



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 17, 2021 – 07:01 PM JST

PDB ID : 6M4I  
Title : Structure of CENP-E motor domain at 1.9 angstrom resolution  
Authors : Shibuya, A.; Yokoyama, H.  
Deposited on : 2020-03-07  
Resolution : 1.90 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](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 : 2.18  
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) : 2.18

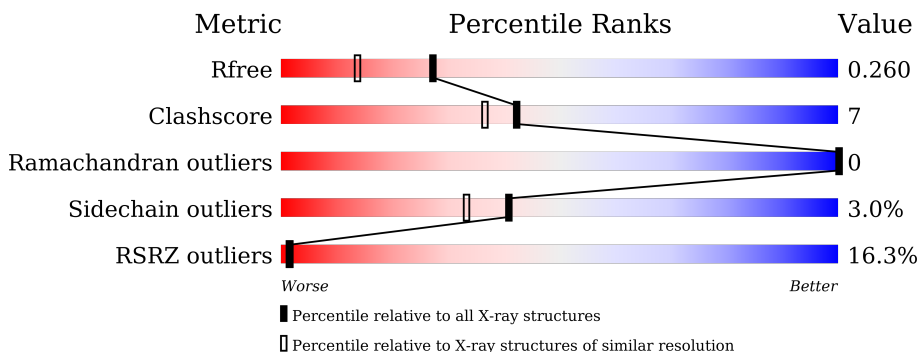
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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  $\geq 3$ , 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	353	<div> <div>10%</div> <div> <div></div> <div>75%</div> <div>11%</div> <div>•</div> <div>13%</div> </div> </div>
1	B	353	<div> <div>18%</div> <div> <div></div> <div>76%</div> <div>10%</div> <div>•</div> <div>14%</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5120 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 Centromere-associated protein E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	0	5	0
			2473	1560	433	469	11			
1	B	303	Total	C	N	O	S	0	4	0
			2451	1548	427	465	11			

There are 30 discrepancies between the modelled and reference sequences:

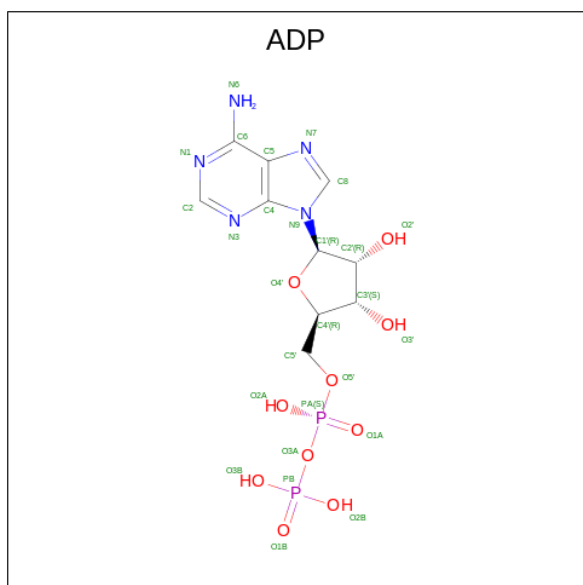
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	initiating methionine	UNP Q02224
A	-4	ASN	-	expression tag	UNP Q02224
A	-3	HIS	-	expression tag	UNP Q02224
A	-2	LYS	-	expression tag	UNP Q02224
A	-1	VAL	-	expression tag	UNP Q02224
A	0	HIS	-	expression tag	UNP Q02224
A	300	PRO	ALA	conflict	UNP Q02224
A	340	GLY	-	expression tag	UNP Q02224
A	341	SER	-	expression tag	UNP Q02224
A	342	HIS	-	expression tag	UNP Q02224
A	343	HIS	-	expression tag	UNP Q02224
A	344	HIS	-	expression tag	UNP Q02224
A	345	HIS	-	expression tag	UNP Q02224
A	346	HIS	-	expression tag	UNP Q02224
A	347	HIS	-	expression tag	UNP Q02224
B	-5	MET	-	initiating methionine	UNP Q02224
B	-4	ASN	-	expression tag	UNP Q02224
B	-3	HIS	-	expression tag	UNP Q02224
B	-2	LYS	-	expression tag	UNP Q02224
B	-1	VAL	-	expression tag	UNP Q02224
B	0	HIS	-	expression tag	UNP Q02224
B	300	PRO	ALA	conflict	UNP Q02224
B	340	GLY	-	expression tag	UNP Q02224
B	341	SER	-	expression tag	UNP Q02224
B	342	HIS	-	expression tag	UNP Q02224

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Chain	Residue	Modelled	Actual	Comment	Reference
B	343	HIS	-	expression tag	UNP Q02224
B	344	HIS	-	expression tag	UNP Q02224
B	345	HIS	-	expression tag	UNP Q02224
B	346	HIS	-	expression tag	UNP Q02224
B	347	HIS	-	expression tag	UNP Q02224

