



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 11:32 am BST

PDB ID : 5FHD
Title : Structure of Bacteroides sp Pif1 complexed with tailed dsDNA resulting in ssDNA bound complex
Authors : Zhou, X.; Ren, W.; Bharath, S.R.; Song, H.
Deposited on : 2015-12-22
Resolution : 2.00 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

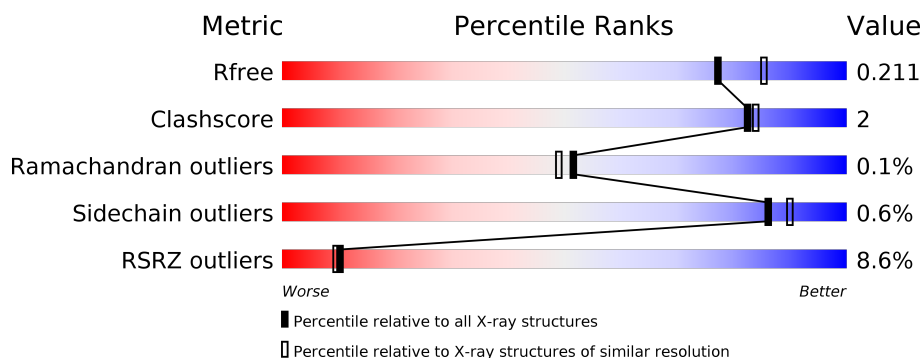
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 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	433	<div> <div>3%</div> <div>90%</div> <div>6%</div> <div>.</div> </div>
1	B	433	<div> <div>13%</div> <div>86%</div> <div>6%</div> <div>8%</div> </div>
2	C	19	<div> <div>5%</div> <div>21%</div> <div>21%</div> <div>58%</div> </div>
2	E	19	<div> <div>32%</div> <div>5%</div> <div>63%</div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7291 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.

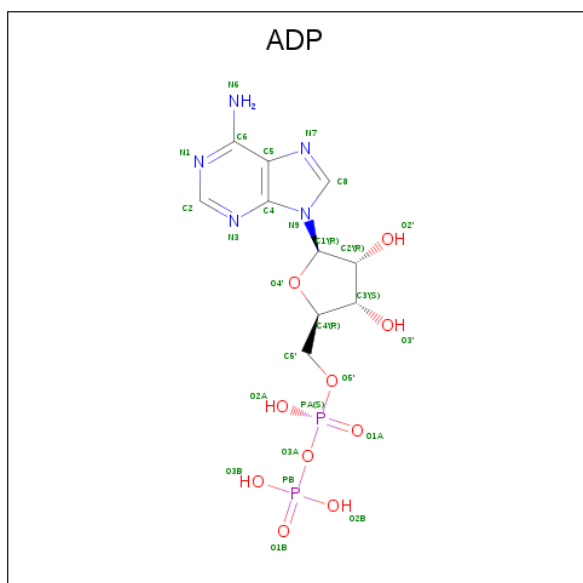
- Molecule 1 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	414	Total	C	N	O	S	0	6	0
			3368	2171	555	624	18			
1	B	400	Total	C	N	O	S	0	1	0
			3114	2006	509	584	15			

- Molecule 2 is a DNA chain called DNA (5'-D(*TP*TP*TP*TP*TP*TP*TP*CP*CP*GP*GP*GP*GP*CP*CP*GP*CP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	8	Total	C	N	O	P	0	0	0
			160	78	21	53	8			
2	E	7	Total	C	N	O	P	0	0	0
			139	69	15	48	7			

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).

