



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 18, 2020 – 11:38 pm BST

PDB ID : 2DFT  
Title : Structure of shikimate kinase from Mycobacterium tuberculosis complexed with ADP and Mg at 2.8 angstroms of resolution  
Authors : Dias, M.V.; Faim, L.M.; Vasconcelos, I.B.; de Oliveira, J.S.; Basso, L.A.; Santos, D.S.; de Azevedo, W.F.  
Deposited on : 2006-03-03  
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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

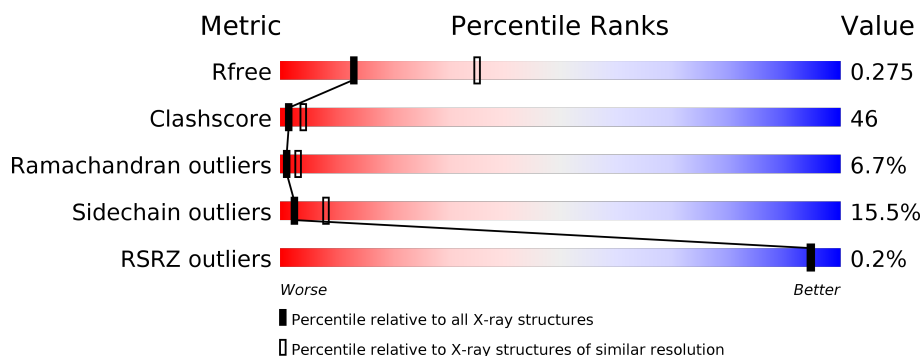
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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  $\geq 3$ , 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	176	<div> <div>33%</div> <div>36%</div> <div>16%</div> <div>•</div> <div>11%</div> </div>
1	B	176	<div> <div>35%</div> <div>37%</div> <div>16%</div> <div>6%</div> <div>6%</div> </div>
1	C	176	<div> <div>24%</div> <div>41%</div> <div>16%</div> <div>6%</div> <div>12%</div> </div>
1	D	176	<div> <div>%</div> <div>32%</div> <div>39%</div> <div>16%</div> <div>7%</div> <div>6%</div> </div>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5152 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 Shikimate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	0	0
			1165	717	232	214	2			
1	B	165	Total	C	N	O	S	0	0	0
			1229	759	244	224	2			
1	C	155	Total	C	N	O	S	0	0	0
			1157	713	230	212	2			
1	D	165	Total	C	N	O	S	0	0	0
			1226	757	244	223	2			

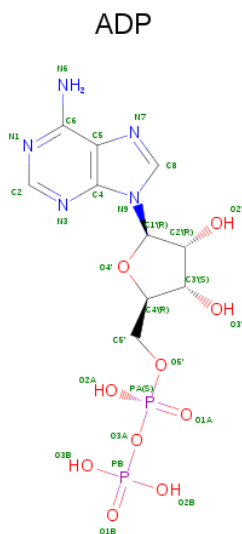
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		
2	D	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>10</sub>P<sub>2</sub>).



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

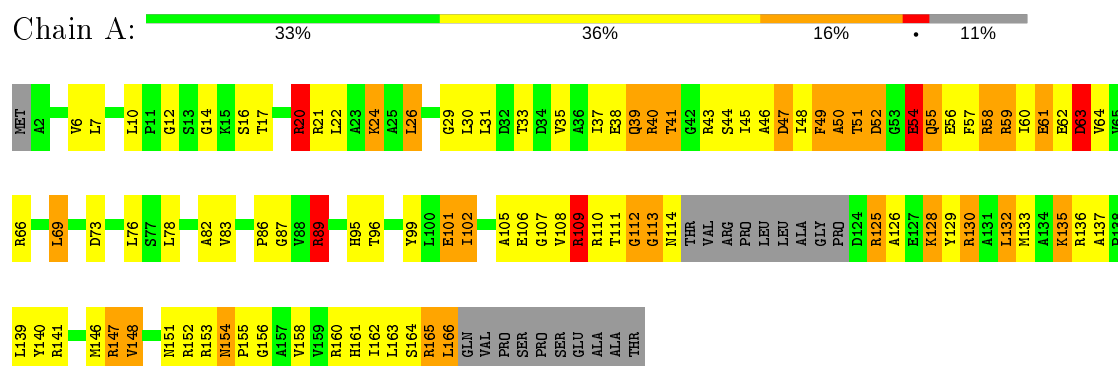
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	60	Total O 60 60	0	0
5	B	85	Total O 85 85	0	0
5	C	51	Total O 51 51	0	0
5	D	64	Total O 64 64	0	0

