

wwPDB X-ray Structure Validation Summary Report (i)

Sep 24, 2023 – 08:42 AM EDT

PDB ID	:	5UMB
Title	:	Crystal structure of ATPase domain of Malaria GRP78 with ADP bound
Authors	:	Chen, Y.; Antoshchenko, T.; Pizarro, J.C.; Song, J.H.; Park, H.
Deposited on		
Resolution	:	2.30 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 A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

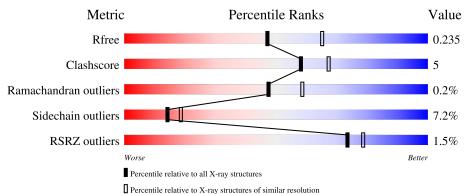
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7(2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.30 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	5042(2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575(2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 <=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	А	378	85%	14%	•
1	В	378	% 85%	14%	•
1	С	378	2% 86 %	12%	•
1	D	378	87%	11%	•

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard



residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PO4	D	502	-	Х	-	-



5UMB

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 12437 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Λ	378	Total	С	Ν	Ο	S	0	0	0
	А	510	2957	1875	506	571	5	0	0	0
1	В	378	Total	С	Ν	Ο	S	0	0	0
	D	510	2960	1876	506	573	5	0	0	0
1	С	377	Total	С	Ν	0	S	0	0	0
	U	511	2951	1870	504	572	5	0	0	0
1	Л	377	Total	С	Ν	0	S	0	0	0
	D	511	2947	1869	505	568	5	0	U	0

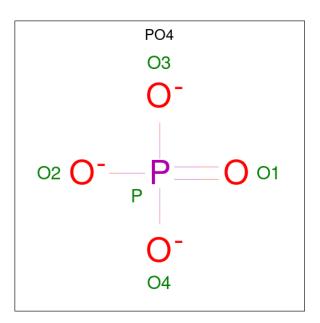
• Molecule 1 is a protein called Chaperone DnaK.

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	Total Mg 1 1	0	0
2	В	1	Total Mg 1 1	0	0
2	С	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0

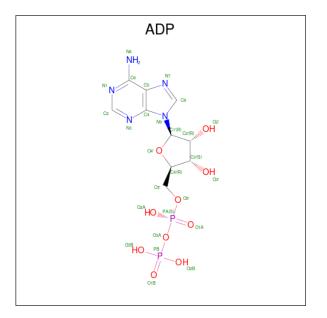
• Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0
3	В	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0
3	С	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0
3	D	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0

• Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).





Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
4	А	1	Total	С	Ν	Ο	Р	0	0	
4	A	1	27	10	5	10	2	0	0	
4	В	1	Total	С	Ν	Ο	Р	0	0	
4	D	1	27	10	5	10	2	0	U	
4	С	1	Total	С	Ν	Ο	Р	0	0	
4	U	1	27	10	5	10	2	0	U	
4	Л	1	Total	С	Ν	Ο	Р	0	0	
4	D	1	27	10	5	10	2	0	U	

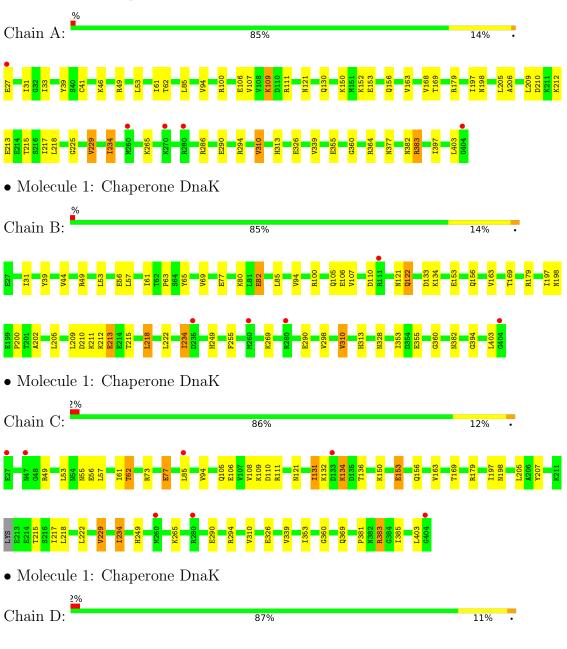
• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	84	Total O 84 84	0	0
5	В	133	Total O 133 133	0	0
5	С	111	Total O 111 111	0	0
5	D	162	Total O 162 162	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: Chaperone DnaK