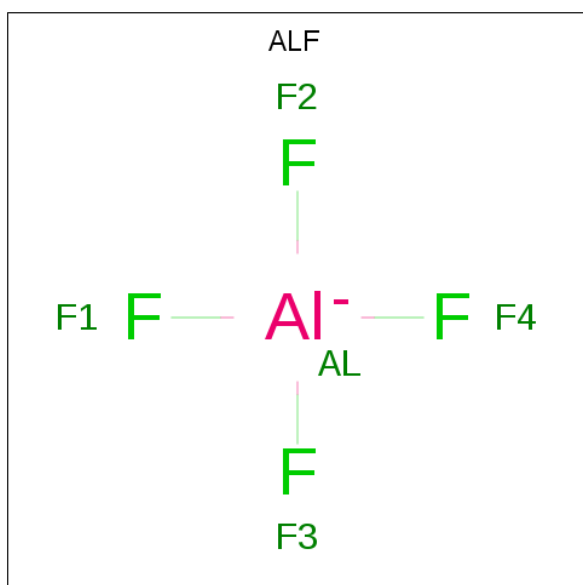


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 5 is TETRAFLUOROALUMINATE ION (three-letter code: ALF) (formula: AlF_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	Al	F	0	0
			5	1	4		
5	B	1	Total	Al	F	0	0
			5	1	4		

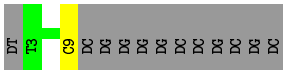
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	235	Total	O	0	0
			235	235		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	13	Total 13	O 13	0	0
6	E	5	Total 5	O 5	0	0
6	B	191	Total 191	O 191	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	48.47Å 149.64Å 185.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.54 – 2.00 29.54 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.1 (29.54-2.00) 99.1 (29.54-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.17 (at 2.00Å)	Xtriage
Refinement program	PHENIX (1.10.1 _2155: ???)	Depositor
R, R_{free}	0.178 , 0.211 0.178 , 0.211	Depositor DCC
R_{free} test set	4546 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7291	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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: ALF, MG, ADP

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	0.64	1/3451 (0.0%)	0.69	0/4660
1	B	0.64	1/3174 (0.0%)	0.70	3/4295 (0.1%)
2	C	1.39	0/176	1.21	0/269
2	E	1.21	0/152	1.35	0/232
All	All	0.69	2/6953 (0.0%)	0.74	3/9456 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	350	PHE	C-N	9.66	1.52	1.34
1	A	411	GLU	CD-OE2	6.16	1.32	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	277	LEU	CA-CB-CG	9.26	136.59	115.30
1	B	175	ARG	NE-CZ-NH1	5.35	122.98	120.30
1	B	175	ARG	NE-CZ-NH2	-5.24	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	3368	0	3405	15	0
1	B	3114	0	3027	17	0
2	C	160	0	94	2	0
2	E	139	0	84	1	0
3	A	27	0	12	0	0
3	B	27	0	12	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
6	A	235	0	0	2	0
6	B	191	0	0	1	0
6	C	13	0	0	0	0
6	E	5	0	0	0	0
All	All	7291	0	6634	33	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 2.

