



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 05:12 PM UTC

PDB ID : 4U0S / pdb_00004u0s
Title : Structure of Eukaryotic fic domain containing protein with ADP
Authors : Cole, A.R.; Katan, M.; Bunney, T.D.
Deposited on : 2014-07-14
Resolution : 2.49 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

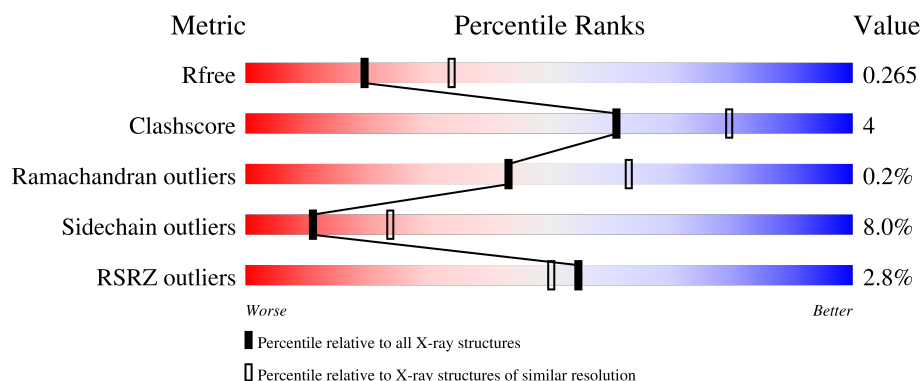
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

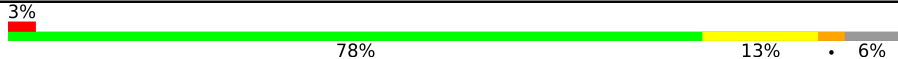
The reported resolution of this entry is 2.49 Å.

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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	344	
1	B	344	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5310 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.

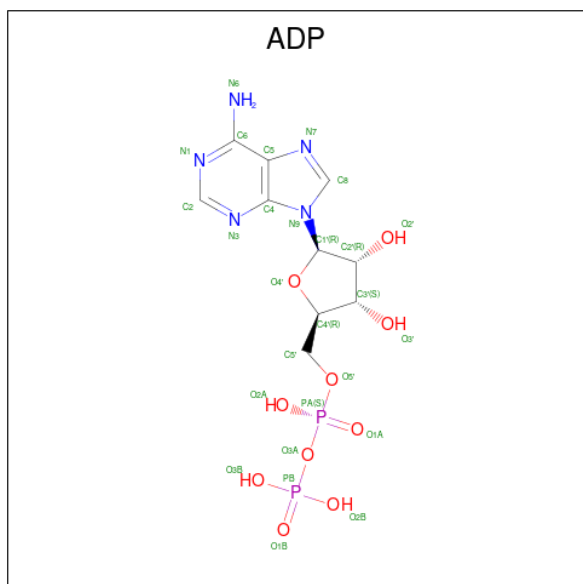
- Molecule 1 is a protein called Adenosine monophosphate-protein transferase FICD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	0	0
			2532	1620	434	467	11			
1	B	324	Total	C	N	O	S	0	0	0
			2527	1615	434	466	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	234	GLY	GLU	engineered mutation	UNP Q9BVA6
B	234	GLY	GLU	engineered mutation	UNP Q9BVA6

- 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		

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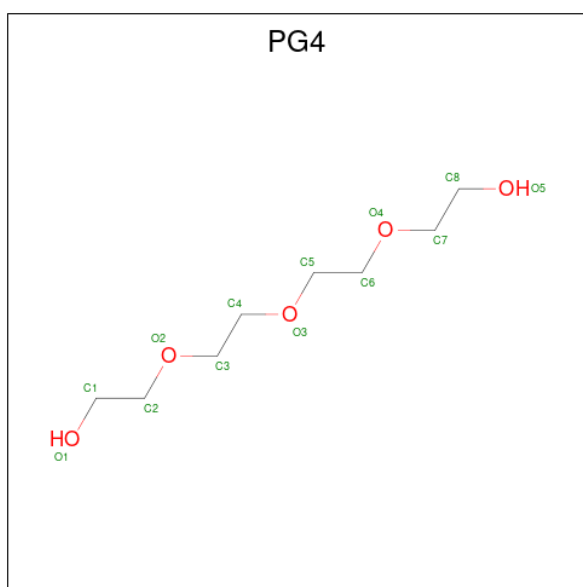
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

