



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:19 PM UTC

PDB ID : 4PKI / pdb_00004pki
Title : Complex of ATP-actin With the C-terminal Actin-Binding Domain of Tropomodulin
Authors : Rao, J.N.; Dominguez, R.
Deposited on : 2014-05-14
Resolution : 2.30 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

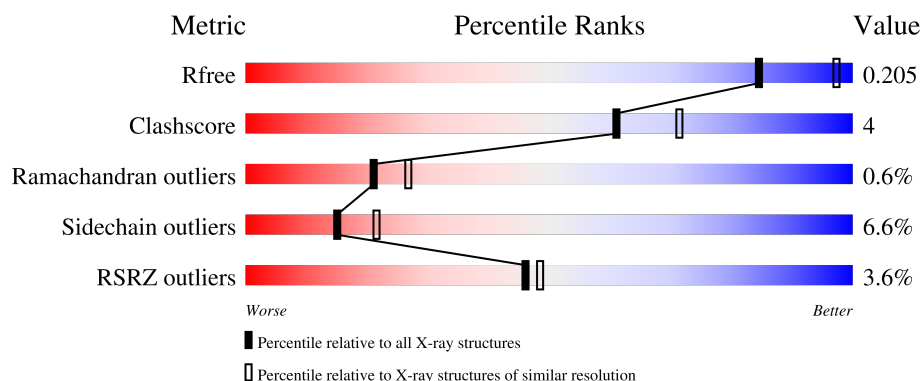
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	377	3% 83% 12% • •
2	G	324	4% 83% 10% • 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10970 atoms, of which 5353 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 Actin, alpha skeletal muscle.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	371	Total	C	H	N	O	S	0	8	0
			5856	1868	2917	490	556	25			

- Molecule 2 is a protein called Gelsolin,Tropomodulin-1 chimera.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	G	306	Total	C	H	N	O	S	0	2	0
			4861	1548	2425	415	465	8			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	177	GLY	-	linker	UNP P06396
G	178	GLY	-	linker	UNP P06396
G	179	SER	-	linker	UNP P06396
G	180	GLY	-	linker	UNP P06396
G	181	GLY	-	linker	UNP P06396
G	182	SER	-	linker	UNP P06396
G	183	GLY	-	linker	UNP P06396
G	184	GLY	-	linker	UNP P06396
G	185	SER	-	linker	UNP P06396

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
3	A	1	Total	C	H	N	O	P	0	0
			42	10	11	5	13	3		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Ca 2 2	0	0
4	G	1	Total Ca 1 1	0	0

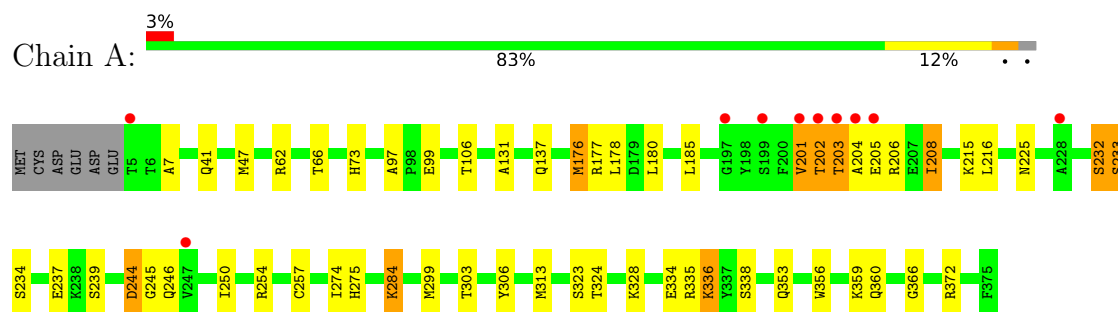
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	113	Total O 113 113	0	0
5	G	95	Total O 95 95	0	0