All (33) 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:277:LEU:HD12	1:B:278:ARG:H	1.39	0.83
1:B:277:LEU:HD12	1:B:278:ARG:N	2.02	0.74
1:A:83:ARG:NH2	6:A:602:HOH:O	2.24	0.70
1:A:111[B]:ARG:HD3	1:A:150:THR:HG22	1.82	0.62
1:A:254:ARG:NH2	1:A:307:ASP:OD1	2.39	0.55
1:A:294:TYR:HB3	1:A:320:ILE:HD13	1.88	0.53
2:E:9:DC:H2'	1:B:91:TYR:CD2	2.43	0.53
1:A:111[A]:ARG:HD3	1:A:150:THR:HG22	1.91	0.52
1:B:14:MET:HE1	1:B:40:TYR:HE2	1.75	0.52
1:A:57:ILE:HD12	1:A:362:LYS:HB3	1.92	0.52
1:A:318:ARG:HG3	1:B:226:TYR:CZ	2.46	0.51
1:B:304:ALA:HB3	1:B:311:THR:HG22	1.92	0.51
1:B:153:GLN:CD	1:B:153:GLN:H	2.14	0.50
1:B:152:LYS:HB3	1:B:153:GLN:NE2	2.27	0.49
1:B:277:LEU:CD1	1:B:278:ARG:H	2.18	0.46
1:B:95:LEU:HD12	6:B:638:HOH:O	2.15	0.46
1:A:3:ASP:N	1:A:3:ASP:OD1	2.49	0.45
1:B:238:ALA:O	1:B:242:LYS:HD3	2.17	0.45
2:C:7:DT:H2'	2:C:8:DC:C6	2.51	0.45
1:A:288:ASN:HA	1:A:295:TYR:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:TYR:OH	6:A:601:HOH:O	2.20	0.44
1:B:269:ILE:HG22	1:B:271:CYS:H	1.83	0.44
1:B:325:TYR:O	1:B:348:THR:HA	2.19	0.43
1:A:60:ILE:HD11	1:A:295:TYR:HB2	2.01	0.42
1:A:295:TYR:N	1:A:298:MET:HE3	2.34	0.42
1:B:52:ALA:HA	1:B:65:VAL:O	2.20	0.41
1:A:147:PRO:HB3	1:A:166:PHE:CE1	2.55	0.41
2:C:4:DT:H73	2:C:5:DT:H73	2.02	0.41
1:B:295:TYR:H	1:B:298:MET:HE3	1.86	0.41
1:B:277:LEU:HD12	1:B:278:ARG:C	2.41	0.41
1:A:295:TYR:H	1:A:298:MET:HE3	1.86	0.41
1:A:269:ILE:HG22	1:A:271:CYS:H	1.87	0.40
1:B:315:ASP:O	1:B:316:ASN:C	2.58	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	414/433 (96%)	408 (99%)	5 (1%)	1 (0%)	47	44
1	B	387/433 (89%)	373 (96%)	14 (4%)	0	100	100
All	All	801/866 (92%)	781 (98%)	19 (2%)	1 (0%)	51	49

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	262	ASP

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	376/390 (96%)	374 (100%)	2 (0%)	88	92
1	B	330/390 (85%)	328 (99%)	2 (1%)	86	90
All	All	706/780 (90%)	702 (99%)	4 (1%)	86	90

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	264	PHE
1	A	419	ARG
1	B	277	LEU
1	B	419	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	153	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 carbohydrates in this entry.

5.6 Ligand geometry

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
3	ADP	A	501	4	24,29,29	1.25	4 (16%)	29,45,45	1.29	5 (17%)
5	ALF	A	503	-	0,4,4	0.00	-	-		
3	ADP	B	501	4	24,29,29	1.05	2 (8%)	29,45,45	1.41	4 (13%)
5	ALF	B	503	-	0,4,4	0.00	-	-		

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	ADP	A	501	4	-	3/12/32/32	0/3/3/3
3	ADP	B	501	4	-	3/12/32/32	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	ADP	PB-O3B	-2.58	1.44	1.54
3	B	501	ADP	C5-C4	2.51	1.47	1.40
3	A	501	ADP	O4'-C1'	2.30	1.44	1.41
3	A	501	ADP	C5-C4	2.24	1.46	1.40
3	A	501	ADP	C2-N3	2.23	1.35	1.32
3	B	501	ADP	C2-N3	2.07	1.35	1.32

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	ADP	O2B-PB-O1B	3.48	124.29	110.68
3	B	501	ADP	N3-C2-N1	-3.41	123.34	128.68
3	A	501	ADP	N3-C2-N1	-2.58	124.65	128.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	ADP	C4-C5-N7	-2.36	106.94	109.40
3	B	501	ADP	C4-C5-N7	-2.34	106.96	109.40
3	A	501	ADP	C1'-N9-C4	-2.32	122.57	126.64
3	A	501	ADP	O3B-PB-O2B	2.27	116.31	107.64
3	B	501	ADP	PA-O3A-PB	-2.12	125.56	132.83
3	A	501	ADP	O2'-C2'-C1'	-2.04	103.32	110.85

There are no chirality outliers.

