



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 22, 2025 – 04:00 AM EDT

PDB ID : 6B4K / pdb\_00006b4k  
Title : Crystal structure of human DDX19B(AMPPNP)  
Authors : Lin, D.H.; Correia, A.R.; Cai, S.W.; Huber, F.M.; Jette, C.A.; Hoelz, A.  
Deposited on : 2017-09-26  
Resolution : 2.20 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0rc1  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.006 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.42

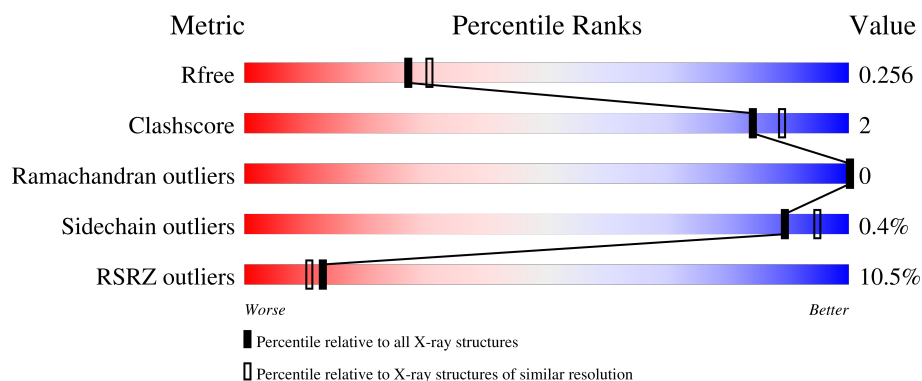
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
$R_{free}$	164625	5791 (2.20-2.20)
Clashscore	180529	6634 (2.20-2.20)
Ramachandran outliers	177936	6560 (2.20-2.20)
Sidechain outliers	177891	6561 (2.20-2.20)
RSRZ outliers	164620	5791 (2.20-2.20)

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	430	<div> <div>8%</div> <div>92%</div> <div>5%</div> </div>
1	B	430	<div> <div>12%</div> <div>92%</div> <div>6%</div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13924 atoms, of which 6898 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 ATP-dependent RNA helicase DDX19B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	419	Total	C	H	N	O	S	0	11	0
			6844	2141	3457	591	635	20			
1	B	420	Total	C	H	N	O	S	0	1	0
			6737	2106	3402	581	628	20			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	50	GLY	-	expression tag	UNP Q9UMR2
A	51	PRO	-	expression tag	UNP Q9UMR2
A	52	HIS	-	expression tag	UNP Q9UMR2
A	53	MET	-	expression tag	UNP Q9UMR2
B	50	GLY	-	expression tag	UNP Q9UMR2
B	51	PRO	-	expression tag	UNP Q9UMR2
B	52	HIS	-	expression tag	UNP Q9UMR2
B	53	MET	-	expression tag	UNP Q9UMR2

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

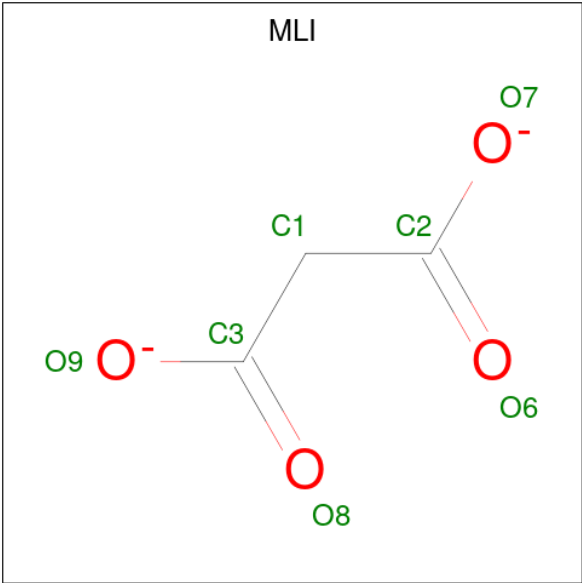
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>6</sub>O<sub>12</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	P	0	0
			48	10	17	6	12	3		
3	B	1	Total	C	H	N	O	P	0	0
			47	10	16	6	12	3		

