



# Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 07:54 am BST

PDB ID : 6SR6  
Title : Crystal structure of the RAC core with a pseudo substrate bound to Ssz1 SBD  
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Deposited on : 2019-09-05  
Resolution : 2.50 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 : 2.11  
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

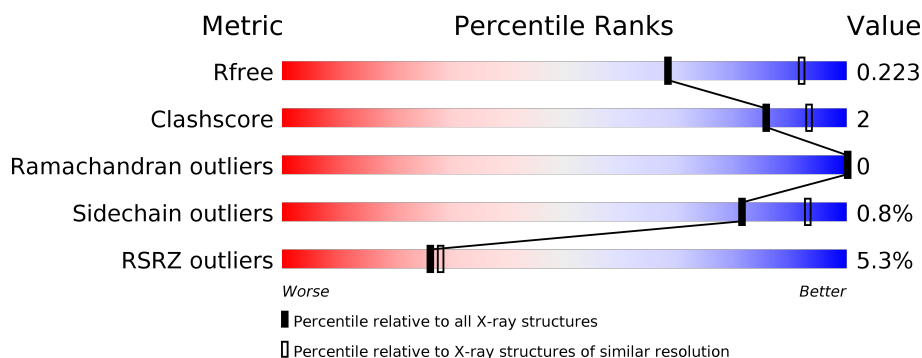
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	589	<div> <div>4%</div> <div>80% 8% 12%</div> </div>
1	C	589	<div> <div>4%</div> <div>84% 5% 11%</div> </div>
2	B	63	<div> <div>16%</div> <div>75% 8% 17%</div> </div>
2	D	63	<div> <div>10%</div> <div>79% 6% 14%</div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8887 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 Putative heat shock protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	519	Total	C	N	O	S	0	0	0
			3942	2470	689	780	3			
1	C	524	Total	C	N	O	S	0	0	0
			3974	2488	696	787	3			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	SER	-	expression tag	UNP G0RZX9
A	-9	ALA	-	expression tag	UNP G0RZX9
A	-8	MET	-	expression tag	UNP G0RZX9
A	-7	GLY	-	expression tag	UNP G0RZX9
A	-6	TRP	-	expression tag	UNP G0RZX9
A	-5	SER	-	expression tag	UNP G0RZX9
A	-4	HIS	-	expression tag	UNP G0RZX9
A	-3	PRO	-	expression tag	UNP G0RZX9
A	-2	GLN	-	expression tag	UNP G0RZX9
A	-1	PHE	-	expression tag	UNP G0RZX9
A	0	GLU	-	expression tag	UNP G0RZX9
A	1	LYS	-	expression tag	UNP G0RZX9
C	-10	SER	-	expression tag	UNP G0RZX9
C	-9	ALA	-	expression tag	UNP G0RZX9
C	-8	MET	-	expression tag	UNP G0RZX9
C	-7	GLY	-	expression tag	UNP G0RZX9
C	-6	TRP	-	expression tag	UNP G0RZX9
C	-5	SER	-	expression tag	UNP G0RZX9
C	-4	HIS	-	expression tag	UNP G0RZX9
C	-3	PRO	-	expression tag	UNP G0RZX9
C	-2	GLN	-	expression tag	UNP G0RZX9
C	-1	PHE	-	expression tag	UNP G0RZX9
C	0	GLU	-	expression tag	UNP G0RZX9
C	1	LYS	-	expression tag	UNP G0RZX9