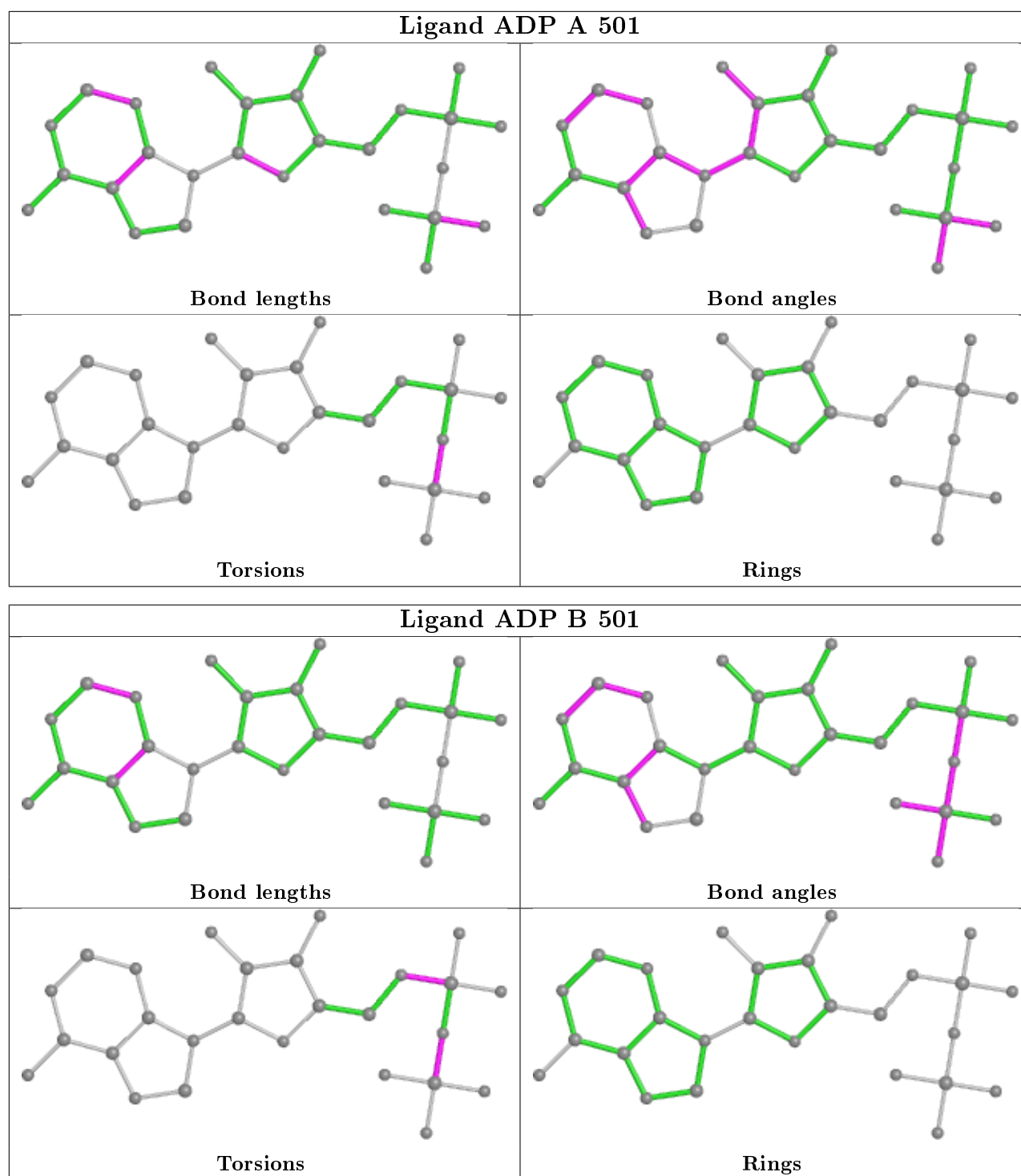
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	501	ADP	PA-O3A-PB-O2B
3	A	501	ADP	PA-O3A-PB-O3B
3	B	501	ADP	PA-O3A-PB-O1B
3	A	501	ADP	PA-O3A-PB-O2B
3	A	501	ADP	PA-O3A-PB-O1B
3	B	501	ADP	C5'-O5'-PA-O3A

