O	2.52	0.42
1:B:163:ILE:HG21	1:B:192:MET:CE	2.50	0.42
1:B:163:ILE:HG21	1:B:192:MET:SD	2.60	0.42
1:A:163:ILE:H	1:A:163:ILE:CD1	2.31	0.41
1:A:241:ARG:NH1	1:A:241:ARG:CB	2.83	0.41
1:B:42:GLY:CA	1:B:288:THR:HG21	2.51	0.41
1:B:86:GLN:HG2	8:B:539:HOH:O	2.21	0.41
1:B:259:GLN:HB3	8:B:551:HOH:O	2.20	0.41
1:A:33:SER:HA	1:A:61:THR:O	2.20	0.41
1:A:37:ARG:CZ	1:A:140[B]:SER:OG	2.69	0.41
1:B:242:GLN:HE22	1:B:273:GLU:HG2	1.85	0.41
1:B:260:LEU:HD11	1:B:264:GLU:OE1	2.20	0.41
1:A:241:ARG:CB	1:A:241:ARG:CZ	2.99	0.41
1:A:241:ARG:HB3	1:A:241:ARG:HH11	1.86	0.40
1:B:76:LEU:O	1:B:128:VAL:HA	2.21	0.40
1:A:13:LEU:HD23	1:A:13:LEU:HA	1.89	0.40
1:B:199:LEU:HD11	1:B:210:LEU:CD2	2.50	0.40

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)
8:A:604:HOH:O	8:B:605:HOH:O[5_554]	2.11	0.09

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	391/392 (100%)	378 (97%)	12 (3%)	1 (0%)	36	36
1	B	390/392 (100%)	375 (96%)	14 (4%)	1 (0%)	36	36

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	781/784 (100%)	753 (96%)	26 (3%)	2 (0%)	36	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	232	ALA
1	A	95	ALA

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	310/309 (100%)	297 (96%)	13 (4%)	26	28
1	B	309/309 (100%)	291 (94%)	18 (6%)	18	16
All	All	619/618 (100%)	588 (95%)	31 (5%)	22	22

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	63	LEU
1	A	86	GLN
1	A	96	GLN
1	A	129[A]	VAL
1	A	129[B]	VAL
1	A	227	VAL
1	A	245	GLU
1	A	264	GLU
1	A	272	LYS
1	A	273	GLU
1	A	356	LEU
1	A	381	SER
1	B	4	LEU
1	B	21	ARG
1	B	49	GLN

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Mol	Chain	Res	Type
1	B	84	GLU
1	B	97	ARG
1	B	163	ILE
1	B	211	VAL
1	B	228	ARG
1	B	230	SER
1	B	233	SER
1	B	248	ARG
1	B	264	GLU
1	B	272	LYS
1	B	288	THR
1	B	338	VAL
1	B	355	LEU
1	B	356	LEU
1	B	382	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	367	HIS
1	A	371	HIS
1	B	229	HIS
1	B	367	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 4 are monoatomic - leaving 11 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GLA	B	402	-	12,12,12	2.80	5 (41%)	17,17,17	1.91	4 (23%)
4	PO4	B	404	-	4,4,4	0.66	0	6,6,6	0.48	0
2	ADP	B	401	5	28,29,29	0.80	1 (3%)	43,45,45	0.73	0
4	PO4	B	405	-	4,4,4	0.67	0	6,6,6	0.46	0
4	PO4	A	405	-	4,4,4	0.80	0	6,6,6	0.45	0
4	PO4	B	403	-	4,4,4	0.70	0	6,6,6	0.48	0
7	GOL	B	407	-	5,5,5	0.18	0	5,5,5	0.48	0
2	ADP	A	401	5	28,29,29	0.65	0	43,45,45	0.68	1 (2%)
4	PO4	A	404	-	4,4,4	0.67	0	6,6,6	0.48	0
4	PO4	A	403	-	4,4,4	0.61	0	6,6,6	0.48	0
3	GLA	A	402	-	12,12,12	2.93	6 (50%)	17,17,17	2.10	5 (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
3	GLA	B	402	-	-	1/2/22/22	0/1/1/1
2	ADP	B	401	5	-	3/16/32/32	0/3/3/3
7	GOL	B	407	-	-	4/4/4/4	-
2	ADP	A	401	5	-	3/16/32/32	0/3/3/3
3	GLA	A	402	-	-	1/2/22/22	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	GLA	O5-C1	-6.02	1.28	1.42
3	A	402	GLA	O5-C1	-5.47	1.29	1.42
3	A	402	GLA	O3-C3	-4.62	1.31	1.43
3	B	402	GLA	O3-C3	-4.49	1.31	1.43
3	A	402	GLA	O5-C5	-4.36	1.33	1.44
3	B	402	GLA	O5-C5	-3.56	1.35	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	ADP	PA-O3A	3.30	1.63	1.59
3	A	402	GLA	C4-C3	3.25	1.60	1.52
3	B	402	GLA	C4-C3	3.20	1.60	1.52
3	A	402	GLA	O1-C1	2.53	1.47	1.39
3	A	402	GLA	C3-C2	2.27	1.58	1.52
3	B	402	GLA	O1-C1	2.27	1.46	1.39