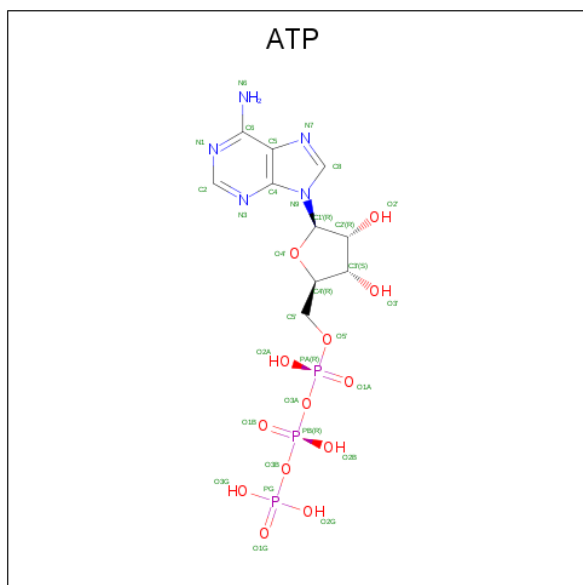
- Molecule 2 is a protein called Putative ribosome associated protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	52	Total	C	N	O	S	0	0	0
			399	255	73	70	1			
2	D	54	Total	C	N	O	S	0	0	0
			417	265	78	73	1			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	GLY	-	expression tag	UNP G0RYD6
B	-1	PRO	-	expression tag	UNP G0RYD6
B	0	ALA	-	expression tag	UNP G0RYD6
D	-2	GLY	-	expression tag	UNP G0RYD6
D	-1	PRO	-	expression tag	UNP G0RYD6
D	0	ALA	-	expression tag	UNP G0RYD6

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula:  $C_{10}H_{16}N_5O_{13}P_3$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

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

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

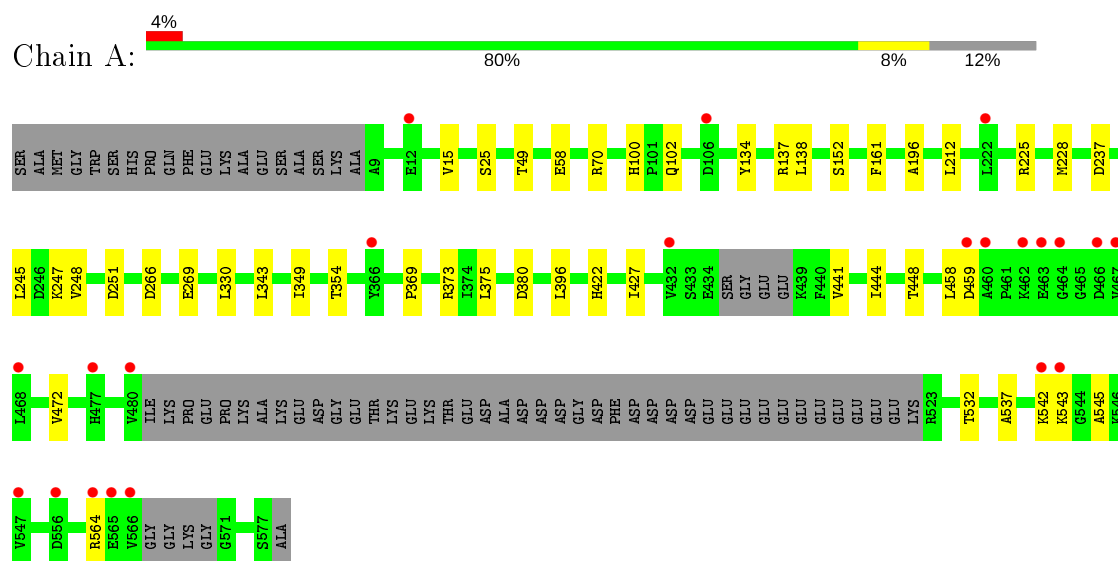
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	27	Total 27	O 27	0	0
5	C	62	Total 62	O 62	0	0
5	D	2	Total 2	O 2	0	0