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	414/433 (95%)	-0.00	15 (3%) 42 42	17, 30, 62, 94	0
1	B	400/433 (92%)	0.43	55 (13%) 2 2	17, 32, 81, 95	0
2	C	8/19 (42%)	0.63	1 (12%) 3 3	29, 41, 60, 78	0
2	E	7/19 (36%)	0.11	0 100 100	30, 34, 64, 76	0
All	All	829/904 (91%)	0.21	71 (8%) 10 9	17, 31, 76, 95	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	327	TRP	10.7
1	B	258	ILE	7.8
1	B	347	CYS	7.4
1	B	259	VAL	6.6
1	B	316	ASN	6.5
1	B	326	THR	6.5
1	B	346	SER	6.4
1	B	303	THR	5.4
1	B	320	ILE	5.3
2	C	10	DG	5.2
1	B	325	TYR	5.1
1	B	257	ASP	5.1
1	B	318	ARG	5.1
1	B	290	SER	5.0
1	B	328	SER	5.0
1	B	272	ASP	5.0
1	B	305	LEU	4.9
1	B	322	PHE	4.9
1	B	348	THR	4.9
1	B	317	GLY	4.9
1	B	291	LEU	4.8

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Mol	Chain	Res	Type	RSRZ
1	B	294	TYR	4.6
1	B	271	CYS	4.5
1	B	268	SER	4.5
1	B	269	ILE	4.4
1	B	311	THR	4.4
1	B	81	TYR	4.4
1	B	319	THR	4.3
1	B	1	MET	4.1
1	A	330	THR	3.7
1	B	302	VAL	3.7
1	B	312	VAL	3.7
1	B	254	ARG	3.6
1	B	301	ILE	3.5
1	B	270	PRO	3.4
1	B	253	ILE	3.4
1	A	221	LYS	3.4
1	B	255	ASN	3.3
1	A	344	ILE	3.2
1	B	288	ASN	3.2
1	A	342	GLU	3.1
1	A	345	GLY	3.0
1	B	281	ALA	3.0
1	B	350	PHE	2.8
1	A	262	ASP	2.7
1	B	256	TYR	2.7
1	B	277	LEU	2.7
1	B	300	GLY	2.7
1	B	226	TYR	2.7
1	A	343	GLU	2.6
1	A	251	GLN	2.6
1	A	259	VAL	2.6
1	B	349	GLN	2.6
1	A	3	ASP	2.6
1	A	222	ILE	2.5
1	B	315	ASP	2.5
1	B	107	ILE	2.5
1	B	91	TYR	2.4
1	B	2	GLU	2.4
1	A	260	ILE	2.3
1	B	308	ASN	2.3
1	B	252	GLU	2.2
1	B	295	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	321	LYS	2.1
1	A	107	ILE	2.1
1	B	94	GLU	2.1
1	A	153	GLN	2.1
1	A	250	GLU	2.1
1	B	351	PRO	2.1
1	B	289	ASP	2.1
1	B	249	GLY	2.1

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 carbohydrates in this entry.