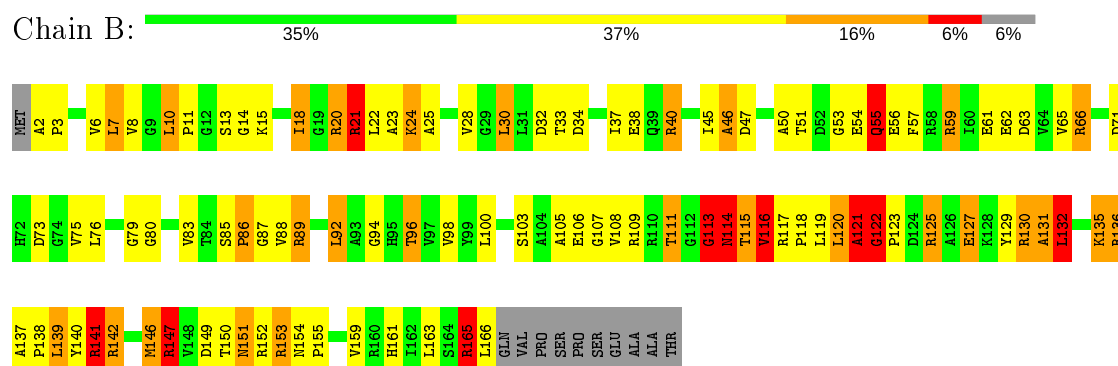
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

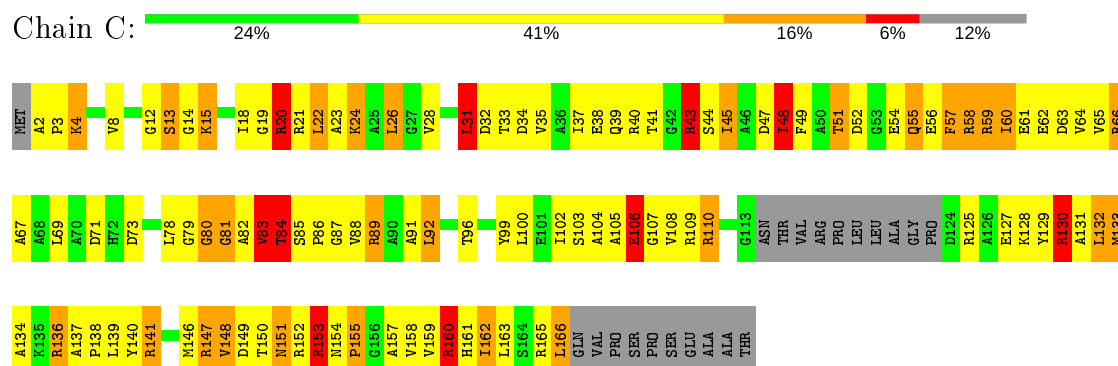
#### • Molecule 1: Shikimate kinase



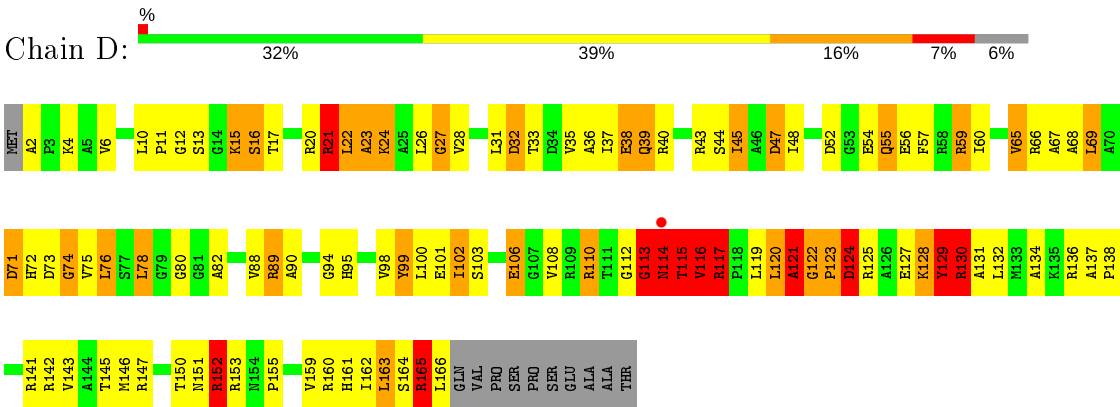
#### • Molecule 1: Shikimate kinase



#### • Molecule 1: Shikimate kinase



● Molecule 1: Shikimate kinase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	60.62Å 62.20Å 170.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	57.17 – 2.80 57.12 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (57.17-2.80) 99.7 (57.12-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.56 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.183 , 0.282 0.184 , 0.275	Depositor DCC
$R_{free}$ test set	832 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	45.0	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 47.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.028 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5152	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ADP, CL

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	1.37	2/1176 (0.2%)	2.04	38/1584 (2.4%)
1	B	1.33	5/1243 (0.4%)	1.92	33/1679 (2.0%)
1	C	1.44	7/1168 (0.6%)	2.03	35/1573 (2.2%)
1	D	1.39	6/1240 (0.5%)	2.01	43/1674 (2.6%)
All	All	1.38	20/4827 (0.4%)	2.00	149/6510 (2.3%)