- Molecule 4 is MALONATE ION (CCD ID: MLI) (formula: C<sub>3</sub>H<sub>2</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			9	3	2	4		
4	A	1	Total	C	H	O	0	0
			9	3	2	4		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	H	O	0	0
			9	3	2	4		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	137	Total	O	0	2
			139	139		
5	B	78	Total	O	0	2
			80	80		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.23Å 45.36Å 127.39Å 90.00° 97.03° 90.00°	Depositor
Resolution (Å)	42.14 – 2.20 42.14 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.0 (42.14-2.20) 99.2 (42.14-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.14	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.04 (at 2.20Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, $R_{free}$	0.207 , 0.248 0.212 , 0.256	Depositor DCC
$R_{free}$ test set	46300 reflections (4.14%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.3	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 57.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	13924	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.64 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.3168e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: MG, ANP, MLI

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	0.28	0/3475	0.46	0/4688
1	B	0.27	0/3387	0.46	0/4571
All	All	0.27	0/6862	0.46	0/9259

There are no bond length outliers.

There are no bond angle outliers.

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	3387	3457	3404	15	0
1	B	3335	3402	3400	17	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	31	17	13	1	0
3	B	31	16	13	0	0
4	A	14	4	4	1	0
4	B	7	2	2	0	0
5	A	139	0	0	1	0
5	B	80	0	0	0	0
All	All	7026	6898	6836	30	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 (30) 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:194:LEU:HD11	1:A:215:ILE:HG12	1.79	0.64
1:A:251:GLN:NE2	5:A:602:HOH:O	2.32	0.61
1:B:443:VAL:HG11	1:B:449:MET:HA	1.89	0.55
1:B:316:SER:N	1:B:444:ASP:OD2	2.42	0.53
1:B:443:VAL:HG11	1:B:449:MET:CA	2.41	0.50
1:A:366:GLU:CB	1:B:203:LEU:HD11	2.42	0.49
1:B:316:SER:CA	1:B:444:ASP:OD2	2.61	0.49
1:B:342:THR:HG22	1:B:344:LYS:H	1.77	0.48
1:B:466:LEU:HD13	1:B:477:ILE:HD11	1.96	0.48
1:A:251:GLN:NE2	4:A:504:MLI:O8	2.47	0.47
1:B:467:ASP:OD2	1:B:469:ASP:OD2	2.31	0.47
1:A:64:LYS:NZ	3:A:502:ANP:O3G	2.45	0.46
1:B:323:ALA:HA	1:B:468:THR:HG21	1.98	0.46
1:B:203:LEU:HD23	1:B:230:PHE:CD2	2.51	0.45
1:B:277:PHE:CZ	1:B:296:LYS:HE2	2.52	0.44
1:A:67:ARG:NH2	1:A:140:SER:OG	2.50	0.44
1:B:66:ILE:O	1:B:380:ARG:NH1	2.50	0.43
1:A:308:LYS:HE2	1:A:310:TYR:OH	2.18	0.43
1:A:310:TYR:CZ	1:A:477:ILE:HG22	2.53	0.42
1:A:82:ARG:NE	1:A:291:ASP:HB3	2.34	0.42
1:B:316:SER:C	1:B:444:ASP:OD2	2.58	0.42
1:B:304:LEU:HD22	1:B:424:LEU:CD1	2.50	0.42
1:B:203:LEU:HD23	1:B:230:PHE:CE2	2.55	0.42
1:A:368:MET:HE1	1:B:201:ASN:CG	2.41	0.41
1:A:474:ILE:O	1:A:477:ILE:HG13	2.21	0.41
1:B:380:ARG:NE	1:B:400:GLU:HG2	2.36	0.41
1:A:72:ASP:OD1	1:A:72:ASP:N	2.54	0.41
1:A:145:THR:HA	1:A:148:PHE:CE2	2.56	0.40
1:A:361:ALA:HB1	1:A:375:VAL:HG13	2.03	0.40
1:A:331:ILE:HD13	1:A:404:VAL:HG21	2.03	0.40