R265 CLU M260 739 K268 739 K268 739 K268 739 K268 846 K369 648 K310 161 W310 165 W310 166 W313 185 K364 198 K364 198 K364 198 K369 114 K179 133 K160 133 K160 133 K160 133 K179 133 K160 133 K179 133 K179 136 K179 136 K179 136 K179 136 K179 136 K179 136 K169 136 K179 136 K169 136 K179 136 K169 136 </tr



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	84.53Å 112.26Å 93.16Å	Depositor
a, b, c, α , β , γ	90.00° 90.13° 90.00°	Depositor
Resolution (Å)	48.08 - 2.30	Depositor
Resolution (A)	48.08 - 2.30	EDS
% Data completeness	99.3 (48.08-2.30)	Depositor
(in resolution range)	99.0(48.08-2.30)	EDS
R _{merge}	(Not available)	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.92 (at 2.29 \text{\AA})$	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
P. P.	0.227 , 0.265	Depositor
R, R_{free}	0.211 , 0.235	DCC
R_{free} test set	3841 reflections $(5.01%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	24.0	Xtriage
Anisotropy	0.999	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.32 , 16.5	EDS
L-test for twinning ²	$< L >=0.48, < L^2>=0.31$	Xtriage
Estimated twinning fraction	0.299 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12437	wwPDB-VP
Average B, all atoms $(Å^2)$	33.0	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 34.44 % 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 6.8264e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

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.



¹Intensities estimated from amplitudes.

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: PO4, 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		
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.57	0/3000	0.69	0/4047	
1	В	0.51	0/3003	0.69	0/4051	
1	С	0.50	0/2993	0.69	0/4037	
1	D	0.53	0/2990	0.68	0/4034	
All	All	0.53	0/11986	0.69	0/16169	

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 (i)

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	А	2957	0	3009	34	0
1	В	2960	0	3011	33	0
1	С	2951	0	2997	33	0
1	D	2947	0	3001	26	0
2	А	1	0	0	0	0
2	В	1	0	0	0	0
2	С	1	0	0	0	0
2	D	1	0	0	0	0
3	А	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	В	5	0	0	0	0
3	С	5	0	0	0	0
3	D	5	0	0	0	0
4	А	27	0	12	2	0
4	В	27	0	12	0	0
4	С	27	0	12	1	0
4	D	27	0	12	2	0
5	А	84	0	0	1	0
5	В	133	0	0	4	0
5	С	111	0	0	0	0
5	D	162	0	0	6	0
All	All	12437	0	12066	115	0

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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 5.

The worst 5 of 115 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:D:174:PHE:HB3	5:D:608:HOH:O	1.46	1.12
1:B:63:PRO:HB2	5:B:604:HOH:O	1.51	1.10
1:D:179:ARG:HG3	5:D:608:HOH:O	1.59	1.02
1:D:209:LEU:HD21	1:D:355:GLU:HG2	1.42	0.97
1:B:80:LYS:HB2	5:B:604:HOH:O	1.72	0.87

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	Analysed Favoured Allowe		Outliers	Percentiles	
1	А	376/378~(100%)	368~(98%)	8 (2%)	0	100 100	

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Mol	Chain	Analysed	Favoured	Favoured Allowed		Percentiles		
1	В	376/378~(100%)	369~(98%)	6(2%)	1 (0%)	41	50	
1	С	373/378~(99%)	365~(98%)	8 (2%)	0	100	100	
1	D	375/378~(99%)	368~(98%)	5 (1%)	2(0%)	29	35	
All	All	1500/1512~(99%)	1470 (98%)	27 (2%)	3 (0%)	47	58	

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All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	47	ASN
1	D	213	GLU
1	В	213	GLU

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric Outlier		Percentiles		
1	А	321/322~(100%)	297~(92%)	24 (8%)	13	17	
1	В	322/322~(100%)	297~(92%)	25~(8%)	12	16	
1	С	321/322~(100%)	295~(92%)	26 (8%)	11	15	
1	D	320/322~(99%)	302~(94%)	18 (6%)	21	29	
All	All	1284/1288~(100%)	1191 (93%)	93~(7%)	14	18	