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

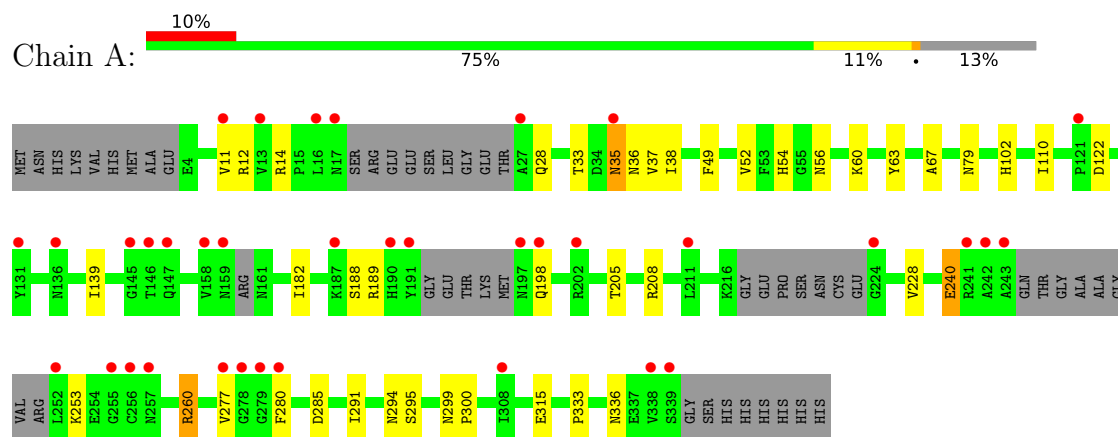
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	100	Total 100	O 100	0	0
4	B	40	Total 40	O 40	0	0

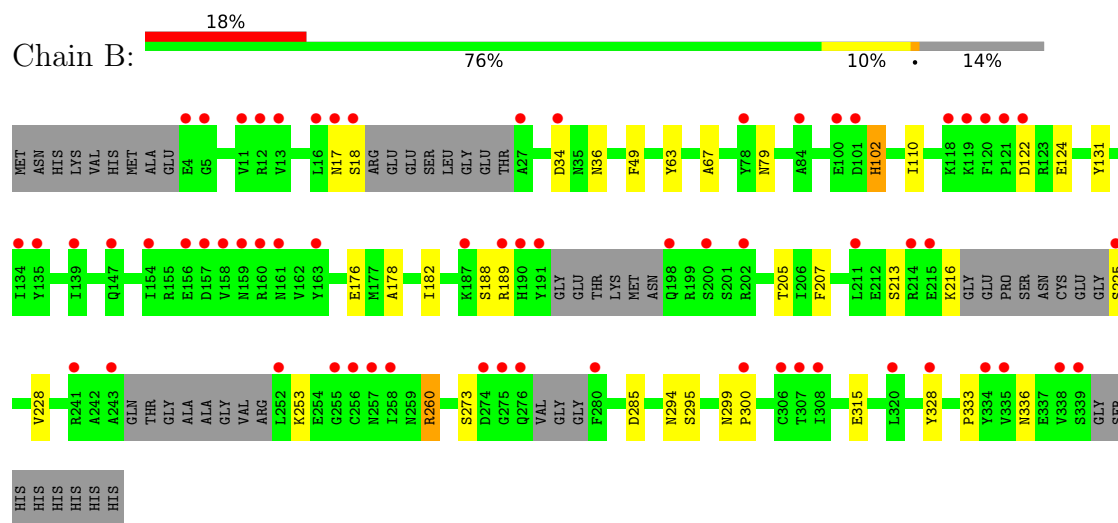
### 3 Residue-property plots [i](#)

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: Centromere-associated protein E



#### • Molecule 1: Centromere-associated protein E



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.83Å 82.84Å 49.37Å 90.00° 100.85° 90.00°	Depositor
Resolution (Å)	19.92 – 1.90 19.92 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (19.92-1.90) 99.9 (19.92-1.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.22 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.217 , 0.255 0.223 , 0.260	Depositor DCC
$R_{free}$ test set	6020 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.7	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 52.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5120	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.91% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ADP, 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	0.77	0/2513	0.90	4/3387 (0.1%)
1	B	0.67	0/2489	0.83	2/3354 (0.1%)
All	All	0.72	0/5002	0.86	6/6741 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	260	ARG	NE-CZ-NH1	7.70	124.15	120.30
1	A	260	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	B	260	ARG	NE-CZ-NH1	5.15	122.88	120.30
1	A	12	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	B	260	ARG	NE-CZ-NH2	-5.03	117.78	120.30
1	A	240	GLU	CB-CA-C	-5.02	100.36	110.40

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	2473	0	2463	40	0
1	B	2451	0	2455	28	0
2	A	27	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	27	0	12	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	100	0	0	7	0
4	B	40	0	0	2	0
All	All	5120	0	4942	68	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 7.