6.4 Ligands [i](#)

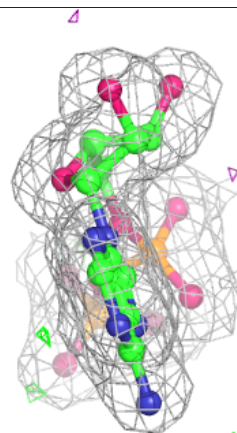
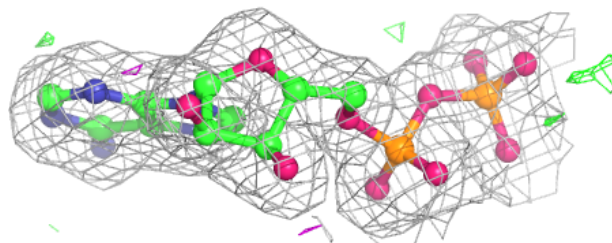
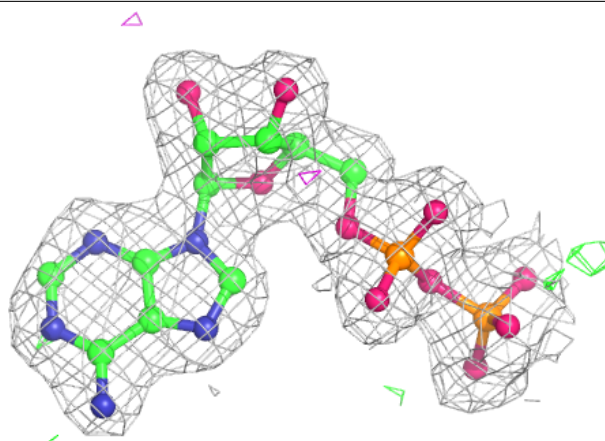
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	MG	B	502	1/1	0.98	0.10	21,21,21,21	0
3	ADP	A	501	27/27	0.98	0.09	15,24,30,38	0
3	ADP	B	501	27/27	0.99	0.08	16,24,29,30	0
4	MG	A	502	1/1	0.99	0.14	20,20,20,20	0
5	ALF	A	503	5/5	0.99	0.13	17,18,19,20	0
5	ALF	B	503	5/5	0.99	0.12	17,18,21,21	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.

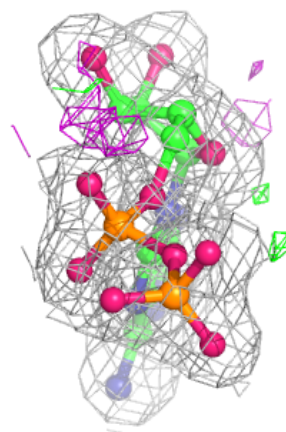
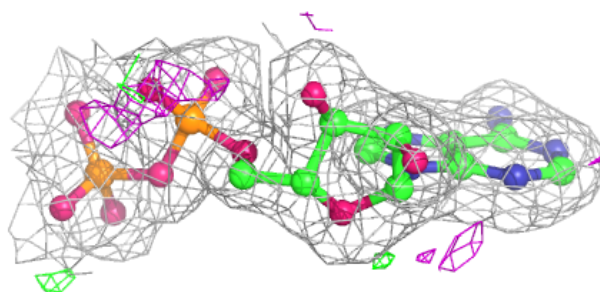
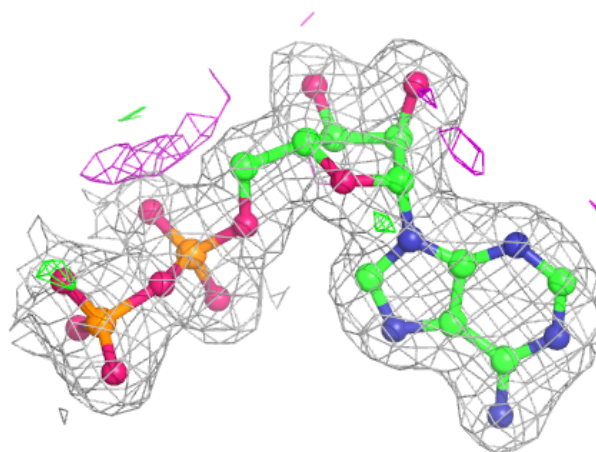
Electron density around ADP A 501:

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



Electron density around ADP B 501:

$2mF_o-DF_c$ (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.