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	GLA	C1-O5-C5	5.43	124.15	113.65
3	A	402	GLA	C1-O5-C5	4.62	122.59	113.65
3	A	402	GLA	C4-C3-C2	-3.86	104.05	110.83
3	B	402	GLA	C4-C3-C2	-3.04	105.50	110.83
3	A	402	GLA	O1-C1-C2	-2.98	100.34	108.98
3	A	402	GLA	C1-C2-C3	2.98	116.42	110.36
3	B	402	GLA	O1-C1-C2	-2.47	101.82	108.98
3	A	402	GLA	O5-C5-C6	2.28	112.10	106.44
3	B	402	GLA	O5-C5-C6	2.09	111.62	106.44
2	A	401	ADP	O2A-PA-O1A	2.07	122.08	112.44

There are no chirality outliers.

All (12) torsion outliers are listed below:

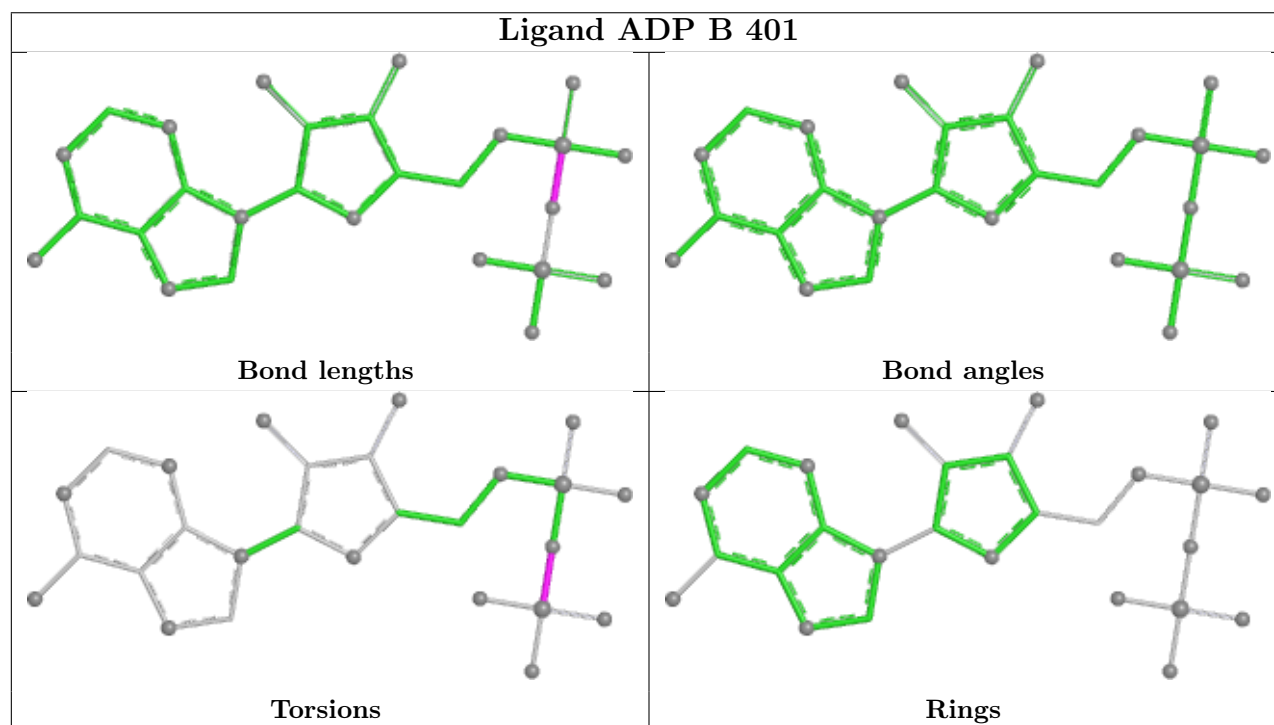
Mol	Chain	Res	Type	Atoms
2	A	401	ADP	PA-O3A-PB-O2B
7	B	407	GOL	O1-C1-C2-O2
7	B	407	GOL	O1-C1-C2-C3
7	B	407	GOL	C1-C2-C3-O3
7	B	407	GOL	O2-C2-C3-O3
3	A	402	GLA	O5-C5-C6-O6
3	B	402	GLA	O5-C5-C6-O6
2	B	401	ADP	PA-O3A-PB-O1B
2	A	401	ADP	PA-O3A-PB-O1B
2	A	401	ADP	PA-O3A-PB-O3B
2	B	401	ADP	PA-O3A-PB-O2B
2	B	401	ADP	PA-O3A-PB-O3B

