



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:39 PM GMT

PDB ID : 4PKI
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

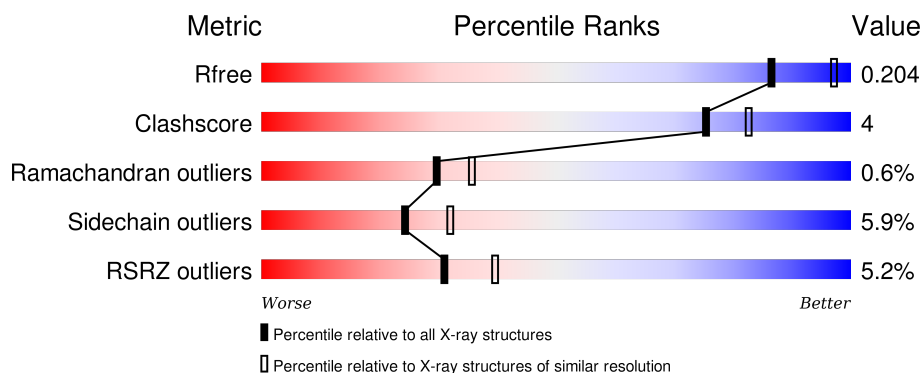
1 Overall quality at a glance

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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>5%</div> <div>84%</div> <div>12%</div> <div>• •</div> </div>
2	G	324	<div> <div>5%</div> <div>85%</div> <div>8%</div> <div>• 6%</div> </div>

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			

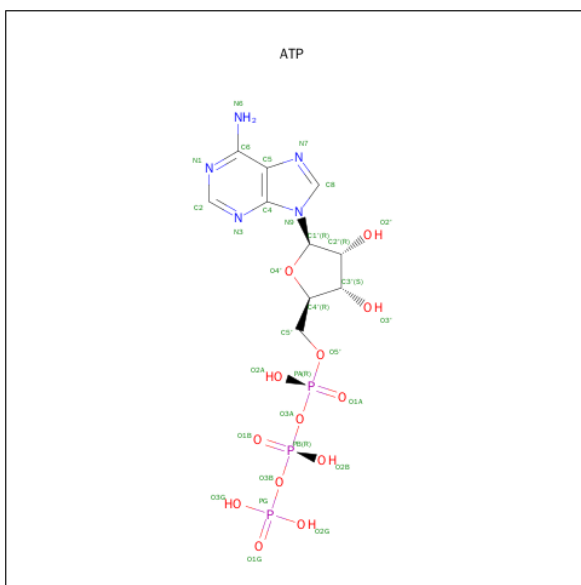
- Molecule 2 is a protein called Gelsolin,Tropomodulin-1 chimera.

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 (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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 (three-letter code: CA) (formula: Ca).

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

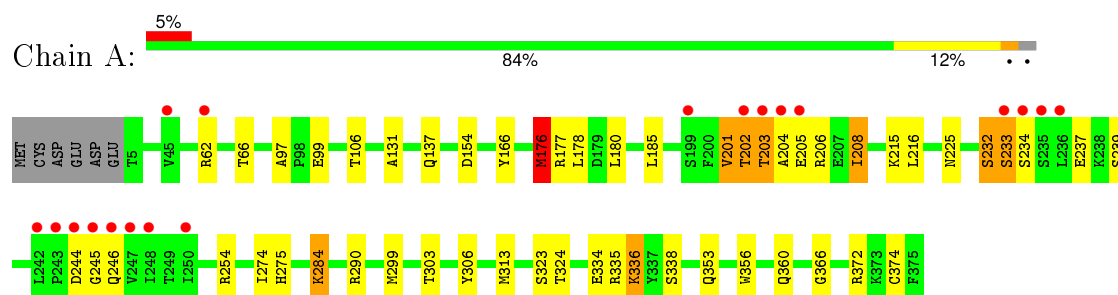
- Molecule 5 is water.

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

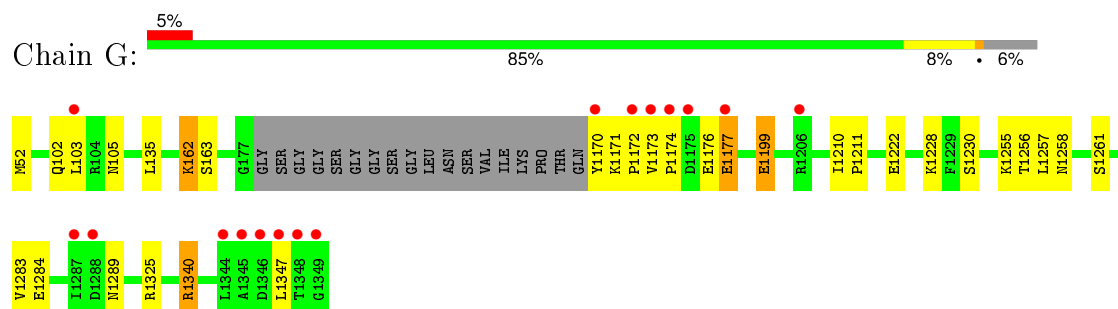
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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 44.88 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.00-2.30) 99.8 (44.88-2.30)	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 0.170 , 0.204	Depositor DCC
R_{free} test set	1997 reflections (4.58%)	DCC
Wilson B-factor (Å ²)	40.1	Xtriage
Anisotropy	0.555	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 43679 reflections	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, 0.375 respectively for untwinned datasets, and 0.333, 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: CA, HIC, ATP

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	0.69	2/3013 (0.1%)	0.74	4/4077 (0.1%)
2	G	0.64	0/2488	0.66	0/3363
All	All	0.67	2/5501 (0.0%)	0.70	4/7440 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	166	TYR	CD1-CE1	5.65	1.47	1.39
1	A	374	CYS	CB-SG	-5.30	1.73	1.81