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	424/430 (99%)	421 (99%)	3 (1%)	0	100	100
1	B	417/430 (97%)	411 (99%)	6 (1%)	0	100	100
All	All	841/860 (98%)	832 (99%)	9 (1%)	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	380/378 (100%)	379 (100%)	1 (0%)	91	96
1	B	371/378 (98%)	369 (100%)	2 (0%)	86	93
All	All	751/756 (99%)	748 (100%)	3 (0%)	89	95

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	277	PHE
1	B	64	LYS
1	B	343	ARG

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 oligosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	MLI	A	503	-	6,6,6	1.27	0	7,7,7	1.04	0
4	MLI	B	503	-	6,6,6	1.21	0	7,7,7	1.09	0
3	ANP	A	502	2	29,33,33	0.77	0	31,52,52	0.72	1 (3%)
3	ANP	B	502	2	29,33,33	0.70	0	31,52,52	0.65	1 (3%)
4	MLI	A	504	-	6,6,6	1.23	0	7,7,7	1.11	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	MLI	A	503	-	-	0/4/4/4	-
4	MLI	B	503	-	-	0/4/4/4	-
3	ANP	A	502	2	-	4/14/38/38	0/3/3/3
3	ANP	B	502	2	-	5/14/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MLI	A	504	-	-	2/4/4/4	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	ANP	C5-C6-N6	2.43	124.01	120.31
3	A	502	ANP	C5-C6-N6	2.38	123.94	120.31

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	ANP	PB-N3B-PG-O1G
3	A	502	ANP	PG-N3B-PB-O1B
3	A	502	ANP	PA-O3A-PB-O2B
3	B	502	ANP	PB-N3B-PG-O1G
3	B	502	ANP	PG-N3B-PB-O1B
4	A	504	MLI	C3-C1-C2-O6
4	A	504	MLI	C3-C1-C2-O7
3	B	502	ANP	PA-O3A-PB-O2B
3	A	502	ANP	PA-O3A-PB-O1B
3	B	502	ANP	PA-O3A-PB-O1B
3	B	502	ANP	PG-N3B-PB-O3A