All (68) 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:38:ILE:HB	4:A:733:HOH:O	1.40	1.18
1:A:54[B]:HIS:ND1	1:A:56:ASN:OD1	2.05	0.89
1:B:299:ASN:O	1:B:333:PRO:HA	1.74	0.87
1:A:299:ASN:O	1:A:333:PRO:HA	1.75	0.86
1:B:182[B]:ILE:HD11	1:B:207:PHE:CE1	2.12	0.84
1:B:182[B]:ILE:HD11	1:B:207:PHE:HE1	1.43	0.81
1:A:54[B]:HIS:CE1	1:A:56:ASN:OD1	2.40	0.73
1:A:38:ILE:CB	4:A:733:HOH:O	2.12	0.70
1:A:291:ILE:O	4:A:701:HOH:O	2.08	0.70
1:B:294:ASN:O	1:B:300:PRO:CD	2.41	0.69
1:A:294:ASN:O	1:A:300:PRO:CD	2.40	0.68
1:A:277:VAL:O	1:A:280:PHE:HD2	1.78	0.67
1:A:11:VAL:HG11	1:A:38:ILE:CD1	2.25	0.67
1:A:60:LYS:H	1:A:102[B]:HIS:HE1	1.46	0.64
1:B:315:GLU:HA	1:B:315:GLU:OE1	2.04	0.58
1:A:33:THR:HG21	1:A:52:VAL:CG1	2.34	0.56
1:B:182[B]:ILE:N	1:B:182[B]:ILE:CD1	2.68	0.56
1:A:110:ILE:HD11	1:A:182:ILE:HD11	1.88	0.56
1:A:299:ASN:N	1:A:300:PRO:CD	2.68	0.56
1:A:295:SER:HA	1:A:300:PRO:CG	2.36	0.55
1:A:189:ARG:HD3	1:A:205:THR:HG21	1.89	0.55
1:B:299:ASN:N	1:B:300:PRO:CD	2.69	0.55
1:A:28:GLN:NE2	4:A:704:HOH:O	2.40	0.54
1:A:295:SER:HA	1:A:300:PRO:HG2	1.90	0.54
1:A:33:THR:HG21	1:A:52:VAL:HG11	1.90	0.54
1:B:299:ASN:O	1:B:299:ASN:OD1	2.26	0.54
1:A:294:ASN:O	1:A:300:PRO:HD3	2.07	0.54
1:A:315:GLU:HA	1:A:315:GLU:OE1	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:ASN:O	1:A:299:ASN:OD1	2.26	0.53
1:B:295:SER:HA	1:B:300:PRO:CG	2.39	0.53
1:B:294:ASN:O	1:B:300:PRO:HD3	2.08	0.52
1:B:178:ALA:O	1:B:182[B]:ILE:HD13	2.09	0.52
1:A:110:ILE:CD1	1:A:182:ILE:CD1	2.89	0.51
1:B:295:SER:HA	1:B:300:PRO:HG2	1.93	0.51
1:A:260:ARG:HD3	1:A:285:ASP:O	2.12	0.50
1:B:260:ARG:HD3	1:B:285:ASP:O	2.12	0.50
1:A:38:ILE:CG1	4:A:733:HOH:O	2.53	0.49
1:B:110:ILE:CD1	1:B:182[A]:ILE:CD1	2.90	0.49
1:B:228:VAL:O	1:B:336:ASN:HB2	2.12	0.49
1:A:228:VAL:O	1:A:336:ASN:HB2	2.12	0.49
1:B:102:HIS:HB2	4:B:737:HOH:O	2.12	0.49
1:B:189:ARG:HD3	1:B:205:THR:HG21	1.94	0.48
1:B:110:ILE:HD11	1:B:182[A]:ILE:HD11	1.95	0.48
1:B:299:ASN:N	1:B:300:PRO:HD3	2.30	0.47
1:B:182[B]:ILE:N	1:B:182[B]:ILE:HD13	2.30	0.46
1:A:33:THR:HG23	1:A:37:VAL:O	2.15	0.46
1:A:139:ILE:CD1	1:A:291:ILE:HD11	2.45	0.46
1:A:208:ARG:NH2	4:A:705:HOH:O	2.42	0.45
1:A:110:ILE:CD1	1:A:182:ILE:HD11	2.46	0.45
1:A:299:ASN:N	1:A:300:PRO:HD3	2.29	0.45
1:B:110:ILE:CD1	1:B:182[A]:ILE:HD11	2.46	0.45
1:A:299:ASN:O	1:A:333:PRO:CA	2.58	0.45
1:B:17:ASN:HB2	4:B:723:HOH:O	2.16	0.45
1:A:63:TYR:CD1	1:A:67:ALA:HB3	2.52	0.45
1:A:33:THR:HG22	1:A:35[A]:ASN:O	2.17	0.44
1:B:63:TYR:CD1	1:B:67:ALA:HB3	2.52	0.44
1:B:273:SER:HB2	1:B:328:TYR:HB2	2.00	0.44
1:A:11:VAL:HG11	1:A:38:ILE:HD11	1.98	0.44
1:B:36[B]:ASN:HB3	1:B:49:PHE:O	2.18	0.43
1:A:198:GLN:O	1:A:198:GLN:HG3	2.17	0.43
1:A:208:ARG:HD2	4:A:713:HOH:O	2.19	0.43
1:A:33:THR:CG2	1:A:52:VAL:HG11	2.50	0.42
1:A:36[A]:ASN:HB3	1:A:49:PHE:O	2.20	0.42
1:A:79:ASN:HB2	1:A:300:PRO:HB3	2.01	0.41
1:B:79:ASN:HB2	1:B:300:PRO:HB3	2.03	0.41
1:B:124:GLU:O	1:B:213:SER:HA	2.19	0.41
1:A:294:ASN:O	1:A:300:PRO:HD2	2.21	0.41
1:B:299:ASN:O	1:B:333:PRO:CA	2.58	0.41