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	154	ASP	CB-CG-OD1	6.73	124.35	118.30
1	A	176[A]	MET	CG-SD-CE	5.91	109.66	100.20
1	A	176[B]	MET	CG-SD-CE	5.91	109.66	100.20
1	A	290	ARG	NE-CZ-NH1	-5.25	117.68	120.30

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	13	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	3	0
5	G	95	0	0	1	0
All	All	5617	5353	5381	42	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 (42) 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
1:A:202:THR:O	1:A:203:THR:OG1	2.15	0.65
2:G:1340:ARG:NH2	5:G:1578:HOH:O	2.20	0.62
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:216:LEU:O	1:A:254:ARG:HD2	2.06	0.55
1:A:176[B]:MET:SD	1:A:284:LYS:NZ	2.77	0.55
1:A:201:VAL:HG23	1:A:201:VAL:O	2.06	0.55
2:G:1177:GLU:N	2:G:1177:GLU:OE1	2.40	0.55
1:A:176[A]:MET:HE2	1:A:177:ARG:H	1.77	0.50
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
1:A:274:ILE:CD1	1:A:313:MET:HE1	2.46	0.45
1:A:244:ASP:OD1	1:A:246:GLN:HG3	2.17	0.45
1:A:324:THR:O	1:A:324:THR:OG1	2.28	0.45
2:G:1283:VAL:HG23	2:G:1284:GLU:HG3	1.99	0.45
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:334:GLU:N	1:A:334:GLU:OE2	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:H	2:G:162:LYS:CD	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:176[A]:MET:HA	1:A:176[A]:MET:HE3	2.02	0.41
1:A:366:GLY:HA2	2:G:1340:ARG:HD2	2.03	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:303:THR:HG22	1:A:303:THR:O	2.22	0.40
1:A:233:SER:O	1:A:234:SER:CB	2.69	0.40
1:A:336:LYS:HB2	5:A:506:HOH:O	2.21	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%)	24	27
2	G	304/324 (94%)	289 (95%)	14 (5%)	1 (0%)	46	57
All	All	680/701 (97%)	647 (95%)	29 (4%)	4 (1%)	30	36

All (4) Ramachandran outliers are listed below:

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

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Mol	Chain	Res	Type
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%)	302 (94%)	19 (6%)	24	32
2	G	266/276 (96%)	248 (93%)	18 (7%)	20	25
All	All	587/595 (99%)	550 (94%)	37 (6%)	24	29

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
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	284	LYS
1	A	299[A]	MET
1	A	299[B]	MET
1	A	323	SER
1	A	336	LYS
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

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Mol	Chain	Res	Type
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

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) 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	105	ASN
2	G	134	GLN

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	8,11,12	1.65	2 (25%)	5,14,16	0.70	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
1	HIC	A	73	1	-	0/4/6/8	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	CZ-NE2	-2.27	1.42	1.48
1	A	73	HIC	CD2-CG	3.46	1.41	1.36

There are no bond angle outliers.

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	24,33,33	1.12	2 (8%)	31,52,52	1.86	7 (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	-	0/18/38/38	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	ATP	O4'-C1'	2.18	1.44	1.41
3	A	401	ATP	C5-C4	3.29	1.47	1.40

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	ATP	N3-C2-N1	-7.07	123.48	128.89
3	A	401	ATP	C2'-C1'-N9	-2.93	109.81	114.29
3	A	401	ATP	C4-C5-N7	-2.33	107.34	109.48
3	A	401	ATP	O5'-PA-O1A	-2.19	101.12	109.62
3	A	401	ATP	PA-O3A-PB	-2.06	126.95	132.73
3	A	401	ATP	C2-N1-C6	2.15	122.60	118.77
3	A	401	ATP	O4'-C1'-N9	2.30	112.92	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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%)	0.25	19 (5%) 32 41	25, 45, 101, 133	0
2	G	306/324 (94%)	0.28	16 (5%) 31 39	27, 54, 95, 145	0
All	All	676/701 (96%)	0.27	35 (5%) 31 39	25, 50, 98, 145	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	G	1348	THR	9.8
2	G	1170	TYR	6.8
2	G	1347	LEU	6.5
2	G	1173	VAL	5.8
1	A	203	THR	5.1
2	G	1349	GLY	5.1
1	A	244	ASP	5.1
2	G	1172	PRO	4.7
1	A	243	PRO	4.4
1	A	204	ALA	4.3
2	G	1346	ASP	4.0
2	G	1344	LEU	3.8
1	A	246	GLN	3.8
1	A	248	ILE	3.8
1	A	242	LEU	3.4
2	G	1345	ALA	3.4
2	G	1175	ASP	3.2
1	A	233	SER	3.1
1	A	235	SER	2.9
2	G	1174	PRO	2.8
1	A	250	ILE	2.8
1	A	247	VAL	2.7
1	A	202	THR	2.6
2	G	1206	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	245	GLY	2.4
1	A	234	SER	2.4
2	G	1177	GLU	2.4
2	G	1288	ASP	2.3
1	A	236	LEU	2.3
1	A	45	VAL	2.3
2	G	1287	ILE	2.2
1	A	62	ARG	2.2
1	A	199	SER	2.2
1	A	205	GLU	2.2
2	G	103	LEU	2.1

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
1	HIC	A	73	11/12	0.96	0.17	-	28,62,88,88	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
4	CA	A	402	1/1	0.99	0.21	1.53	37,37,37,37	0
4	CA	A	403	1/1	0.99	0.19	0.97	41,41,41,41	0
3	ATP	A	401	31/31	0.98	0.17	0.96	32,41,53,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	CA	G	1401	1/1	0.99	0.17	0.95	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.