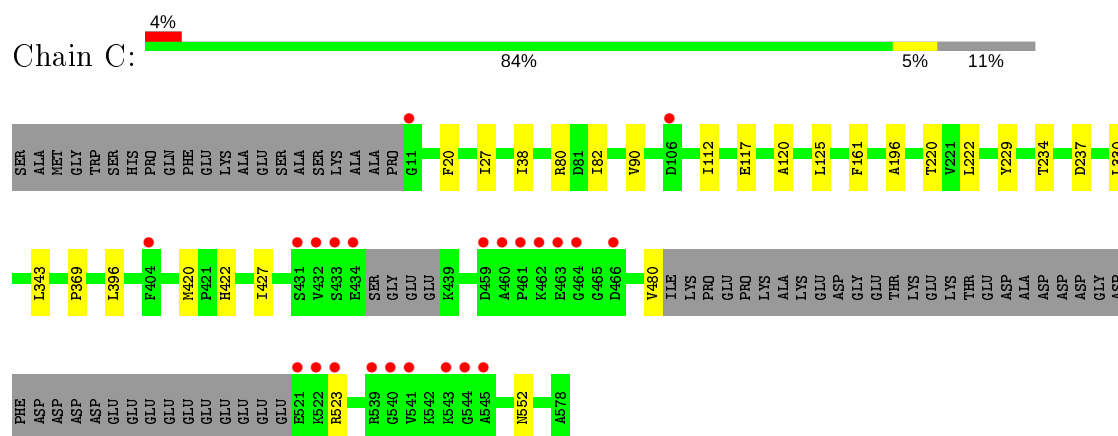
### 3 Residue-property plots [i](#)

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: Putative heat shock protein

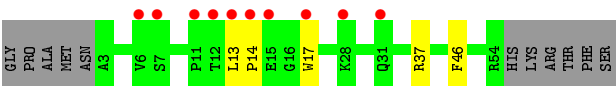


- Molecule 1: Putative heat shock protein

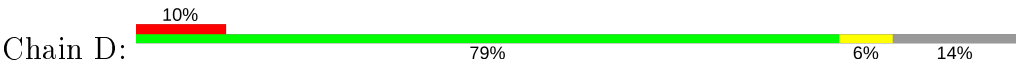


- Molecule 2: Putative ribosome associated protein





● Molecule 2: Putative ribosome associated protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	51.98Å 258.45Å 53.02Å 90.00° 100.09° 90.00°	Depositor
Resolution (Å)	33.71 – 2.50 48.40 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.8 (33.71-2.50) 98.9 (48.40-2.50)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.37 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, $R_{free}$	0.189 , 0.223 0.189 , 0.223	Depositor DCC
$R_{free}$ test set	2292 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.0	Xtriage
Anisotropy	0.473	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 35.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.031 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8887	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% 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, ATP

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	0.25	0/4002	0.43	0/5437
1	C	0.25	0/4034	0.43	0/5477
2	B	0.25	0/409	0.42	0/555
2	D	0.26	0/428	0.42	0/581
All	All	0.25	0/8873	0.43	0/12050