There are no symmetry-related clashes.

## 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	299/353 (85%)	290 (97%)	9 (3%)	0	100	100
1	B	295/353 (84%)	285 (97%)	10 (3%)	0	100	100
All	All	594/706 (84%)	575 (97%)	19 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	274/307 (89%)	267 (97%)	7 (3%)	46	39
1	B	273/307 (89%)	262 (96%)	11 (4%)	31	22
All	All	547/614 (89%)	529 (97%)	18 (3%)	41	29

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	35[A]	ASN
1	A	35[B]	ASN
1	A	122	ASP
1	A	188	SER
1	A	240	GLU
1	A	253	LYS
1	B	18	SER

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Mol	Chain	Res	Type
1	B	34[A]	ASP
1	B	34[B]	ASP
1	B	102	HIS
1	B	122	ASP
1	B	131	TYR
1	B	176	GLU
1	B	188	SER
1	B	216	LYS
1	B	225	SER
1	B	253	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z  > 2$
2	ADP	B	600	3	24,29,29	0.75	1 (4%)	29,45,45	0.77	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	ADP	A	600	3	24,29,29	0.83	1 (4%)	29,45,45	0.85	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
2	ADP	B	600	3	-	2/12/32/32	0/3/3/3
2	ADP	A	600	3	-	2/12/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	600	ADP	C8-N7	-2.06	1.31	1.34
2	B	600	ADP	C8-N7	-2.03	1.31	1.34

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

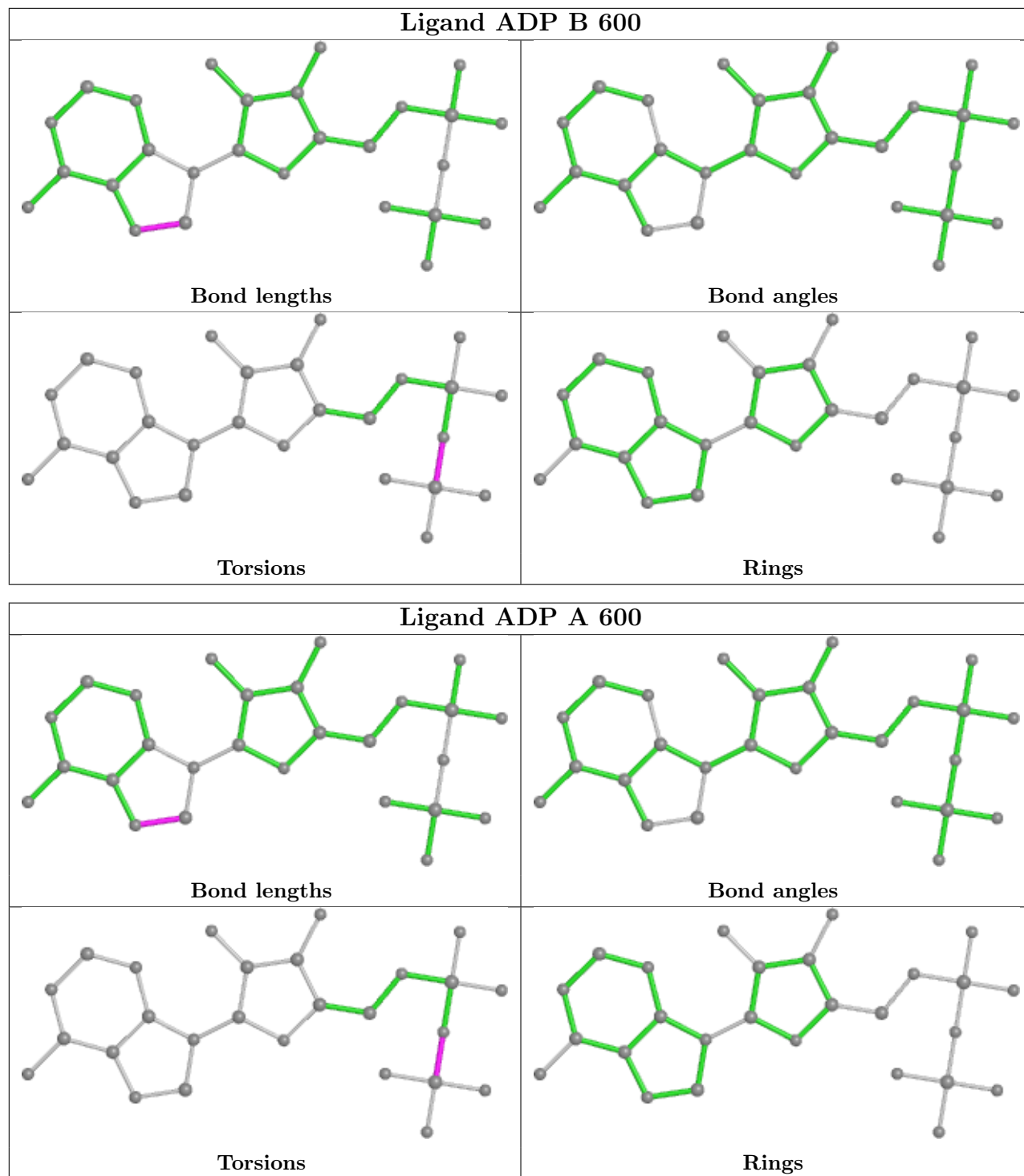
Mol	Chain	Res	Type	Atoms
2	A	600	ADP	PA-O3A-PB-O2B
2	B	600	ADP	PA-O3A-PB-O2B
2	A	600	ADP	PA-O3A-PB-O1B
2	B	600	ADP	PA-O3A-PB-O3B