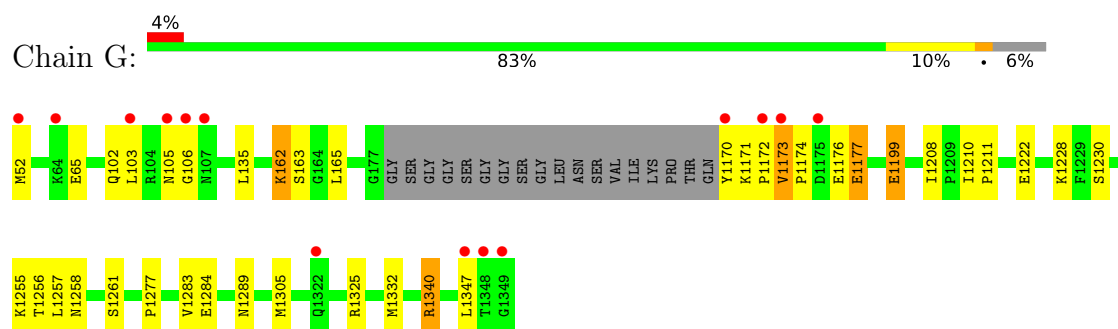
3 Residue-property plots [i](#)

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: Actin, alpha skeletal muscle



- Molecule 2: Gelsolin,Tropomodulin-1 chimera



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.39Å 81.21Å 170.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.00 – 2.30 45.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.00-2.30) 99.9 (45.00-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.168 , 0.201 0.174 , 0.205	Depositor DCC
R_{free} test set	2000 reflections (4.57%)	wwPDB-VP
Wilson B-factor (Å ²)	40.1	Xtriage
Anisotropy	0.555	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10970	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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: ATP, CA, HIC

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	0.84	1/3013 (0.0%)	0.90	3/4077 (0.1%)
2	G	0.73	0/2488	0.91	5/3363 (0.1%)
All	All	0.79	1/5501 (0.0%)	0.91	8/7440 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	7	ALA	CA-CB	-5.10	1.45	1.53

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	65	GLU	N-CA-C	7.72	114.94	108.07
2	G	1173	VAL	N-CA-C	-6.76	103.65	109.19
2	G	106	GLY	N-CA-C	-5.54	108.01	115.21
2	G	1208	ILE	CA-C-N	-5.34	114.28	119.78
2	G	1208	ILE	C-N-CA	-5.34	114.28	119.78
1	A	244	ASP	N-CA-C	5.30	119.37	113.01
1	A	257	CYS	CA-C-N	-5.08	114.81	120.45
1	A	257	CYS	C-N-CA	-5.08	114.81	120.45

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	2939	2917	2938	30	0
2	G	2436	2425	2431	16	0
3	A	31	11	12	0	0
4	A	2	0	0	0	0
4	G	1	0	0	0	0
5	A	113	0	0	4	0
5	G	95	0	0	1	0
All	All	5617	5353	5381	45	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 4.