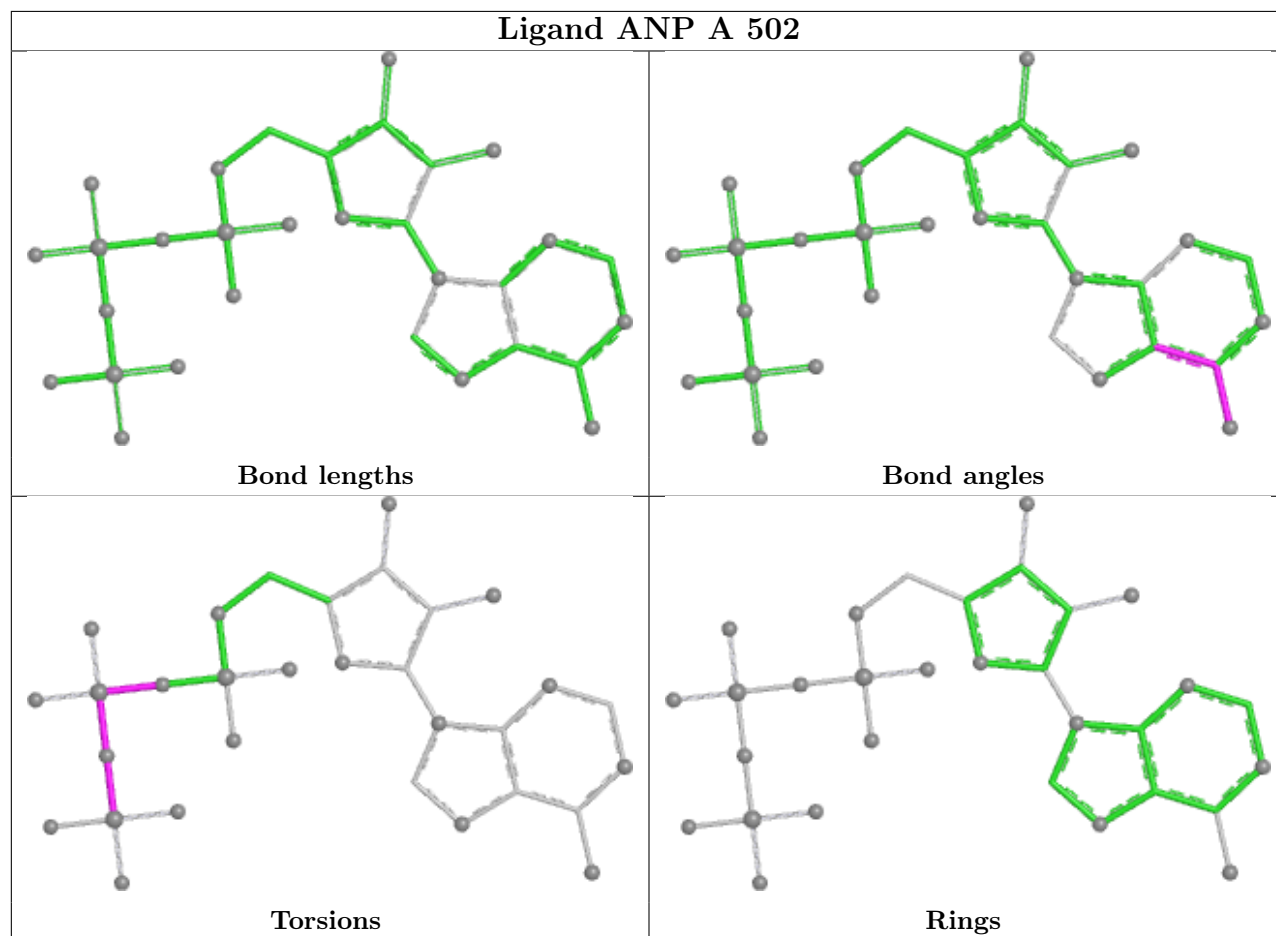
There are no ring outliers.