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

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, 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	306/353 (86%)	0.76	36 (11%) 4 5	28, 45, 92, 130	0
1	B	303/353 (85%)	1.25	63 (20%) 1 1	43, 64, 114, 138	0
All	All	609/706 (86%)	1.01	99 (16%) 1 1	28, 56, 105, 138	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	256	CYS	11.1
1	A	158	VAL	10.2
1	B	243	ALA	10.2
1	B	191	TYR	9.8
1	B	158	VAL	9.1
1	A	191	TYR	9.0
1	A	256	CYS	8.9
1	B	338	VAL	8.0
1	B	17	ASN	8.0
1	A	27	ALA	7.7
1	B	147	GLN	7.7
1	A	255	GLY	7.5
1	B	339	SER	7.1
1	B	280	PHE	6.3
1	B	241	ARG	6.2
1	A	252	LEU	6.2
1	A	198	GLN	6.1
1	B	198	GLN	6.1
1	B	334	TYR	6.1
1	B	27	ALA	5.9
1	A	339	SER	5.9
1	B	121	PRO	5.7
1	B	18	SER	5.5
1	B	160	ARG	5.4

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Mol	Chain	Res	Type	RSRZ
1	B	190	HIS	5.3
1	B	159	ASN	5.2
1	A	280	PHE	5.1
1	B	252	LEU	5.0
1	A	190	HIS	5.0
1	A	278	GLY	4.8
1	B	16	LEU	4.5
1	B	101	ASP	4.5
1	A	277	VAL	4.3
1	B	275	GLY	4.3
1	B	308	ILE	4.3
1	A	16	LEU	4.3
1	A	17	ASN	4.0
1	A	243	ALA	3.9
1	A	224	GLY	3.9
1	A	279	GLY	3.8
1	B	335	VAL	3.4
1	B	163	TYR	3.4
1	B	276	GLN	3.3
1	B	255	GLY	3.3
1	B	225	SER	3.3
1	A	308	ILE	3.3
1	B	257	ASN	3.3
1	B	156	GLU	3.3
1	B	307	THR	3.2
1	A	197	ASN	3.2
1	B	5	GLY	3.2
1	B	187	LYS	3.2
1	B	211	LEU	3.2
1	A	145	GLY	3.2
1	A	211	LEU	3.1
1	B	100	GLU	3.1
1	A	147	GLN	3.1
1	B	154	ILE	3.1
1	B	157	ASP	3.0
1	B	258	ILE	3.0
1	A	257	ASN	3.0
1	B	119	LYS	2.9
1	B	200	SER	2.9
1	B	215	GLU	2.9
1	B	328	TYR	2.9
1	A	146	THR	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	35[A]	ASN	2.8
1	A	11	VAL	2.8
1	A	13	VAL	2.8
1	B	134	ILE	2.7
1	B	11	VAL	2.7
1	A	136	ASN	2.7
1	B	13	VAL	2.6
1	A	159	ASN	2.6
1	B	202	ARG	2.6
1	B	122	ASP	2.6
1	B	120	PHE	2.5
1	B	118	LYS	2.5
1	A	121	PRO	2.5
1	A	202	ARG	2.5
1	B	189	ARG	2.5
1	B	4	GLU	2.4
1	A	242	ALA	2.4
1	B	84	ALA	2.4
1	B	274	ASP	2.3
1	B	161	ASN	2.3
1	B	306	CYS	2.2
1	B	139	ILE	2.2
1	B	12	ARG	2.2
1	B	135	TYR	2.2
1	A	131	TYR	2.1
1	A	338	VAL	2.1
1	B	34[A]	ASP	2.1
1	A	241	ARG	2.1
1	B	214	ARG	2.1
1	B	78	TYR	2.1
1	B	300	PRO	2.1
1	A	187	LYS	2.1
1	B	320	LEU	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](#)