All (45) 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:A:201:VAL:O	1:A:202:THR:OG1	2.09	0.66
2:G:1340:ARG:NH2	5:G:1578:HOH:O	2.20	0.62
1:A:202:THR:O	1:A:203:THR:OG1	2.15	0.61
1:A:372:ARG:NH2	5:A:501:HOH:O	2.32	0.59
1:A:244:ASP:N	1:A:245:GLY:HA2	2.18	0.59
1:A:131:ALA:HB1	1:A:356:TRP:HB3	1.87	0.56
1:A:176[B]:MET:SD	1:A:284:LYS:NZ	2.77	0.55
1:A:216:LEU:O	1:A:254:ARG:HD2	2.06	0.55
1:A:201:VAL:O	1:A:201:VAL:HG23	2.06	0.55
2:G:1177:GLU:OE1	2:G:1177:GLU:N	2.40	0.55
1:A:176[A]:MET:HE2	1:A:177:ARG:H	1.73	0.53
1:A:274:ILE:CD1	1:A:313:MET:HE1	2.40	0.51
1:A:178:LEU:HG	1:A:180:LEU:HB3	1.95	0.48
1:A:62:ARG:HE	1:A:204:ALA:HA	1.78	0.48
1:A:97:ALA:HB1	1:A:99:GLU:OE2	2.14	0.47
2:G:1261:SER:HA	2:G:1289:ASN:O	2.14	0.47
1:A:202:THR:HG22	1:A:203:THR:N	2.30	0.46
1:A:176[A]:MET:HA	1:A:176[A]:MET:CE	2.45	0.46
2:G:1277:PRO:HG3	2:G:1305:MET:HE1	1.98	0.46
1:A:244:ASP:OD1	1:A:246:GLN:HG3	2.17	0.45
2:G:1283:VAL:HG23	2:G:1284:GLU:HG3	1.99	0.45
2:G:165:LEU:C	2:G:165:LEU:HD12	2.42	0.44
1:A:232:SER:OG	1:A:233:SER:N	2.50	0.44
1:A:205:GLU:O	1:A:208:ILE:HG22	2.18	0.44
1:A:324:THR:O	1:A:324:THR:OG1	2.28	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:GLU:OE2	1:A:334:GLU:N	2.46	0.44
2:G:1210:ILE:N	2:G:1211:PRO:CD	2.81	0.44
1:A:335:ARG:HA	1:A:338:SER:OG	2.18	0.43
2:G:162:LYS:CD	2:G:162:LYS:H	2.30	0.43
2:G:1199:GLU:HG3	2:G:1228:LYS:HB2	2.00	0.43
1:A:185:LEU:HD23	1:A:306:TYR:OH	2.18	0.43
2:G:1228:LYS:NZ	2:G:1256:THR:HG21	2.34	0.42
2:G:1230:SER:CB	2:G:1258:ASN:HB3	2.49	0.42
2:G:1173:VAL:HG22	2:G:1174:PRO:HD2	2.02	0.42
1:A:106:THR:HB	1:A:137:GLN:HG3	2.02	0.42
1:A:366:GLY:HA2	2:G:1340:ARG:HD2	2.03	0.41
2:G:1332:MET:HA	2:G:1332:MET:HE2	2.02	0.41
2:G:1230:SER:HB3	2:G:1258:ASN:HB3	2.01	0.41
1:A:275:HIS:CD2	1:A:275:HIS:H	2.38	0.41
2:G:1173:VAL:CG2	2:G:1174:PRO:HD2	2.51	0.41
1:A:359:LYS:NZ	5:A:556:HOH:O	2.54	0.41
1:A:303:THR:O	1:A:303:THR:HG22	2.22	0.40
1:A:336:LYS:HB2	5:A:506:HOH:O	2.21	0.40
1:A:233:SER:O	1:A:234:SER:CB	2.69	0.40
1:A:324:THR:HG23	5:A:509:HOH:O	2.21	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	376/377 (100%)	358 (95%)	15 (4%)	3 (1%)	16	20
2	G	304/324 (94%)	289 (95%)	14 (5%)	1 (0%)	36	46
All	All	680/701 (97%)	647 (95%)	29 (4%)	4 (1%)	21	27

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	202	THR
2	G	1172	PRO
1	A	203	THR
1	A	232	SER

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	321/319 (101%)	298 (93%)	23 (7%)	13	18
2	G	266/276 (96%)	248 (93%)	18 (7%)	14	20
All	All	587/595 (99%)	546 (93%)	41 (7%)	15	19