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

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

- Molecule 4 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	8	5		
4	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	113	Total	O	0	0
			113	113		
5	B	56	Total	O	0	0
			56	56		

- Molecule 1: Adenosine monophosphate-protein transferase FICD



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.04Å 76.01Å 92.03Å 90.00° 107.56° 90.00°	Depositor
Resolution (Å)	31.86 – 2.49 31.86 – 2.49	Depositor EDS
% Data completeness (in resolution range)	98.8 (31.86-2.49) 98.7 (31.86-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 2.48Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.207 , 0.245 0.227 , 0.265	Depositor DCC
R_{free} test set	1647 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	42.2	Xtriage
Anisotropy	0.511	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 76.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5310	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.34% 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: ADP, PG4, MG

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	0.88	0/2586	1.57	19/3515 (0.5%)
1	B	0.89	0/2581	1.58	15/3510 (0.4%)
All	All	0.89	0/5167	1.57	34/7025 (0.5%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	427	ASP	N-CA-C	-9.59	98.16	110.53
1	A	369	ASN	CA-C-N	6.72	127.39	120.00
1	A	369	ASN	C-N-CA	6.72	127.39	120.00
1	A	185	PRO	CA-C-N	6.36	129.08	120.44
1	A	185	PRO	C-N-CA	6.36	129.08	120.44
1	A	263	GLU	CA-C-N	6.34	129.43	120.42
1	A	263	GLU	C-N-CA	6.34	129.43	120.42
1	B	263	GLU	CA-C-N	6.12	129.11	120.42
1	B	263	GLU	C-N-CA	6.12	129.11	120.42
1	A	391	ILE	N-CA-C	-5.68	99.05	107.51
1	B	427	ASP	CA-C-O	-5.63	115.47	121.94
1	A	139	PHE	CA-CB-CG	5.61	119.41	113.80
1	B	427	ASP	CA-CB-CG	5.60	118.20	112.60
1	A	422	THR	N-CA-CB	5.59	118.43	110.16
1	B	188	GLU	CA-C-N	5.50	127.92	120.44
1	B	188	GLU	C-N-CA	5.50	127.92	120.44
1	A	355	ALA	CA-C-N	5.49	127.58	120.44
1	A	355	ALA	C-N-CA	5.49	127.58	120.44
1	B	139	PHE	CA-CB-CG	5.46	119.26	113.80
1	B	422	THR	N-CA-CB	5.46	118.23	110.16
1	B	379	LEU	N-CA-C	-5.45	105.25	111.14
1	A	398	ASP	CA-C-N	5.37	127.79	120.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	398	ASP	C-N-CA	5.37	127.79	120.54
1	A	146	PHE	CA-C-N	5.25	125.77	119.94
1	A	146	PHE	C-N-CA	5.25	125.77	119.94
1	B	302	ASP	CA-CB-CG	5.20	117.80	112.60
1	B	398	ASP	CA-C-N	5.14	127.48	120.54
1	B	398	ASP	C-N-CA	5.14	127.48	120.54
1	A	161	TYR	CA-C-N	5.10	127.11	120.28
1	A	161	TYR	C-N-CA	5.10	127.11	120.28
1	A	302	ASP	CA-CB-CG	5.09	117.69	112.60
1	A	278	LEU	N-CA-C	5.09	119.48	113.12
1	B	430	LEU	CA-C-N	5.06	131.21	121.54
1	B	430	LEU	C-N-CA	5.06	131.21	121.54