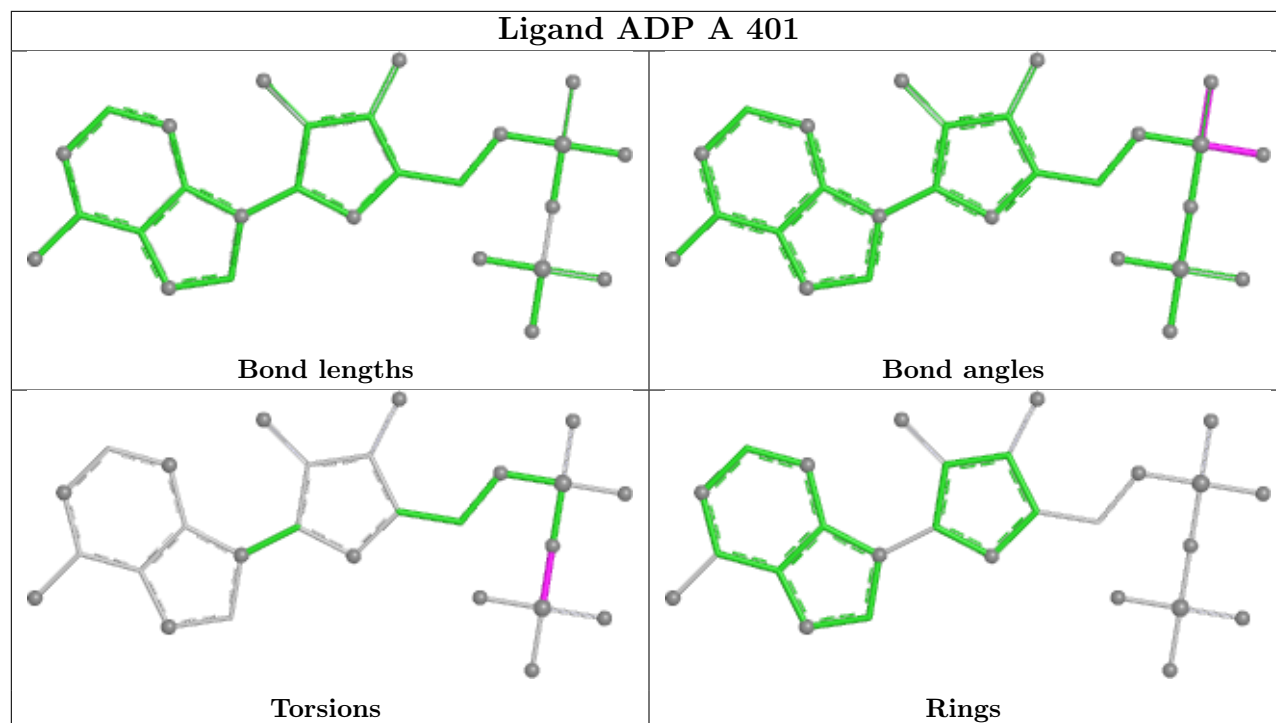
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	407	GOL	2	0
4	A	403	PO4	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

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	391/392 (99%)	0.25	14 (3%) 46 48	17, 49, 75, 117	2 (0%)
1	B	391/392 (99%)	0.29	17 (4%) 40 41	23, 49, 85, 110	1 (0%)
All	All	782/784 (99%)	0.27	31 (3%) 42 44	17, 49, 81, 117	3 (0%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	231	LEU	6.5
1	A	231	LEU	6.4
1	B	232	ALA	6.2
1	B	233	SER	4.4
1	B	229	HIS	4.1
1	A	228	ARG	3.9
1	B	230	SER	3.6
1	B	2	ALA	3.6
1	A	249	ALA	3.6
1	A	2	ALA	3.3
1	B	228	ARG	3.1
1	A	232	ALA	3.1
1	A	233	SER	2.9
1	B	348	PHE	2.9
1	B	234	SER	2.6
1	A	230	SER	2.5
1	B	214	SER	2.5
1	A	272	LYS	2.5
1	B	286	ARG	2.4
1	B	213	LEU	2.4
1	A	81	GLY	2.3
1	A	229	HIS	2.3
1	A	370	GLU	2.3
1	A	300	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	366	ARG	2.3
1	B	261	GLU	2.2
1	A	84	GLU	2.1
1	B	263	LEU	2.0
1	B	301	ARG	2.0
1	B	268	ASP	2.0
1	B	339	TYR	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](#)