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	47	MET
1	A	66	THR
1	A	176[A]	MET
1	A	176[B]	MET
1	A	201	VAL
1	A	206	ARG
1	A	208	ILE
1	A	215	LYS
1	A	225	ASN
1	A	233	SER
1	A	237	GLU
1	A	239	SER
1	A	250	ILE
1	A	284	LYS
1	A	299[A]	MET
1	A	299[B]	MET
1	A	323	SER
1	A	328	LYS
1	A	336	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	353	GLN
1	A	360[A]	GLN
1	A	360[B]	GLN
2	G	52	MET
2	G	102	GLN
2	G	103	LEU
2	G	105	ASN
2	G	135	LEU
2	G	162	LYS
2	G	163	SER
2	G	1170	TYR
2	G	1171	LYS
2	G	1176	GLU
2	G	1177	GLU
2	G	1199	GLU
2	G	1222	GLU
2	G	1255	LYS
2	G	1257	LEU
2	G	1325	ARG
2	G	1340	ARG
2	G	1347	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	ASN
1	A	162	ASN
1	A	225	ASN
1	A	275	HIS
1	A	280	ASN
2	G	1204	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	HIC	A	73	1	10,11,12	1.50	2 (20%)	9,14,16	1.66	1 (11%)

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
1	HIC	A	73	1	-	0/5/6/8	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	73	HIC	CD2-CG	2.95	1.41	1.36
1	A	73	HIC	CD2-NE2	2.22	1.40	1.37

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	73	HIC	NE2-CE1-ND1	-4.25	111.04	112.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 3 are monoatomic - leaving 1 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	ATP	A	401	4	32,33,33	1.38	3 (9%)	48,52,52	1.83	11 (22%)

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	ATP	A	401	4	-	2/22/38/38	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	ATP	C5-C4	4.95	1.47	1.39
3	A	401	ATP	C8-N7	2.47	1.36	1.31
3	A	401	ATP	C5-C6	2.37	1.47	1.41

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	ATP	C5-C4-N3	-5.56	119.06	126.72
3	A	401	ATP	N3-C4-N9	4.79	135.31	127.17
3	A	401	ATP	C2-N3-C4	3.40	120.15	111.83
3	A	401	ATP	N3-C2-N1	-3.37	123.48	128.58
3	A	401	ATP	C4-N9-C8	3.16	109.06	105.74
3	A	401	ATP	C4-C5-N7	-2.84	107.34	110.58
3	A	401	ATP	O4'-C1'-N9	2.52	112.92	108.09
3	A	401	ATP	C2-N1-C6	2.36	122.60	118.73
3	A	401	ATP	C5-N7-C8	2.26	107.00	103.45
3	A	401	ATP	O2A-PA-O3A	2.14	113.05	107.27
3	A	401	ATP	N9-C8-N7	-2.09	110.98	113.94

There are no chirality outliers.