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	2532	0	2457	22	0
1	B	2527	0	2444	23	0
2	A	27	0	12	0	0
2	B	27	0	12	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	13	0	18	3	0
4	B	13	0	18	3	0
5	A	113	0	0	0	0
5	B	56	0	0	0	0
All	All	5310	0	4961	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:254:PRO:HB3	1:B:254:PRO:HB3	1.63	0.79
1:A:258:LEU:HD13	1:B:265:ILE:HD11	1.81	0.61
1:B:389:ILE:HG21	1:B:422:THR:HB	1.83	0.60
1:A:265:ILE:HD11	1:B:258:LEU:HD13	1.84	0.60
1:A:389:ILE:HG21	1:A:422:THR:HB	1.82	0.60
1:A:310:ARG:HG2	1:A:366:ILE:HD13	1.86	0.58
1:A:406:ALA:HB1	1:A:411:VAL:HG12	1.86	0.57
1:B:230:THR:HG21	1:B:267:MET:HE2	1.86	0.57
1:A:230:THR:HG21	1:A:267:MET:HE2	1.87	0.56
1:B:310:ARG:HG2	1:B:366:ILE:HD13	1.87	0.55
1:B:184:LEU:HB3	1:B:185:PRO:HD3	1.89	0.54
1:A:130:MET:HA	1:A:130:MET:HE2	1.89	0.54
1:A:358:LYS:HB3	4:A:503:PG4:H21	1.91	0.53
1:B:358:LYS:HE3	4:B:503:PG4:H61	1.90	0.53
1:B:184:LEU:HB3	1:B:185:PRO:CD	2.38	0.53
1:B:356:HIS:C	1:B:356:HIS:CD2	2.87	0.52
1:A:184:LEU:HB3	1:A:185:PRO:HD3	1.93	0.51
1:A:113:ALA:HB2	1:A:128:LEU:HB3	1.92	0.51
1:B:430:LEU:O	1:B:431:PHE:HB3	2.11	0.49
1:B:358:LYS:HG2	4:B:503:PG4:H31	1.96	0.48
1:A:113:ALA:HA	1:A:128:LEU:HD13	1.97	0.47
1:A:222:THR:HA	1:A:225:HIS:HB3	1.99	0.45
1:A:110:LEU:HD13	1:A:142:ALA:HA	1.98	0.45
1:A:358:LYS:NZ	4:A:503:PG4:O1	2.49	0.45
1:B:161:TYR:HD1	1:B:410:ASP:OD2	2.00	0.45
1:B:358:LYS:HG2	4:B:503:PG4:H21	1.98	0.44
1:B:382:MET:SD	1:B:388:PRO:HG3	2.58	0.44
1:B:360:VAL:HG13	2:B:501:ADP:N3	2.33	0.43
1:B:110:LEU:HD13	1:B:142:ALA:HA	1.99	0.43
1:A:298:LEU:HD13	1:A:302:ASP:HB3	2.00	0.43
1:A:382:MET:SD	1:A:388:PRO:HG3	2.58	0.43
1:B:238:LEU:HD23	1:B:256:LYS:HG2	2.01	0.43
1:B:139:PHE:CD2	1:B:142:ALA:HB2	2.54	0.43
1:B:222:THR:HA	1:B:225:HIS:HB3	2.01	0.43
1:A:238:LEU:HD23	1:A:256:LYS:HG2	2.00	0.42
1:A:139:PHE:CD2	1:A:142:ALA:HB2	2.54	0.42
1:B:298:LEU:HD13	1:B:302:ASP:HB3	2.02	0.42
1:B:205:VAL:HG23	1:B:430:LEU:HD22	2.02	0.42
1:A:122:ARG:NH2	1:A:152:GLU:OE1	2.53	0.42
1:B:122:ARG:NH1	1:B:149:PHE:CE1	2.88	0.42
1:A:175:LYS:O	1:A:179:ASN:ND2	2.33	0.41
1:A:358:LYS:HG2	4:A:503:PG4:H31	2.04	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	319/344 (93%)	309 (97%)	10 (3%)	0	100	100
1	B	320/344 (93%)	308 (96%)	11 (3%)	1 (0%)	36	55
All	All	639/688 (93%)	617 (97%)	21 (3%)	1 (0%)	43	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	431	PHE

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	262/301 (87%)	239 (91%)	23 (9%)	9	20
1	B	261/301 (87%)	242 (93%)	19 (7%)	13	27
All	All	523/602 (87%)	481 (92%)	42 (8%)	11	24

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	LEU
1	A	122	ARG

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Mol	Chain	Res	Type
1	A	123	GLU
1	A	126	GLN
1	A	128	LEU
1	A	130	MET
1	A	135	MET
1	A	177	LEU
1	A	208	ILE
1	A	218	VAL
1	A	230	THR
1	A	240	LEU
1	A	241	SER
1	A	256	LYS
1	A	271	MET
1	A	282	ILE
1	A	287	ILE
1	A	298	LEU
1	A	316	VAL
1	A	342	GLU
1	A	360	VAL
1	A	363	HIS
1	A	422	THR
1	B	135	MET
1	B	177	LEU
1	B	208	ILE
1	B	220	GLU
1	B	240	LEU
1	B	241	SER
1	B	256	LYS
1	B	271	MET
1	B	282	ILE
1	B	287	ILE
1	B	298	LEU
1	B	316	VAL
1	B	342	GLU
1	B	360	VAL
1	B	363	HIS
1	B	422	THR
1	B	430	LEU
1	B	431	PHE
1	B	434	THR