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	3
1	B	0	4
1	C	0	4
1	D	0	6
All	All	0	17

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	2	ALA	CA-CB	7.69	1.68	1.52
1	B	83	VAL	CB-CG1	-6.63	1.39	1.52
1	B	127	GLU	CG-CD	6.31	1.61	1.51
1	D	116	VAL	CB-CG2	6.00	1.65	1.52
1	D	129	TYR	CB-CG	-5.91	1.42	1.51

The worst 5 of 149 bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	125	ARG	NE-CZ-NH1	-15.45	112.58	120.30
1	C	20	ARG	NE-CZ-NH2	-13.97	113.31	120.30
1	A	20	ARG	NE-CZ-NH2	-13.57	113.52	120.30
1	D	21	ARG	NE-CZ-NH2	-12.74	113.93	120.30
1	D	130	ARG	NE-CZ-NH1	12.73	126.66	120.30

There are no chirality outliers.

5 of 17 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	109	ARG	Mainchain
1	A	111	THR	Peptide
1	A	113	GLY	Peptide
1	B	113	GLY	Peptide
1	B	114	ASN	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	1165	0	1208	110	0
1	B	1229	0	1282	114	1
1	C	1157	0	1202	123	1
1	D	1226	0	1275	110	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
4	A	27	0	12	6	0
4	B	27	0	12	2	0
4	C	27	0	12	8	0
4	D	27	0	12	6	0
5	A	60	0	0	31	2
5	B	85	0	0	47	1
5	C	51	0	0	38	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	64	0	0	33	1
All	All	5152	0	5015	457	4

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 46.

The worst 5 of 457 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:165:ARG:HH11	1:B:165:ARG:CG	1.34	1.37
1:A:59:ARG:HB2	5:A:278:HOH:O	1.15	1.31
1:C:157:ALA:HB3	5:C:411:HOH:O	1.25	1.29
1:A:125:ARG:HA	5:A:387:HOH:O	1.21	1.28
1:B:116:VAL:CG1	5:B:373:HOH:O	1.66	1.25

All (4) 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)
5:A:436:HOH:O	5:B:441:HOH:O[4_456]	1.71	0.49
1:B:135:LYS:CD	5:A:436:HOH:O[4_556]	1.93	0.27
1:C:96:THR:OG1	1:D:130:ARG:NH2[3_655]	2.04	0.16
5:C:287:HOH:O	5:D:286:HOH:O[3_555]	2.07	0.13

## 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	152/176 (86%)	132 (87%)	14 (9%)	6 (4%)	3 10
1	B	163/176 (93%)	133 (82%)	17 (10%)	13 (8%)	1 2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	151/176 (86%)	129 (85%)	14 (9%)	8 (5%)	2	6
1	D	163/176 (93%)	136 (83%)	12 (7%)	15 (9%)	1	1
All	All	629/704 (89%)	530 (84%)	57 (9%)	42 (7%)	1	3

5 of 42 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	47	ASP
1	B	46	ALA
1	B	115	THR
1	B	121	ALA
1	B	132	LEU

### 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	116/132 (88%)	97 (84%)	19 (16%)	2	7
1	B	123/132 (93%)	108 (88%)	15 (12%)	5	15
1	C	115/132 (87%)	93 (81%)	22 (19%)	1	4
1	D	122/132 (92%)	104 (85%)	18 (15%)	3	9
All	All	476/528 (90%)	402 (84%)	74 (16%)	2	8

5 of 74 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	165	ARG
1	C	43	ARG
1	D	127	GLU
1	C	4	LYS
1	C	26	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	151	ASN
1	D	55	GLN
1	C	161	HIS
1	B	39	GLN
1	D	39	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 [i](#)

Of 11 ligands modelled in this entry, 7 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	ADP	A	180	2	24,29,29	1.16	2 (8%)	29,45,45	2.54	13 (44%)
4	ADP	C	182	2	24,29,29	0.95	1 (4%)	29,45,45	2.61	13 (44%)
4	ADP	B	181	2	24,29,29	1.04	2 (8%)	29,45,45	2.67	9 (31%)
4	ADP	D	183	2	24,29,29	0.74	0	29,45,45	2.17	8 (27%)

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	ADP	A	180	2	-	5/12/32/32	0/3/3/3
4	ADP	C	182	2	-	3/12/32/32	0/3/3/3
4	ADP	B	181	2	-	3/12/32/32	0/3/3/3
4	ADP	D	183	2	-	5/12/32/32	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	180	ADP	C2'-C1'	-3.04	1.49	1.53
4	B	181	ADP	C8-N7	2.46	1.39	1.34
4	B	181	ADP	C2-N3	2.12	1.35	1.32
4	A	180	ADP	C5-N7	-2.06	1.32	1.39
4	C	182	ADP	PB-O3B	-2.03	1.47	1.54