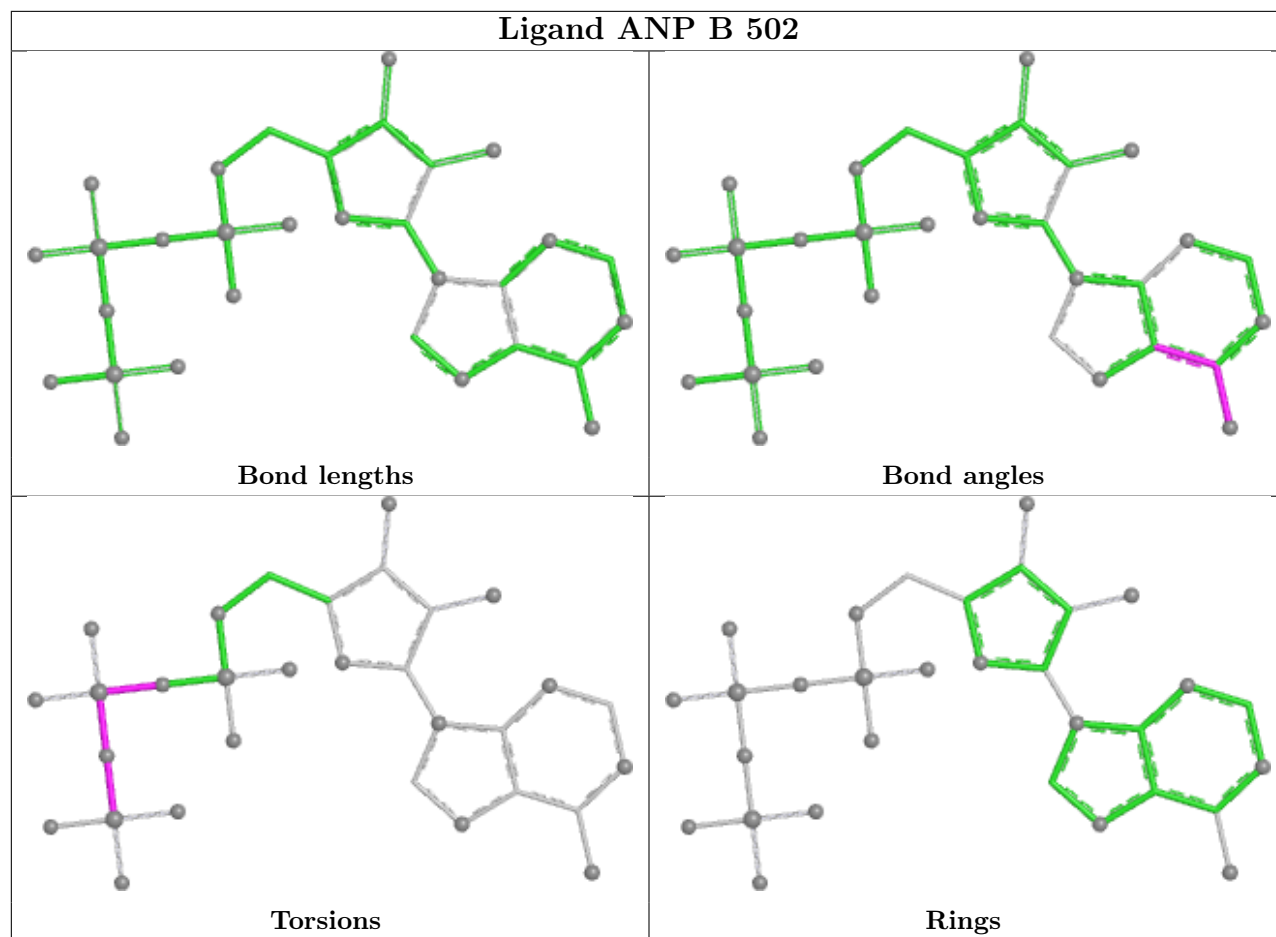
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	ANP	1	0
4	A	504	MLI	1	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

### 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	419/430 (97%)	0.50	36 (8%) 18 15	20, 68, 135, 171	7 (1%)
1	B	420/430 (97%)	0.78	52 (12%) 9 7	33, 86, 181, 270	1 (0%)
All	All	839/860 (97%)	0.64	88 (10%) 13 11	20, 75, 164, 270	8 (0%)