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	383	GLN
1	B	226	HIS
1	B	275	ASN
1	B	383	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\# Z > 2$
4	PG4	A	503	-	12,12,12	0.70	0	11,11,11	0.50	0
2	ADP	B	501	3	28,29,29	0.42	0	43,45,45	0.73	1 (2%)
2	ADP	A	501	3	28,29,29	0.61	1 (3%)	43,45,45	0.77	1 (2%)
4	PG4	B	503	-	12,12,12	0.64	0	11,11,11	0.29	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
4	PG4	A	503	-	-	4/10/10/10	-
2	ADP	B	501	3	-	8/16/32/32	0/3/3/3
2	ADP	A	501	3	-	1/16/32/32	0/3/3/3
4	PG4	B	503	-	-	7/10/10/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	ADP	PB-O3B	-2.03	1.47	1.54

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	ADP	O2A-PA-O5'	2.28	117.90	107.57
2	B	501	ADP	O2'-C2'-C3'	-2.17	104.85	111.82

There are no chirality outliers.

All (20) torsion outliers are listed below:

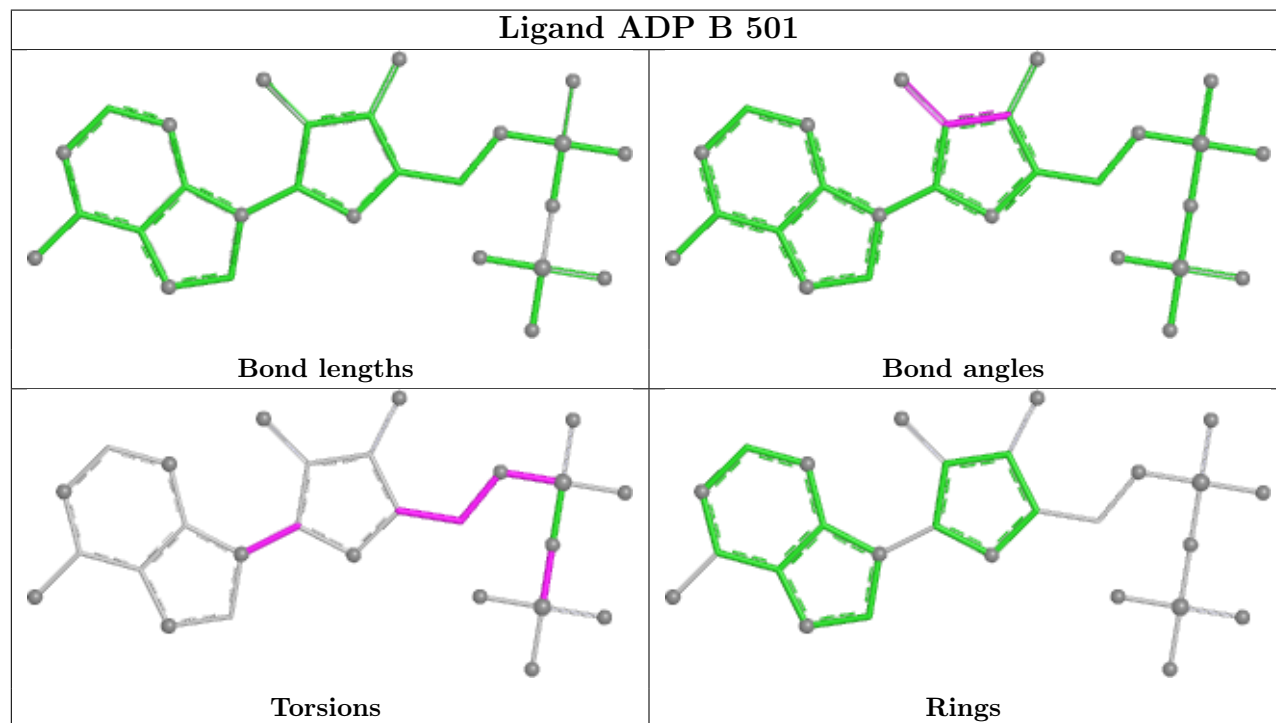
Mol	Chain	Res	Type	Atoms
2	B	501	ADP	PA-O3A-PB-O2B
2	B	501	ADP	C5'-O5'-PA-O1A
4	A	503	PG4	O2-C3-C4-O3
4	B	503	PG4	O2-C3-C4-O3
4	A	503	PG4	O3-C5-C6-O4
2	B	501	ADP	C3'-C4'-C5'-O5'
4	B	503	PG4	O3-C5-C6-O4
4	B	503	PG4	O4-C7-C8-O5
2	B	501	ADP	O4'-C4'-C5'-O5'
2	B	501	ADP	PA-O3A-PB-O3B
4	A	503	PG4	C5-C6-O4-C7
4	B	503	PG4	C6-C5-O3-C4
4	B	503	PG4	C1-C2-O2-C3
4	B	503	PG4	C8-C7-O4-C6
4	B	503	PG4	O1-C1-C2-O2
2	B	501	ADP	C5'-O5'-PA-O3A
2	B	501	ADP	C4'-C5'-O5'-PA
4	A	503	PG4	C3-C4-O3-C5
2	B	501	ADP	C2'-C1'-N9-C8
2	A	501	ADP	C2'-C1'-N9-C8