The worst 5 of 43 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	181	ADP	C4-C5-N7	-8.25	100.81	109.40
4	D	183	ADP	N3-C2-N1	-7.09	117.60	128.68
4	B	181	ADP	C5-C6-N6	6.07	129.57	120.35
4	C	182	ADP	N3-C2-N1	-5.87	119.51	128.68
4	C	182	ADP	C2'-C3'-C4'	5.58	113.48	102.64

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

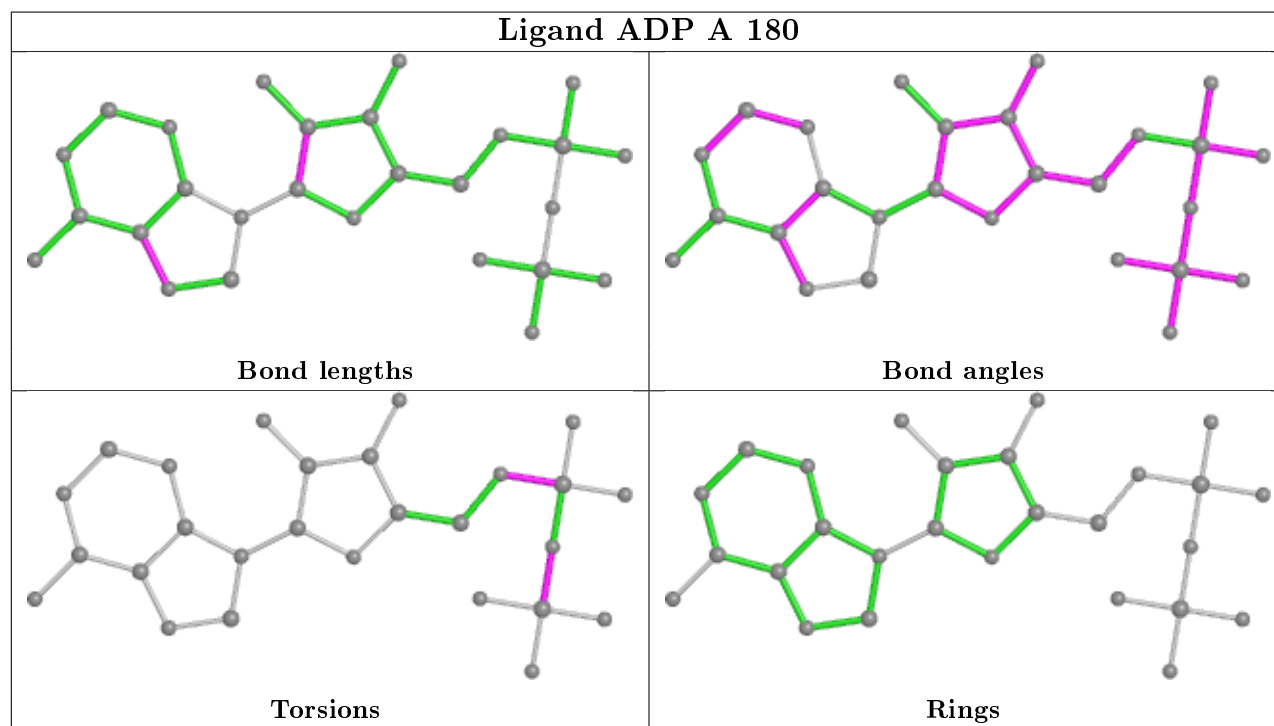
Mol	Chain	Res	Type	Atoms
4	A	180	ADP	C5'-O5'-PA-O1A
4	B	181	ADP	C5'-O5'-PA-O1A
4	C	182	ADP	C5'-O5'-PA-O1A
4	D	183	ADP	PA-O3A-PB-O2B
4	B	181	ADP	PA-O3A-PB-O1B