 $5~{\rm of}~93$ residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	С	105	GLN
1	С	326	GLU
1	С	109	LYS
1	С	215	THR
1	D	46	LYS

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 35 such side chains are listed below:



Mol	Chain	Res	Type
1	D	121	ASN
1	D	126	ASN
1	D	198	ASN
1	В	165	ASN
1	В	156	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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	Bo	ond leng	ths	Bond angles		
MOI	Type	Ullalli	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
4	ADP	D	503	2	24,29,29	0.91	1 (4%)	$29,\!45,\!45$	1.19	4 (13%)
3	PO4	В	502	2	4,4,4	2.65	3 (75%)	6,6,6	0.97	0
4	ADP	В	503	2	24,29,29	0.79	0	29,45,45	1.14	2 (6%)
3	PO4	С	502	2	4,4,4	2.48	1 (25%)	$6,\!6,\!6$	0.79	0
4	ADP	С	503	2	24,29,29	0.86	1 (4%)	29,45,45	0.89	2 (6%)
3	PO4	А	502	-	4,4,4	2.03	1 (25%)	6,6,6	0.86	0
3	PO4	D	502	2	4,4,4	2.59	2 (50%)	6,6,6	1.65	2 (33%)
4	ADP	А	503	2	24,29,29	0.72	0	29,45,45	0.91	2(6%)



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	D	503	2	-	1/12/32/32	0/3/3/3
4	ADP	В	503	2	-	0/12/32/32	0/3/3/3
4	ADP	С	503	2	-	1/12/32/32	0/3/3/3
4	ADP	А	503	2	-	0/12/32/32	0/3/3/3

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\operatorname{Observed}(\operatorname{\AA})$	$\mathrm{Ideal}(\mathrm{\AA})$
3	В	502	PO4	P-01	4.09	1.60	1.50
3	С	502	PO4	P-01	4.01	1.60	1.50
3	D	502	PO4	P-01	3.96	1.60	1.50
4	D	503	ADP	PB-O2B	-2.69	1.44	1.54
3	А	502	PO4	P-01	2.35	1.56	1.50

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
4	D	503	ADP	PA-O3A-PB	3.43	144.61	132.83
3	D	502	PO4	O4-P-O1	-2.95	100.10	110.89
4	D	503	ADP	O2B-PB-O3A	2.88	114.30	104.64
3	D	502	PO4	O4-P-O2	2.55	116.15	107.97
4	В	503	ADP	PA-O3A-PB	2.41	141.11	132.83

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	С	503	ADP	C3'-C4'-C5'-O5'
4	D	503	ADP	PB-O3A-PA-O1A

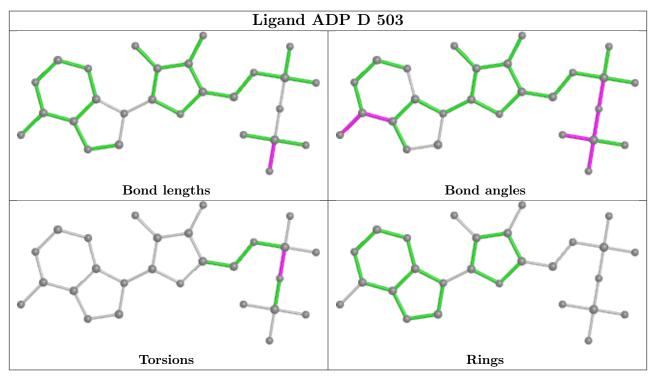
There are no ring outliers.

3 monomers are involved in 5 short contacts:

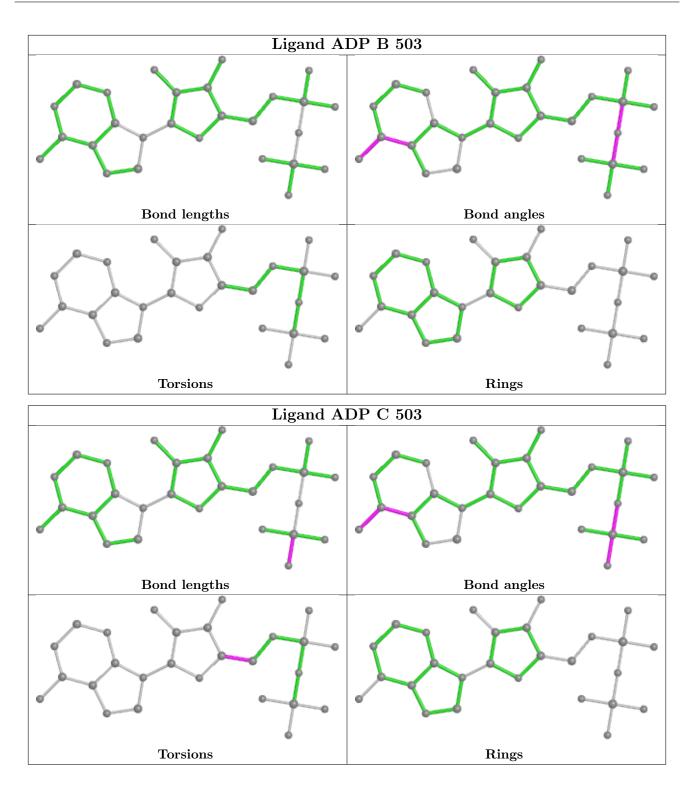
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	503	ADP	2	0
4	С	503	ADP	1	0
4	А	503	ADP	2	0



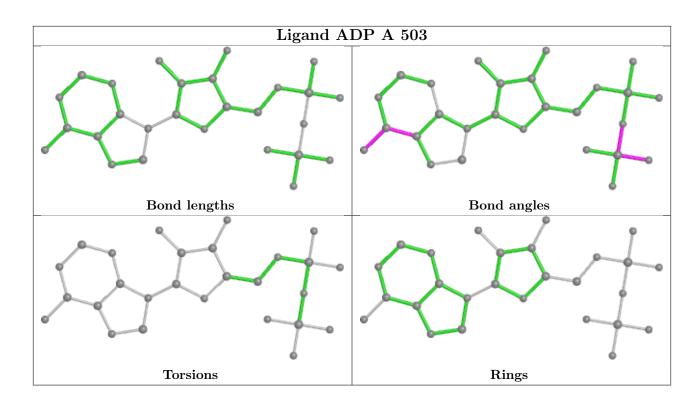
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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 sufficient 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^{th} 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	$\langle RSRZ \rangle$	# RSRZ > 2	$OWAB(Å^2)$	Q<0.9
1	А	378/378~(100%)	-0.19	5 (1%) 77 81	12, 29, 60, 73	5 (1%)
1	В	378/378~(100%)	-0.18	5 (1%) 77 81	14, 30, 57, 82	5(1%)
1	С	377/378~(99%)	-0.16	7 (1%) 66 73	13, 30, 58, 73	5 (1%)
1	D	377/378~(99%)	-0.14	6 (1%) 72 77	14, 28, 57, 78	5 (1%)
All	All	1510/1512~(99%)	-0.17	23 (1%) 73 79	12, 29, 58, 82	20 (1%)

The worst 5 of 23 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	404	GLY	4.4
1	D	280	ARG	3.8
1	С	404	GLY	3.7
1	D	260	MET	3.6
1	В	280	ARG	3.5

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.

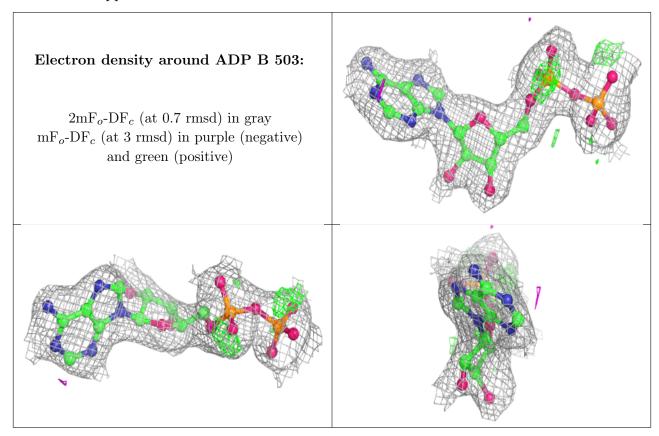
6.4 Ligands (i)

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^{th} 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.