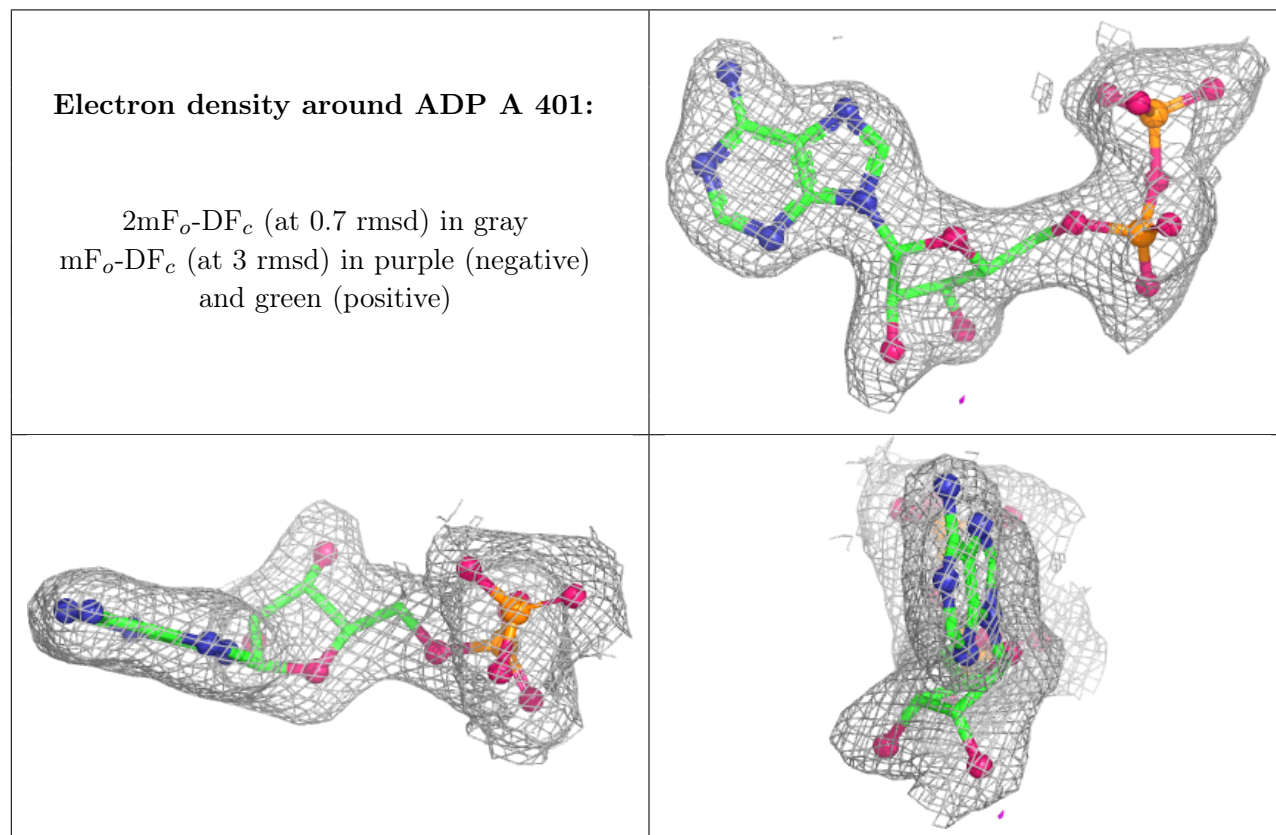
There are no oligosaccharides in this entry.