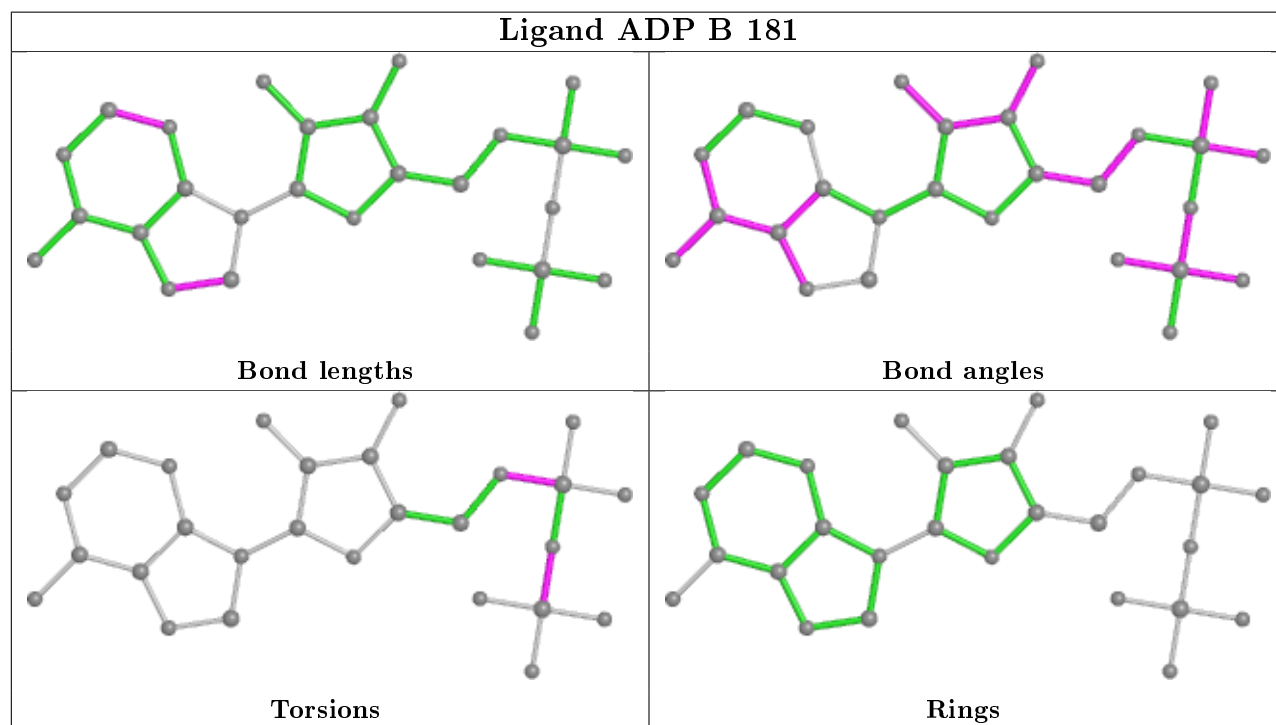
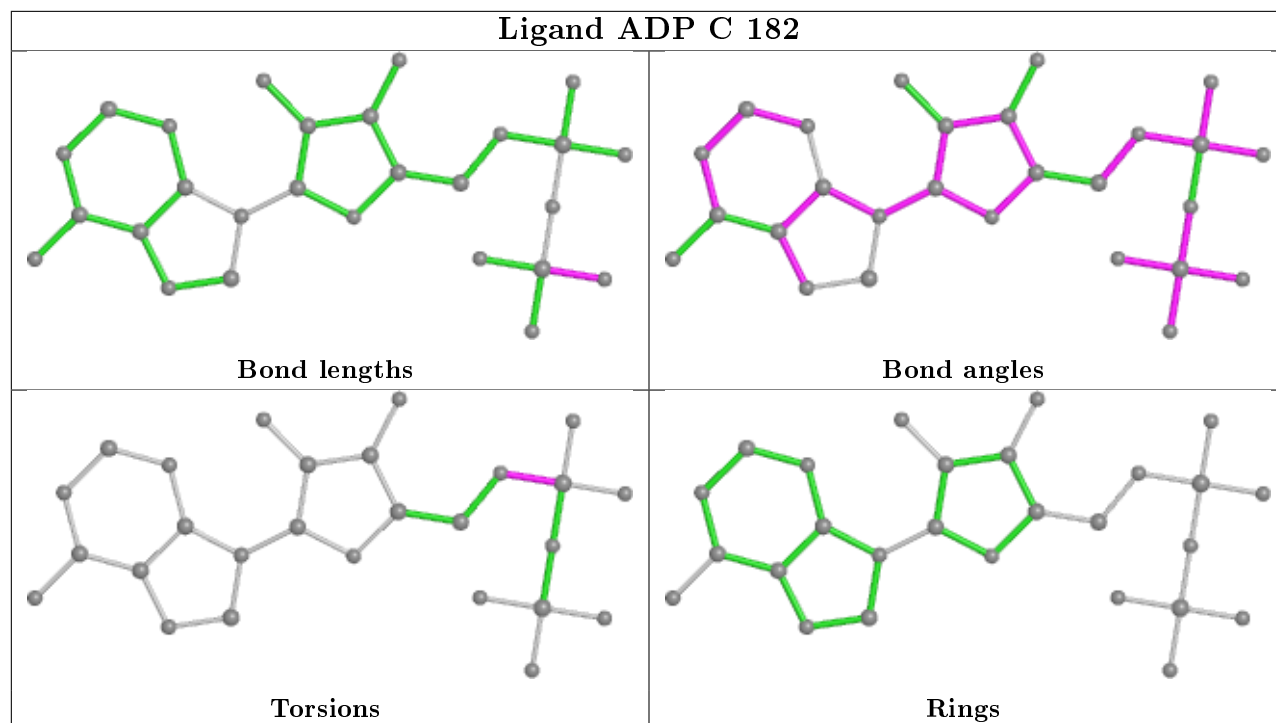
There are no ring outliers.