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	348[A]	TRP	5.1
1	B	428	GLY	4.7
1	A	89[A]	TYR	4.5
1	B	189	TYR	4.3
1	A	277	PHE	4.3
1	A	428	GLY	4.2
1	B	126	MET	4.2
1	B	277	PHE	4.0
1	B	324	LEU	3.9
1	A	334	ALA	3.8
1	B	295	ILE	3.8
1	B	165	CYS	3.7
1	B	330	ALA	3.6
1	B	474	ILE	3.3
1	A	434	GLY	3.2
1	B	427	ILE	3.1
1	A	328	TYR	3.0
1	A	203	LEU	3.0
1	A	152	MET	3.0
1	A	207	GLN	3.0
1	B	402	VAL	2.9
1	B	386	VAL	2.8
1	A	206	GLY	2.8
1	B	125	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	402	VAL	2.8
1	A	478	ALA	2.8
1	B	468	THR	2.8
1	B	460	ASN	2.8
1	B	262	ARG	2.7
1	B	396	GLY	2.7
1	B	451	ILE	2.7
1	B	88	LEU	2.7
1	A	234	LYS	2.7
1	A	331	ILE	2.6
1	B	84	PRO	2.6
1	A	190	PRO	2.6
1	A	209	ILE	2.6
1	B	208	LYS	2.6
1	A	335	GLN	2.5
1	A	330	ALA	2.5
1	B	321	PHE	2.5
1	B	304	LEU	2.5
1	B	316	SER	2.5
1	B	329	GLY	2.5
1	B	360	VAL	2.5
1	B	471	LEU	2.5
1	A	378[A]	ARG	2.4
1	B	332	THR	2.4
1	B	302	GLU	2.4
1	B	303	THR	2.4
1	B	346	ALA	2.4
1	B	466	LEU	2.4
1	B	172	TYR	2.3
1	A	327	LEU	2.3
1	B	408	PHE	2.3
1	B	72	ASP	2.3
1	B	357	GLY	2.3
1	A	125	LEU	2.3
1	B	313	LEU	2.3
1	B	250	THR	2.3
1	B	82	ARG	2.3
1	A	137	GLN	2.2
1	A	401	GLN	2.2
1	B	315	SER	2.2
1	B	412	VAL	2.2
1	A	304	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	358	HIS	2.2
1	B	462	LYS	2.2
1	B	331	ILE	2.2
1	A	468	THR	2.2
1	B	86	SER	2.1
1	A	357	GLY	2.1
1	B	209	ILE	2.1
1	A	341	HIS	2.1
1	A	202	LYS	2.1
1	B	311	TYR	2.1
1	A	88[A]	LEU	2.1
1	A	303	THR	2.1
1	B	398	ASP	2.1
1	B	478	ALA	2.1
1	B	399	VAL	2.1
1	A	444	ASP	2.1
1	B	333	ILE	2.1
1	B	350	ALA	2.1
1	A	82	ARG	2.0
1	A	346	ALA	2.0
1	B	206	GLY	2.0
1	A	251	GLN	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 [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.

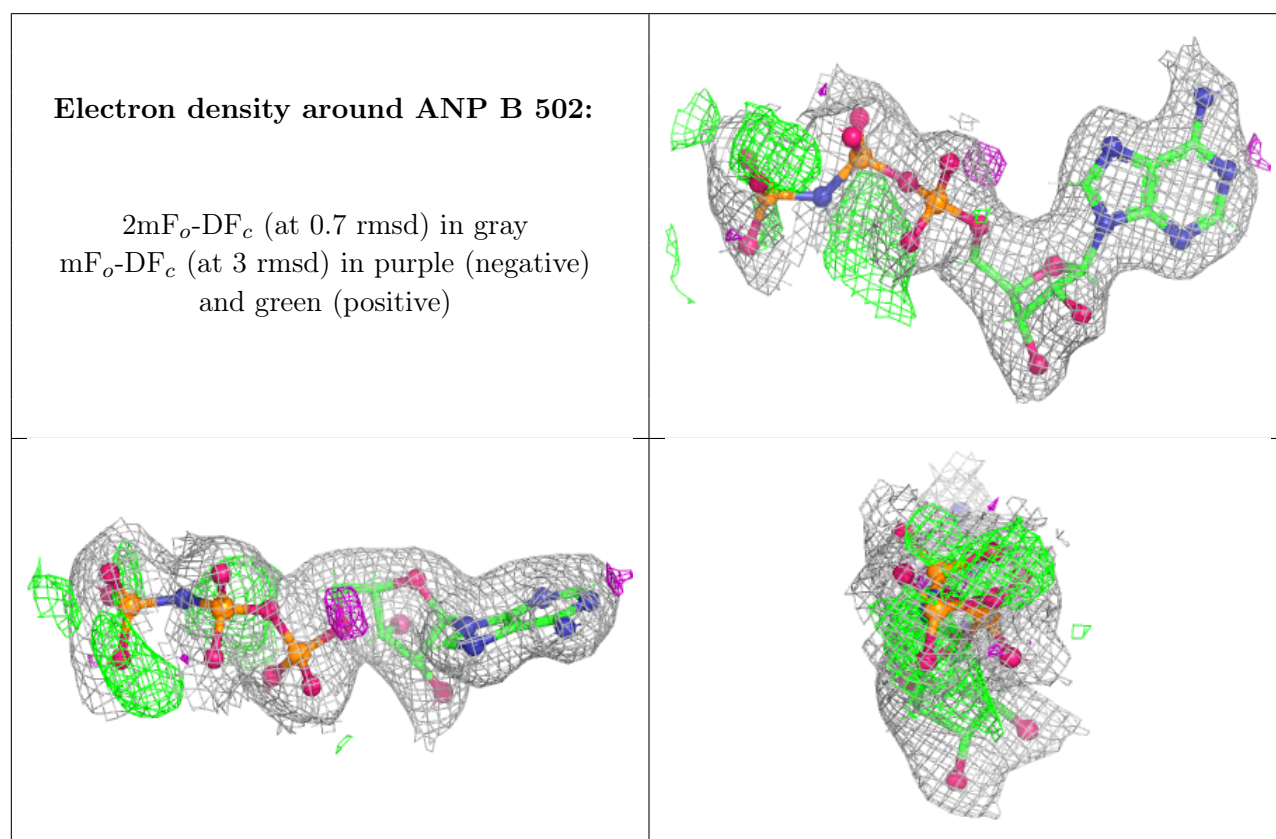
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	MG	A	501	1/1	0.88	0.15	54,54,54,54	0