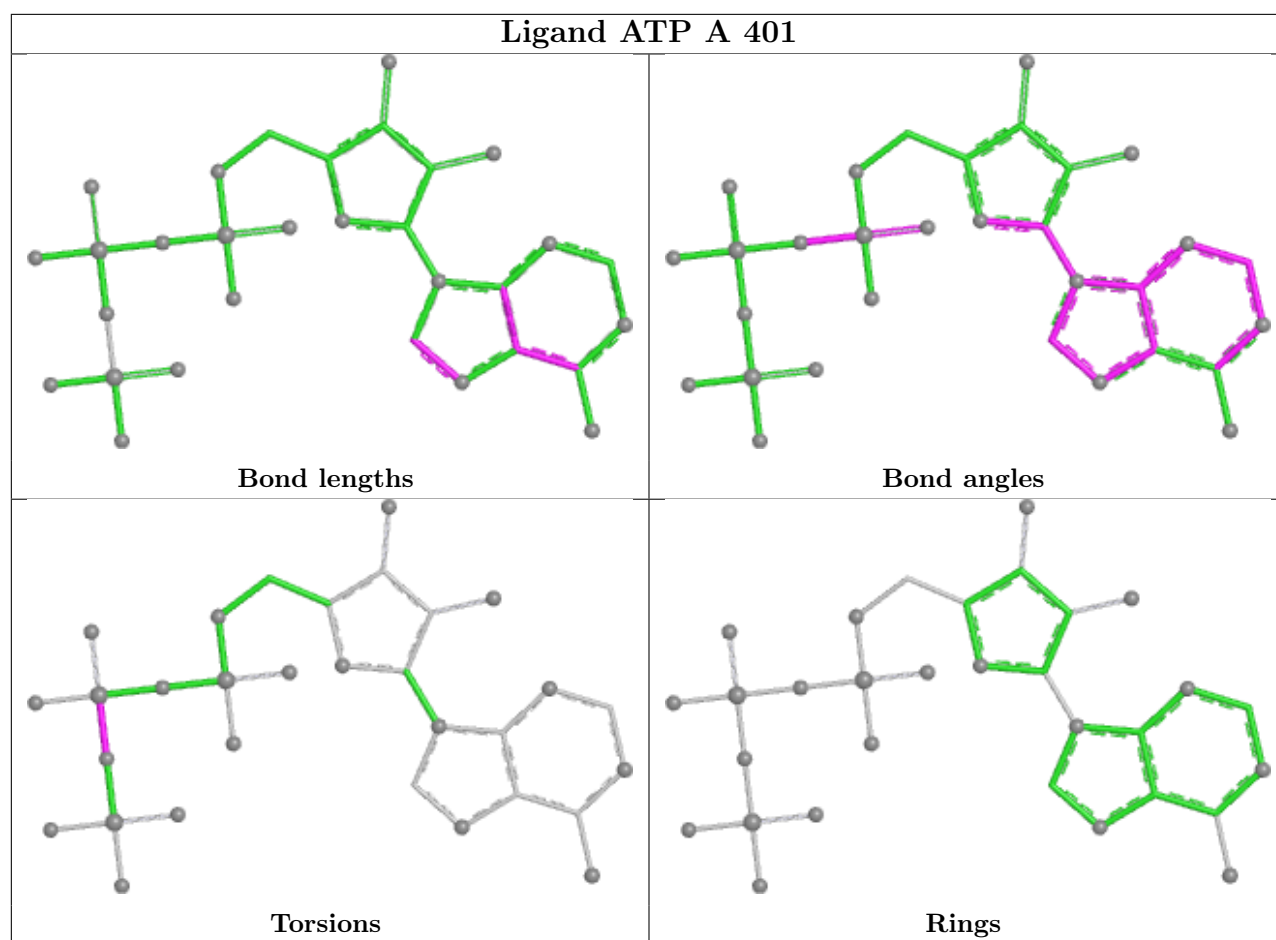
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	ATP	PG-O3B-PB-O1B
3	A	401	ATP	PG-O3B-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

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	370/377 (98%)	-0.07	10 (2%) 56 58	17, 49, 111, 145	8 (2%)
2	G	306/324 (94%)	0.05	14 (4%) 37 39	29, 59, 104, 160	2 (0%)
All	All	676/701 (96%)	-0.02	24 (3%) 46 48	17, 54, 109, 160	10 (1%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	THR	5.4
2	G	1172	PRO	4.6
2	G	1170	TYR	4.5
1	A	204	ALA	4.2
2	G	103	LEU	4.1
1	A	203	THR	3.9
1	A	202	THR	3.7
2	G	52	MET	3.6
2	G	1173	VAL	3.5
1	A	199	SER	3.4
1	A	201	VAL	3.3
2	G	106	GLY	3.2
1	A	228	ALA	2.9
2	G	1175	ASP	2.9
1	A	197	GLY	2.9
2	G	1349	GLY	2.8
2	G	1348	THR	2.7
1	A	205	GLU	2.7
2	G	64	LYS	2.6
2	G	107	ASN	2.6
2	G	1347	LEU	2.4
2	G	1322	GLN	2.4
1	A	247	VAL	2.3
2	G	105	ASN	2.2

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	HIC	A	73	11/12	0.95	0.12	28,62,88,88	0