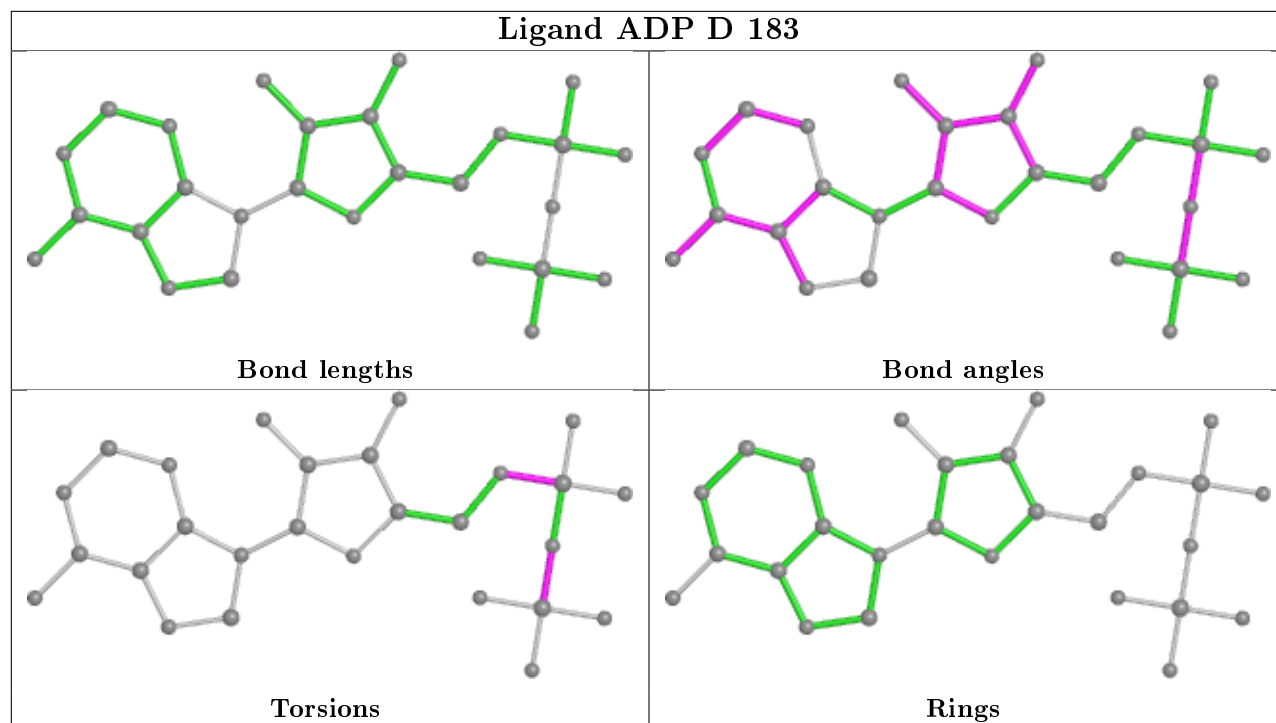
4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	180	ADP	6	0
4	C	182	ADP	8	0
4	B	181	ADP	2	0
4	D	183	ADP	6	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	156/176 (88%)	-0.45	0	100 100	12, 32, 53, 68	0
1	B	165/176 (93%)	-0.50	0	100 100	9, 28, 46, 62	0
1	C	155/176 (88%)	-0.35	0	100 100	18, 34, 61, 83	0
1	D	165/176 (93%)	-0.34	1 (0%)	89 86	20, 36, 54, 68	0
All	All	641/704 (91%)	-0.41	1 (0%)	95 94	9, 33, 53, 83	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	114	ASN	2.9

### 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, 95<sup>th</sup> 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.