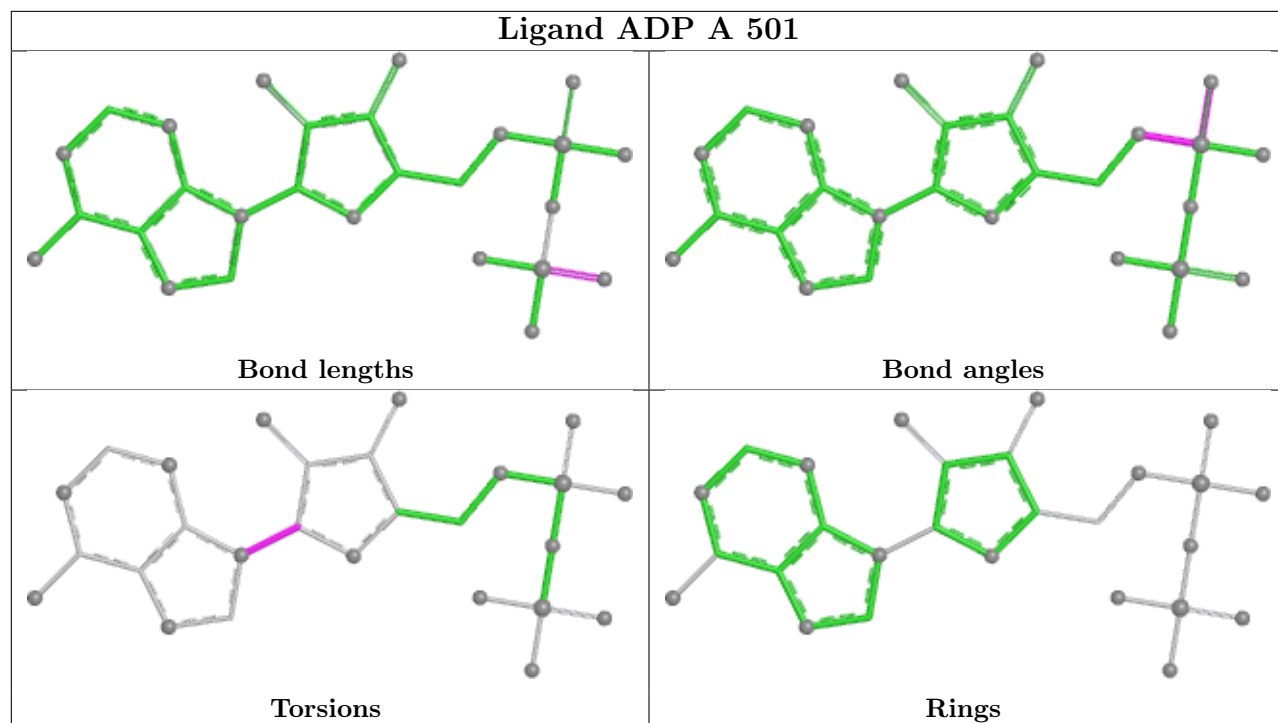
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	503	PG4	3	0
2	B	501	ADP	1	0
4	B	503	PG4	3	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	323/344 (93%)	0.20	10 (3%) 51 47	29, 50, 110, 128	0
1	B	324/344 (94%)	0.51	8 (2%) 58 54	37, 66, 111, 123	0
All	All	647/688 (94%)	0.35	18 (2%) 55 50	29, 60, 111, 128	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	108	ALA	4.5
1	A	108	ALA	4.5
1	B	205	VAL	3.8
1	A	214	ALA	3.7
1	A	109	ALA	3.6
1	B	109	ALA	3.4
1	A	208	ILE	3.3
1	B	140	VAL	3.2
1	B	112	GLN	2.9
1	A	215	LEU	2.8
1	A	137	PRO	2.6
1	A	213	SER	2.5
1	B	123	GLU	2.4
1	B	208	ILE	2.2
1	A	433	THR	2.2
1	B	352	ALA	2.1
1	A	212	ASN	2.1
1	A	165	ARG	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](#)