6.3 Carbohydrates [i](#)

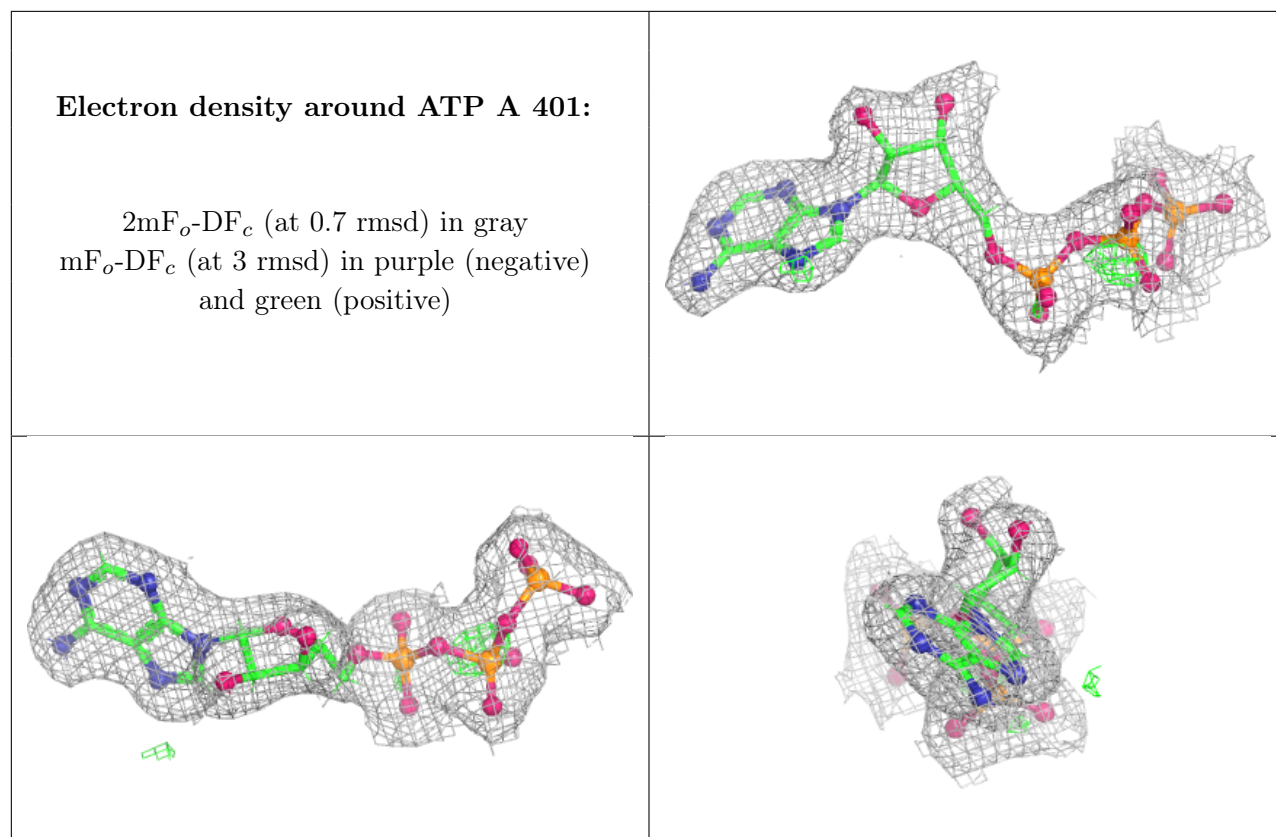
There are no oligosaccharides in this entry.

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	ATP	A	401	31/31	0.99	0.06	32,41,53,55	0
4	CA	A	403	1/1	0.99	0.03	41,41,41,41	0
4	CA	A	402	1/1	1.00	0.04	37,37,37,37	0
4	CA	G	1401	1/1	1.00	0.03	40,40,40,40	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.