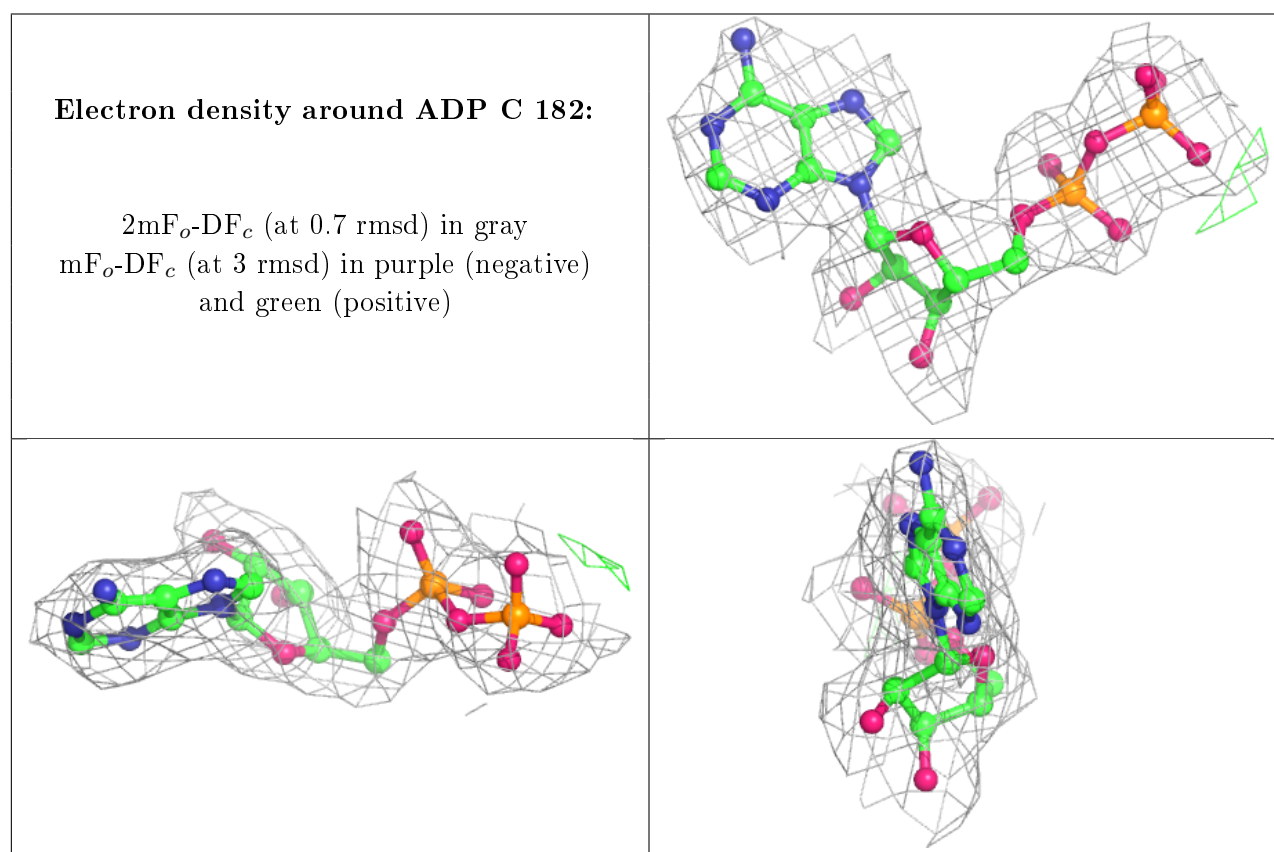
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	MG	C	202	1/1	0.79	0.21	29,29,29,29	0

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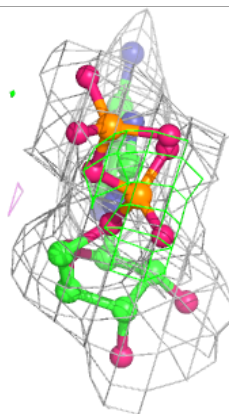
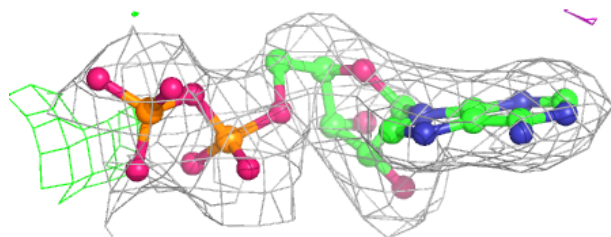
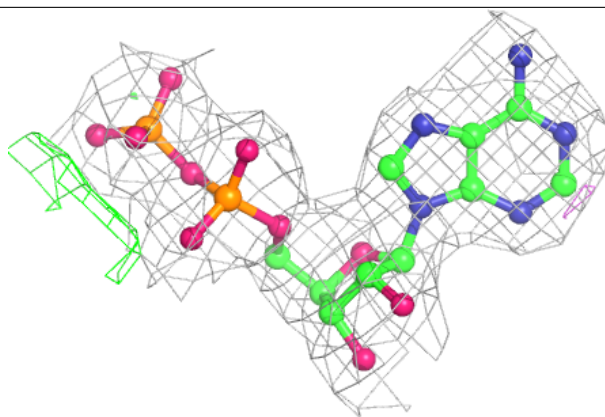
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MG	D	208	1/1	0.89	0.21	41,41,41,41	0
2	MG	A	190	1/1	0.90	0.23	22,22,22,22	0
3	CL	A	193	1/1	0.91	0.09	46,46,46,46	0
2	MG	B	196	1/1	0.92	0.20	15,15,15,15	0
3	CL	D	211	1/1	0.95	0.12	34,34,34,34	0
4	ADP	C	182	27/27	0.97	0.13	7,29,43,54	0
4	ADP	A	180	27/27	0.98	0.11	10,16,25,32	0
4	ADP	D	183	27/27	0.98	0.13	17,33,49,52	0
4	ADP	B	181	27/27	0.99	0.13	10,23,32,36	0
3	CL	B	199	1/1	0.99	0.10	26,26,26,26	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.