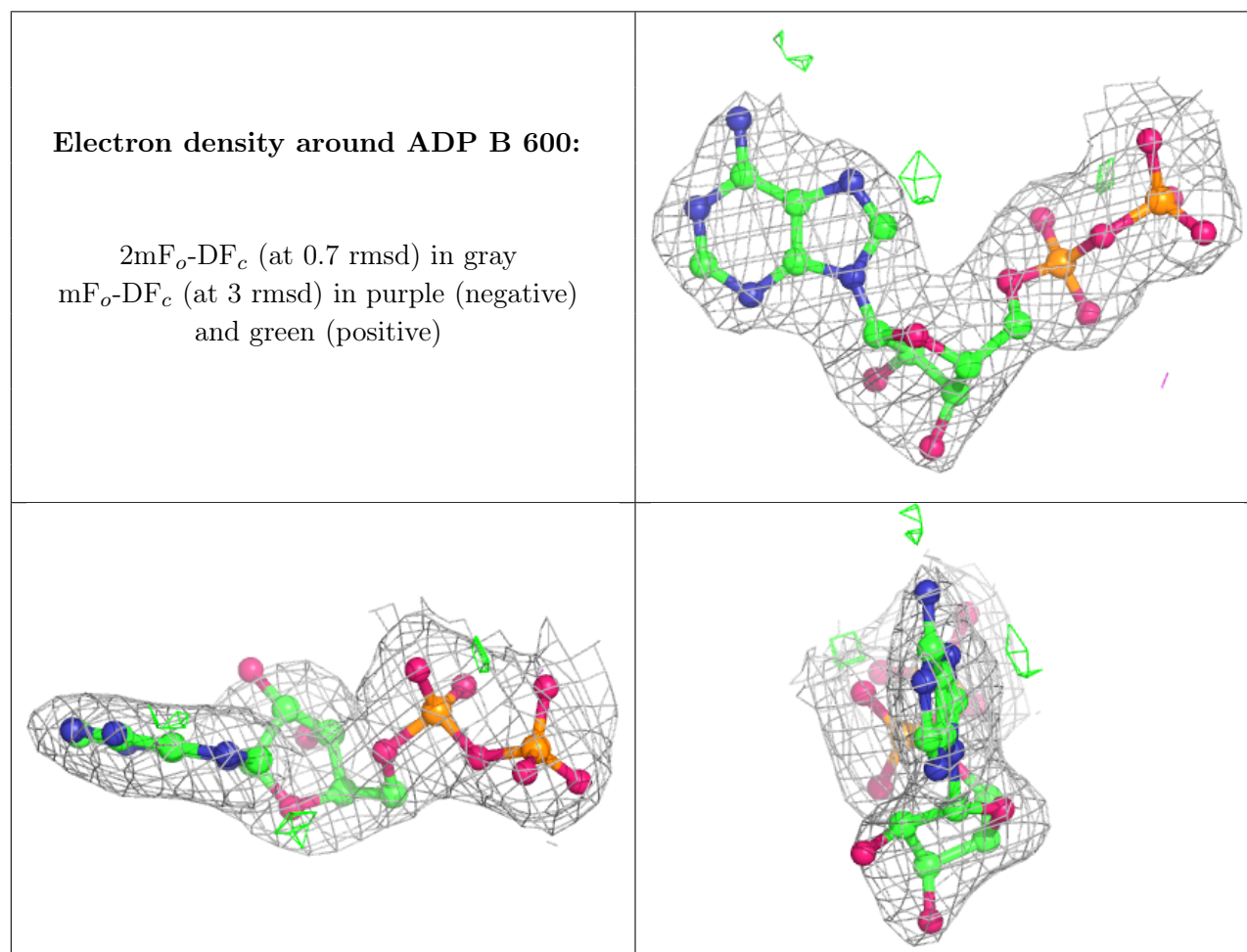
There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

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.