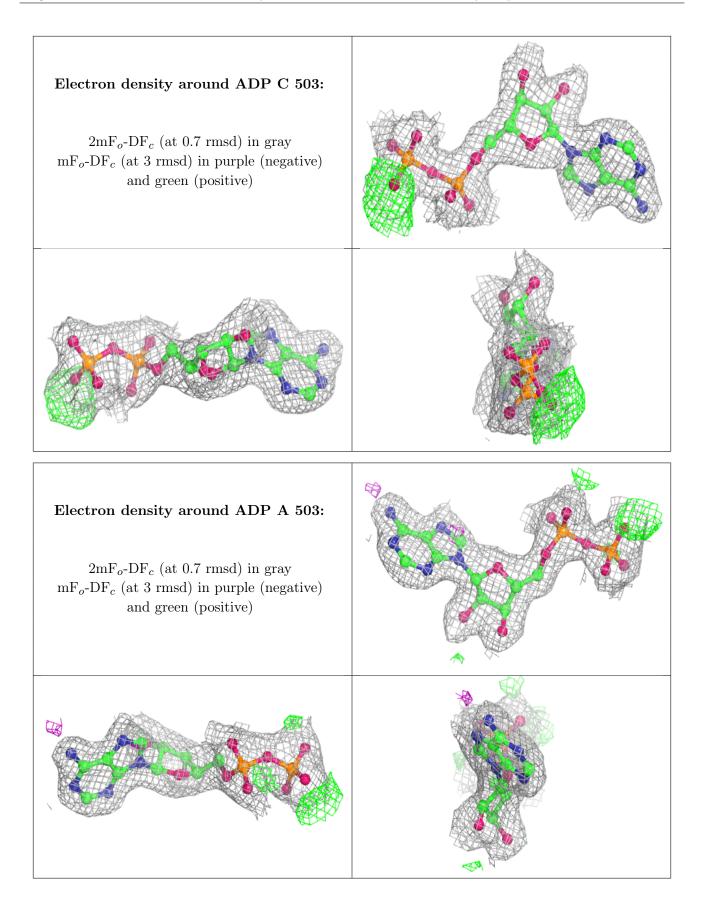


Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	$Q{<}0.9$
2	MG	А	501	1/1	0.88	0.11	22,22,22,22	0
2	MG	С	501	1/1	0.89	0.13	26, 26, 26, 26	0
3	PO4	С	502	5/5	0.94	0.17	$25,\!29,\!29,\!29$	0
2	MG	В	501	1/1	0.95	0.10	21,21,21,21	0
4	ADP	В	503	27/27	0.96	0.14	24,35,43,45	0
4	ADP	С	503	27/27	0.96	0.11	8,26,31,32	0
3	PO4	D	502	5/5	0.97	0.10	8,12,13,21	0
4	ADP	А	503	27/27	0.97	0.11	$6,\!15,\!22,\!26$	0
4	ADP	D	503	27/27	0.97	0.12	$10,\!18,\!22,\!27$	0
2	MG	D	501	1/1	0.98	0.08	3,3,3,3	0
3	PO4	А	502	5/5	0.98	0.09	15,18,24,24	0
3	PO4	В	502	5/5	0.98	0.11	4,10,15,22	0

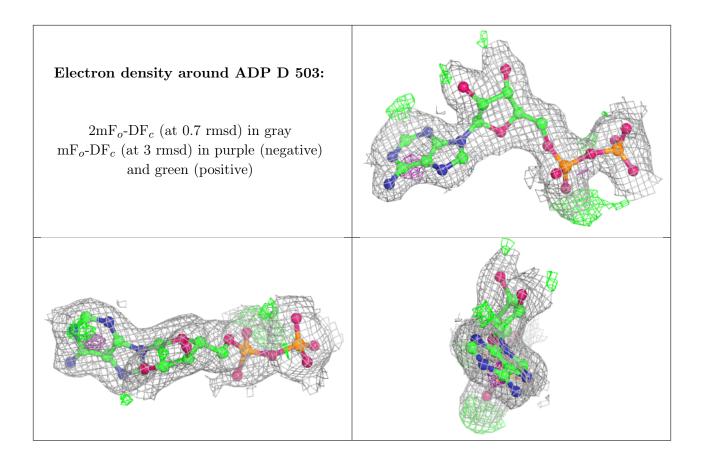
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.











6.5 Other polymers (i)

There are no such residues in this entry.