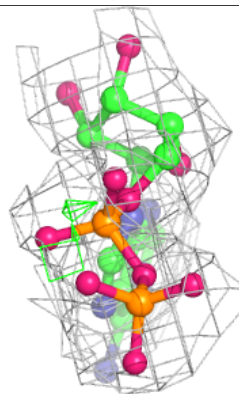
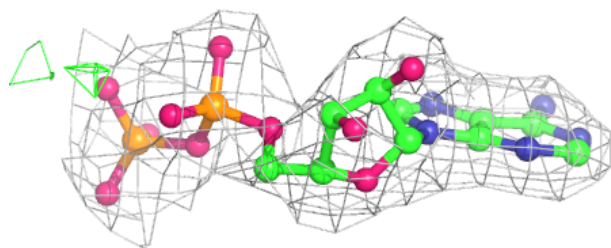
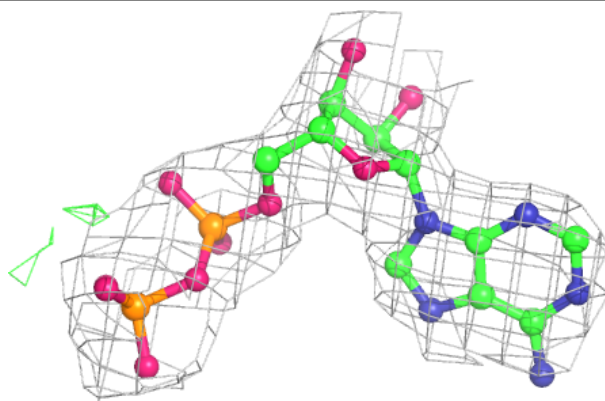


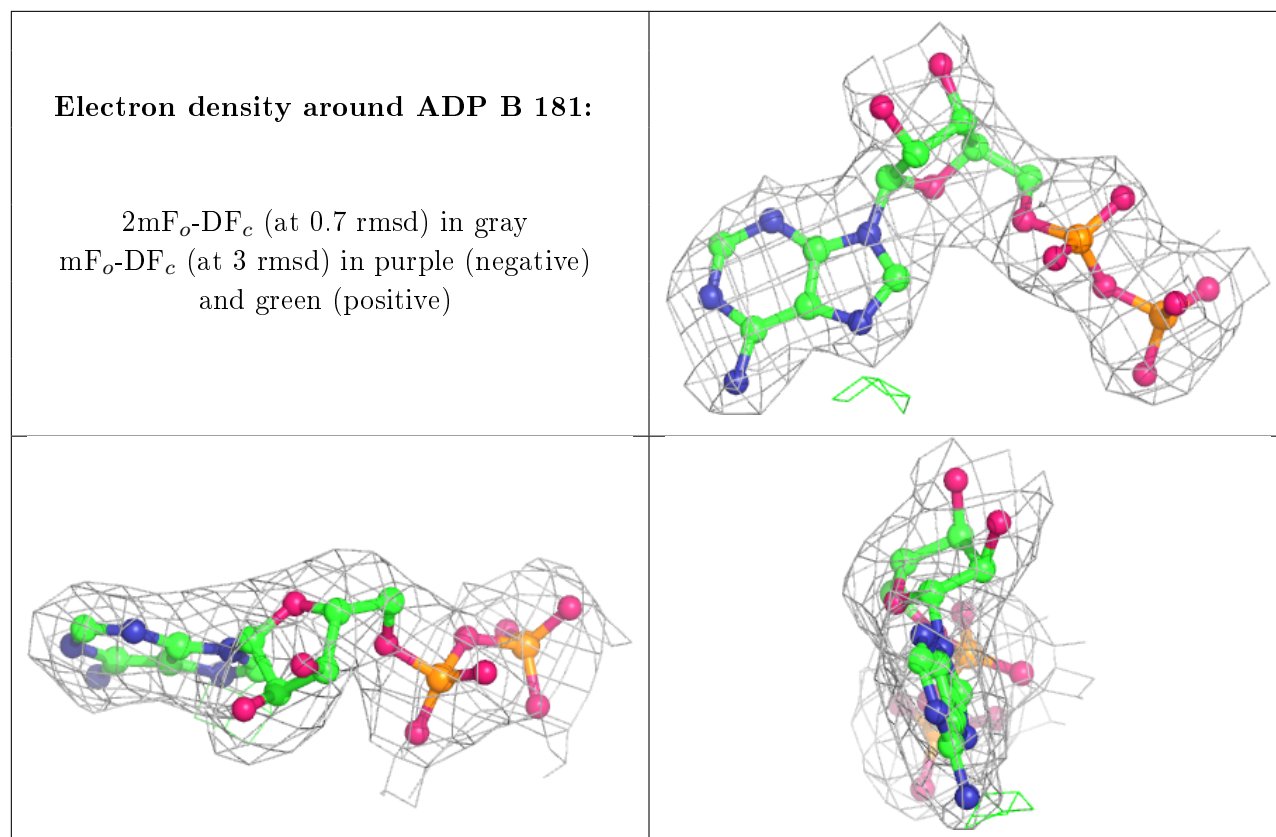
**Electron density around ADP A 180:**

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

**Electron density around ADP D 183:**

$2mF_o-DF_c$  (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.