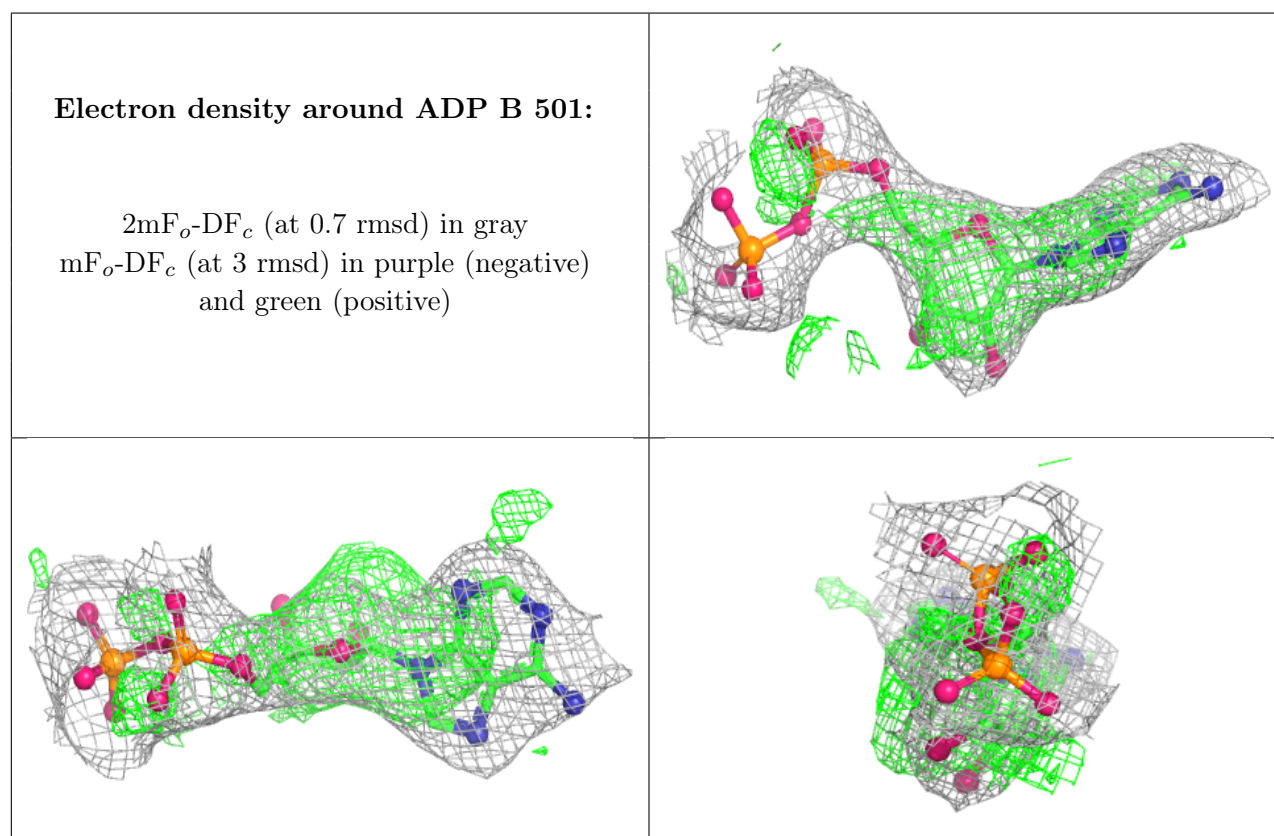
There are no oligosaccharides in this entry.