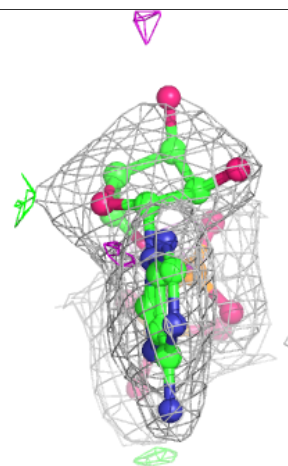
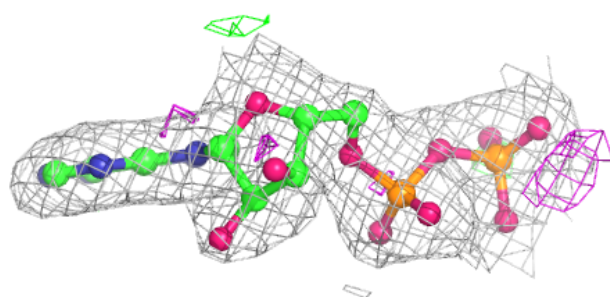
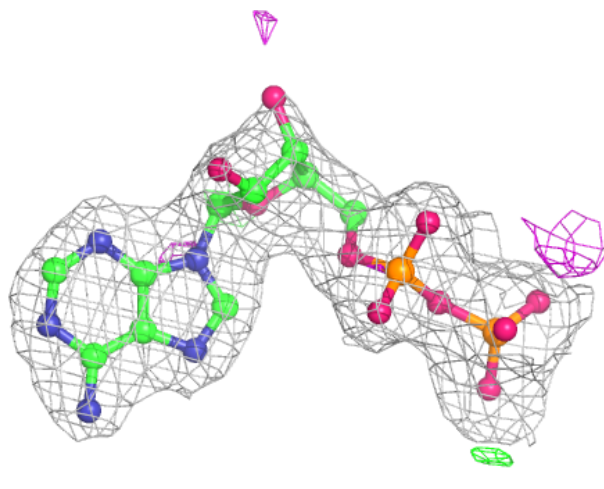
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	ADP	B	600	27/27	0.95	0.13	52,60,70,84	0
2	ADP	A	600	27/27	0.96	0.10	33,41,57,72	0
3	MG	A	601	1/1	0.96	0.11	45,45,45,45	0
3	MG	B	601	1/1	0.97	0.03	51,51,51,51	0

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.