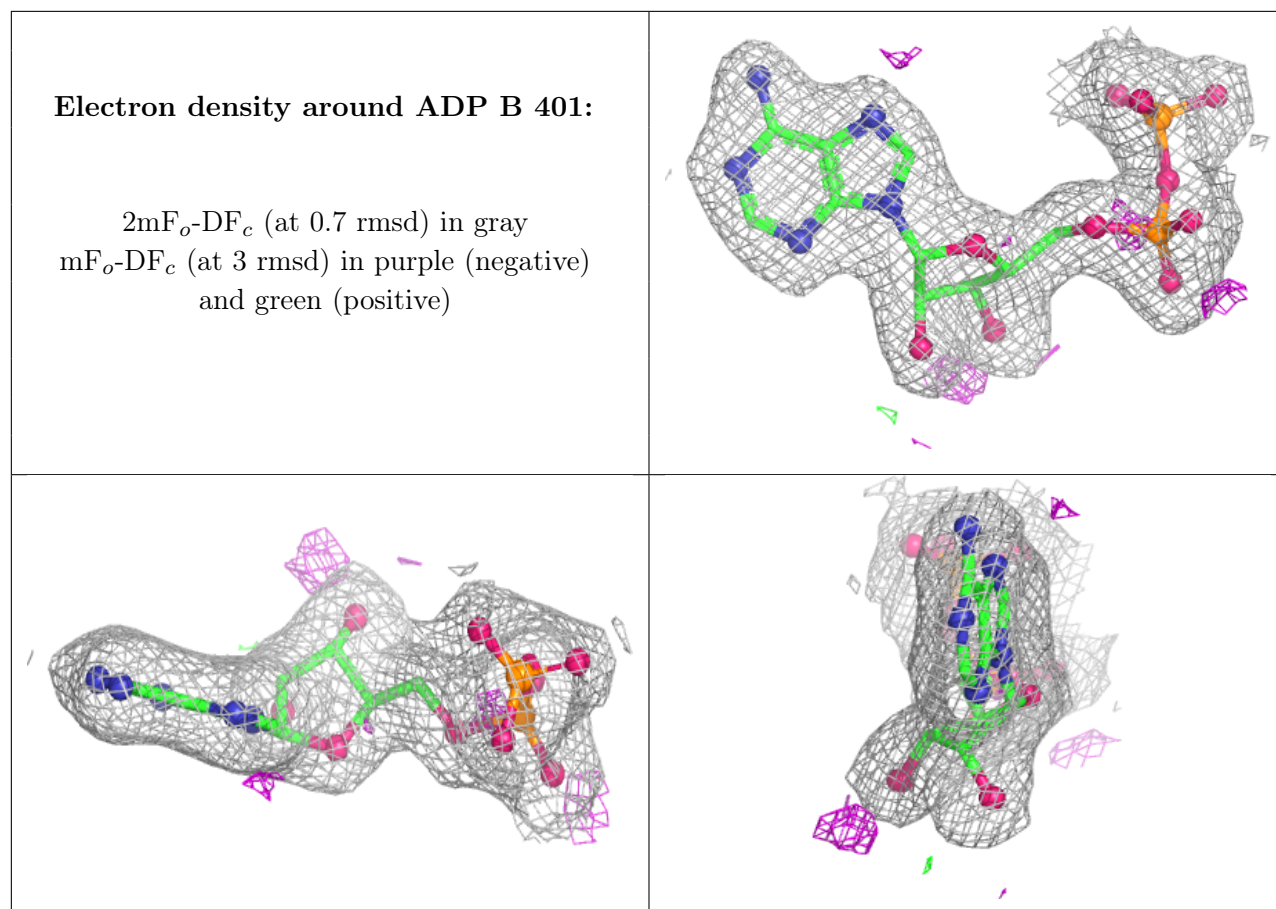
6.4 Ligands [i](#)

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
4	PO4	B	404	5/5	0.64	0.12	105,105,118,122	0
4	PO4	A	405	5/5	0.67	0.11	107,107,111,112	0
4	PO4	B	403	5/5	0.70	0.12	112,118,119,120	0
4	PO4	B	405	5/5	0.75	0.10	124,127,129,132	0
4	PO4	A	403	5/5	0.81	0.10	74,86,88,89	0
4	PO4	A	404	5/5	0.85	0.08	84,93,94,95	0
7	GOL	B	407	6/6	0.87	0.21	42,51,61,68	0
6	NA	B	408	1/1	0.93	0.06	50,50,50,50	0
5	MG	A	406	1/1	0.95	0.07	65,65,65,65	0
3	GLA	A	402	12/12	0.97	0.06	40,44,45,45	0
5	MG	B	406	1/1	0.97	0.08	48,48,48,48	0
2	ADP	A	401	27/27	0.98	0.05	36,40,43,44	0
6	NA	A	407	1/1	0.98	0.07	48,48,48,48	0
3	GLA	B	402	12/12	0.98	0.05	33,36,39,40	0
2	ADP	B	401	27/27	0.98	0.05	31,33,35,37	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 ⓘ

There are no such residues in this entry.