*Continued on next page...*

*Continued from previous page...*

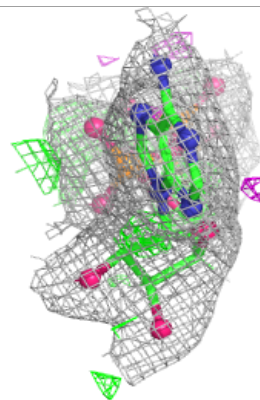
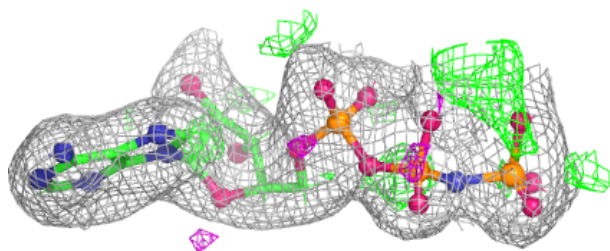
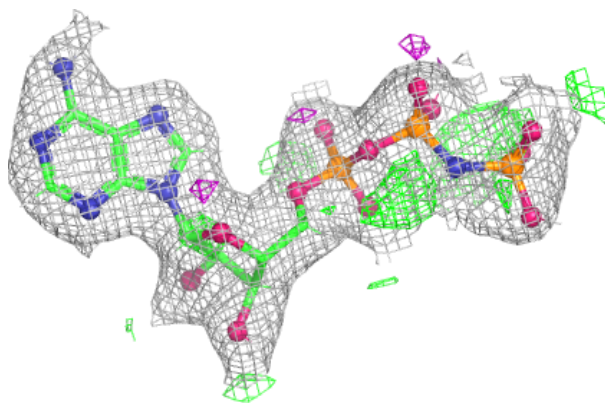
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	MLI	A	504	7/7	0.89	0.19	57,87,110,113	9
2	MG	B	501	1/1	0.91	0.22	69,69,69,69	0
4	MLI	A	503	7/7	0.92	0.08	43,50,60,64	0
3	ANP	B	502	31/31	0.93	0.10	40,45,88,121	0
3	ANP	A	502	31/31	0.95	0.09	37,46,73,111	0
4	MLI	B	503	7/7	0.95	0.06	43,50,54,57	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 ANP A 502:**

$2mF_o-DF_c$  (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.