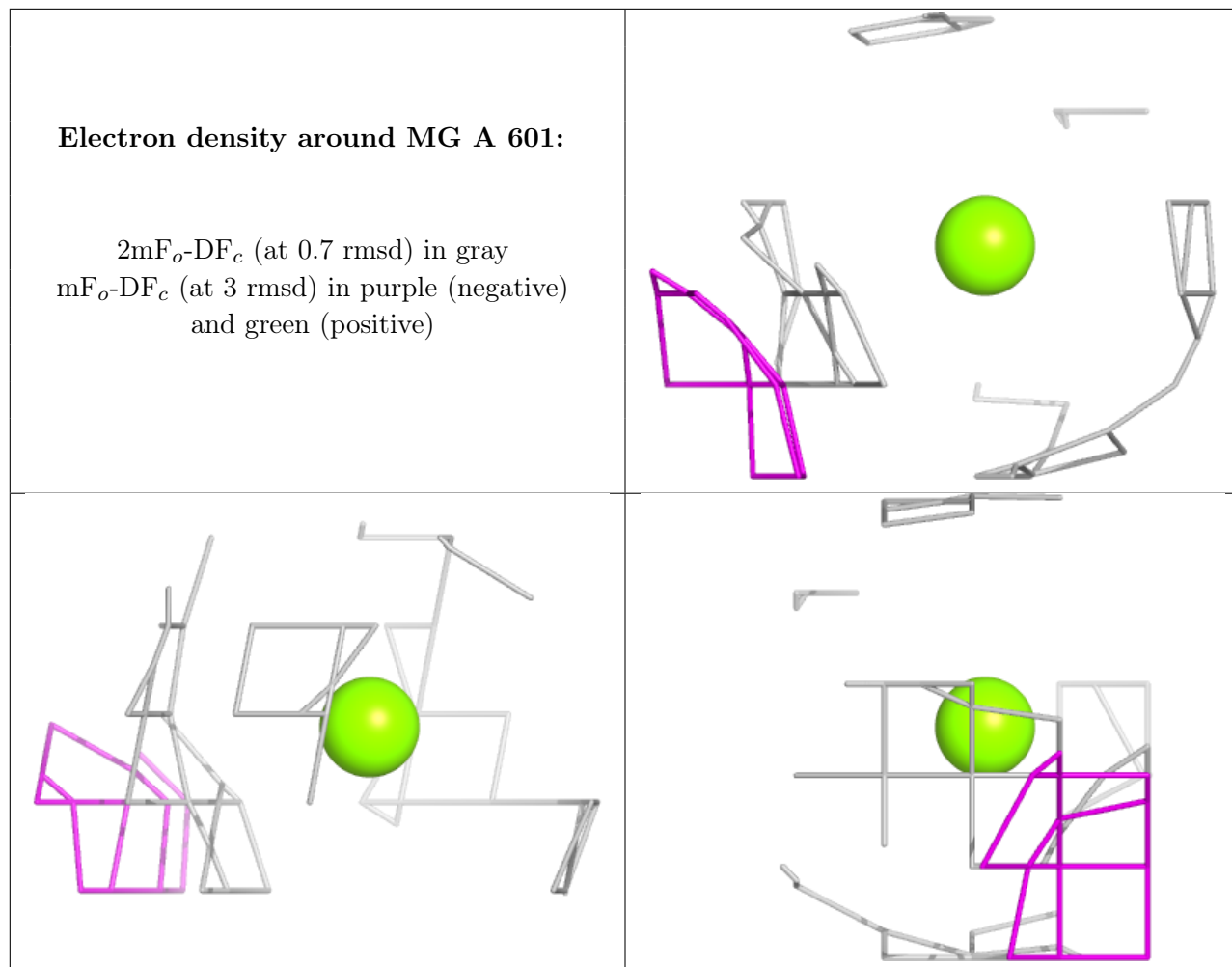
**Electron density around ADP A 600:**

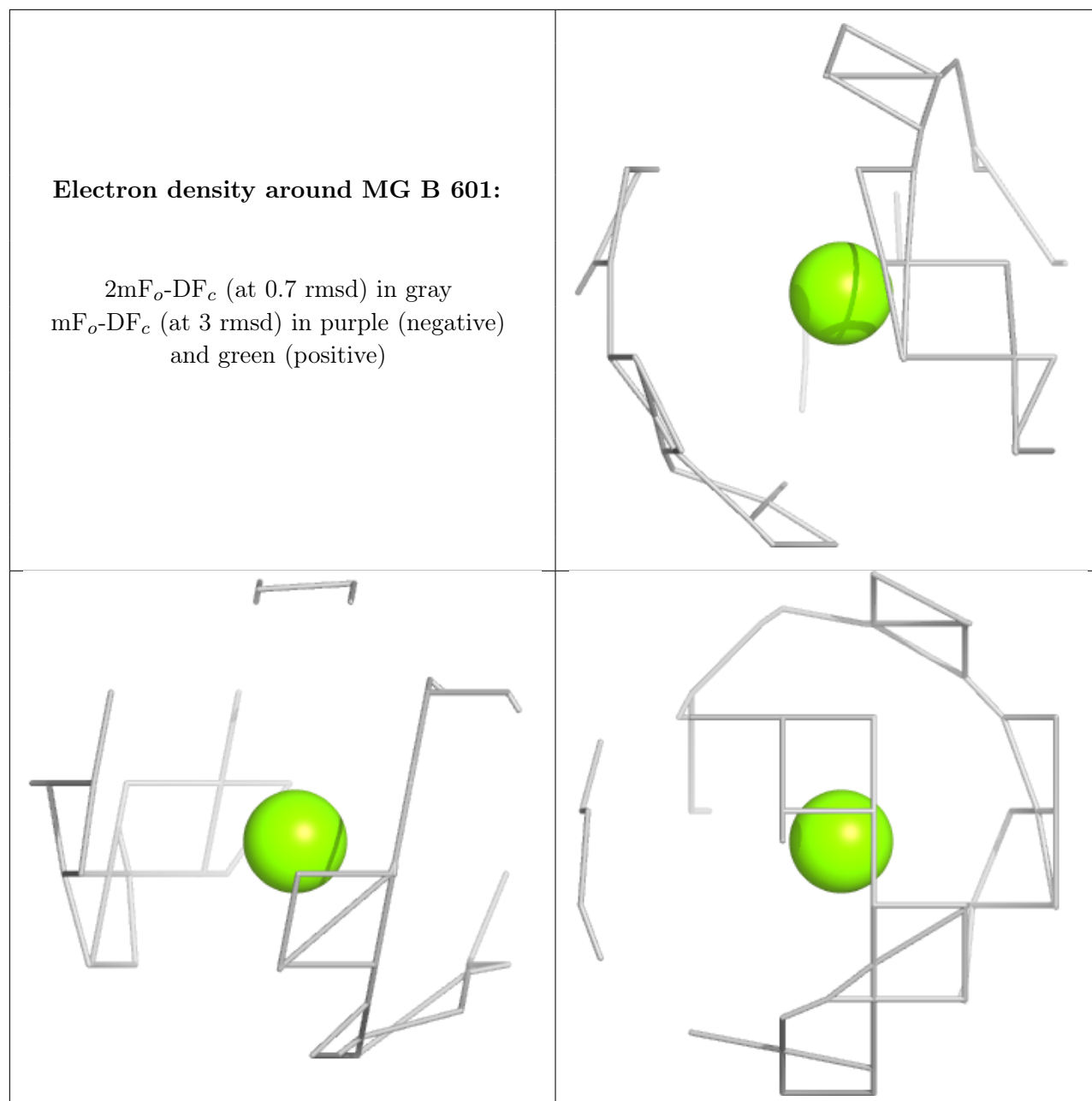
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around MG A 601:**

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





## 6.5 Other polymers ⓘ

There are no such residues in this entry.