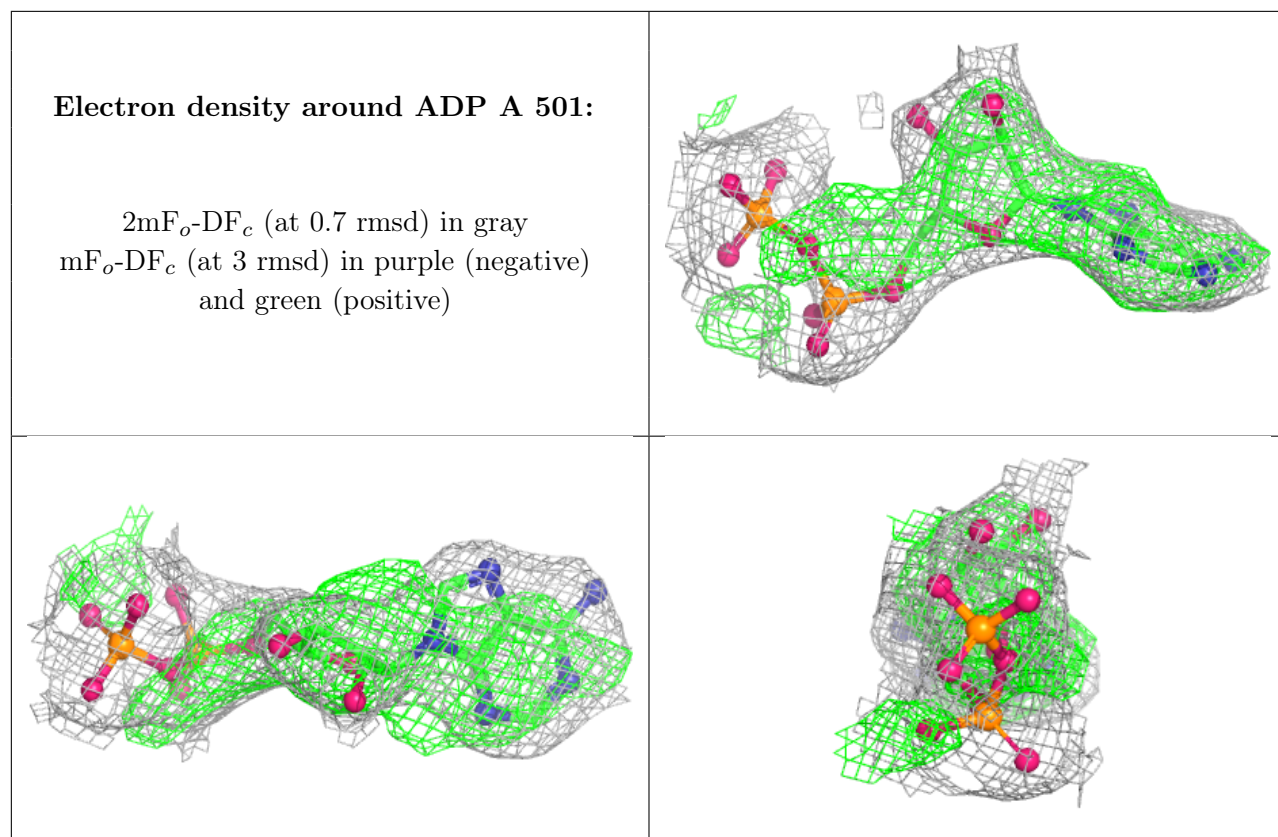
6.4 Ligands [i](#)

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	PG4	B	503	13/13	0.78	0.17	77,79,81,81	0
4	PG4	A	503	13/13	0.86	0.13	57,60,65,66	0
2	ADP	B	501	27/27	0.92	0.14	15,26,43,54	21
2	ADP	A	501	27/27	0.95	0.11	3,11,32,39	21
3	MG	B	502	1/1	0.95	0.11	49,49,49,49	0
3	MG	A	502	1/1	0.97	0.15	39,39,39,39	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.