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	A	3942	0	3967	24	0
1	C	3974	0	4002	13	0
2	B	399	0	410	4	0
2	D	417	0	423	3	0
3	A	31	0	12	0	0
3	C	31	0	12	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	A	27	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	62	0	0	0	0
5	D	2	0	0	0	0
All	All	8887	0	8826	39	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 (39) 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:C:220:THR:HG23	1:C:234:THR:HG22	1.67	0.77
2:B:13:LEU:HD12	2:B:14:PRO:HD2	1.79	0.64
1:C:222:LEU:HD23	1:C:229:TYR:HB3	1.82	0.60
1:A:212:LEU:HB3	1:A:354:THR:HB	1.84	0.60
1:A:228:MET:HE3	2:B:46:PHE:HA	1.84	0.59
1:C:343:LEU:HD23	1:C:369:PRO:HG3	1.85	0.59
1:A:196:ALA:HB3	1:A:396:LEU:HD13	1.86	0.57
1:A:373:ARG:NH1	1:A:380:ASP:OD2	2.36	0.57
1:C:27:ILE:HG13	1:C:38:ILE:HB	1.87	0.57
1:C:480:VAL:HG22	1:C:523:ARG:HG2	1.88	0.56
1:C:196:ALA:HB3	1:C:396:LEU:HD13	1.87	0.55
1:C:112:ILE:HD11	1:C:125:LEU:HD12	1.90	0.54
1:A:343:LEU:HD23	1:A:369:PRO:HG3	1.91	0.52
1:A:349:ILE:HG12	1:A:375:LEU:HD12	1.93	0.51
1:A:58:GLU:O	1:A:137:ARG:NH2	2.45	0.50
1:A:15:VAL:HG12	1:A:152:SER:HB2	1.96	0.48
1:A:237:ASP:HB3	1:A:330:LEU:HD11	1.96	0.48
1:C:117:GLU:HG2	1:C:120:ALA:HB2	1.96	0.48
1:C:237:ASP:HB2	1:C:330:LEU:HD21	1.97	0.47
1:A:237:ASP:HB2	1:A:330:LEU:HD21	1.96	0.47
1:A:247:LYS:NZ	1:A:251:ASP:OD2	2.49	0.46
1:C:20:PHE:O	1:C:80:ARG:NH1	2.48	0.46
1:A:427:ILE:HB	1:A:444:ILE:HB	1.99	0.45
1:A:49:THR:HA	1:A:134:TYR:CZ	2.51	0.45
1:C:427:ILE:HD11	2:D:29:LEU:HD11	1.98	0.44
1:C:422:HIS:CE1	2:D:37:ARG:HB2	2.52	0.44
1:A:542:LYS:HG2	1:A:543:LYS:N	2.33	0.44
1:A:25:SER:HB3	1:A:138:LEU:HD11	1.99	0.43
1:C:82:ILE:HD12	1:C:90:VAL:HG13	2.00	0.43
1:A:441:VAL:HG11	1:A:458:LEU:HD21	1.99	0.43
1:A:422:HIS:CE1	2:B:37:ARG:HB2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:40:GLU:O	2:D:42:VAL:N	2.51	0.42
1:A:266:ASP:HB3	1:A:269:GLU:HG2	2.01	0.42
1:A:542:LYS:HB3	1:A:545:ALA:HB2	2.02	0.41
1:A:537:ALA:HB2	2:B:17:TRP:CZ3	2.55	0.41
1:A:245:LEU:O	1:A:248:VAL:HG22	2.21	0.41
1:A:70:ARG:NH2	5:A:801:HOH:O	2.18	0.41
1:A:472:VAL:HG12	1:A:532:THR:HG22	2.03	0.41
1:A:225:ARG:HG3	1:A:448:THR:HA	2.03	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	511/589 (87%)	502 (98%)	9 (2%)	0	100	100
1	C	518/589 (88%)	512 (99%)	6 (1%)	0	100	100
2	B	50/63 (79%)	50 (100%)	0	0	100	100
2	D	52/63 (82%)	49 (94%)	3 (6%)	0	100	100
All	All	1131/1304 (87%)	1113 (98%)	18 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	427/483 (88%)	422 (99%)	5 (1%)	71	88
1	C	429/483 (89%)	426 (99%)	3 (1%)	84	94
2	B	42/51 (82%)	42 (100%)	0	100	100
2	D	44/51 (86%)	44 (100%)	0	100	100
All	All	942/1068 (88%)	934 (99%)	8 (1%)	81	93

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	HIS
1	A	102	GLN
1	A	161	PHE
1	A	459	ASP
1	A	564	ARG
1	C	161	PHE
1	C	420	MET
1	C	552	ASN

Some 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 carbohydrates 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
3	ATP	C	700	4	26,33,33	0.92	1 (3%)	31,52,52	1.31	2 (6%)
3	ATP	A	700	4	26,33,33	0.92	1 (3%)	31,52,52	1.40	4 (12%)

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	ATP	C	700	4	-	5/18/38/38	0/3/3/3
3	ATP	A	700	4	-	2/18/38/38	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	700	ATP	C5-C4	2.45	1.47	1.40
3	C	700	ATP	C5-C4	2.33	1.47	1.40

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	700	ATP	N3-C2-N1	-3.33	123.47	128.68
3	C	700	ATP	N3-C2-N1	-3.29	123.54	128.68
3	A	700	ATP	C4-C5-N7	-2.67	106.61	109.40
3	C	700	ATP	C4-C5-N7	-2.59	106.70	109.40
3	A	700	ATP	PB-O3B-PG	-2.54	124.11	132.83
3	A	700	ATP	PA-O3A-PB	-2.13	125.51	132.83

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	700	ATP	PB-O3A-PA-O1A
3	C	700	ATP	PB-O3B-PG-O3G
3	C	700	ATP	PB-O3A-PA-O2A

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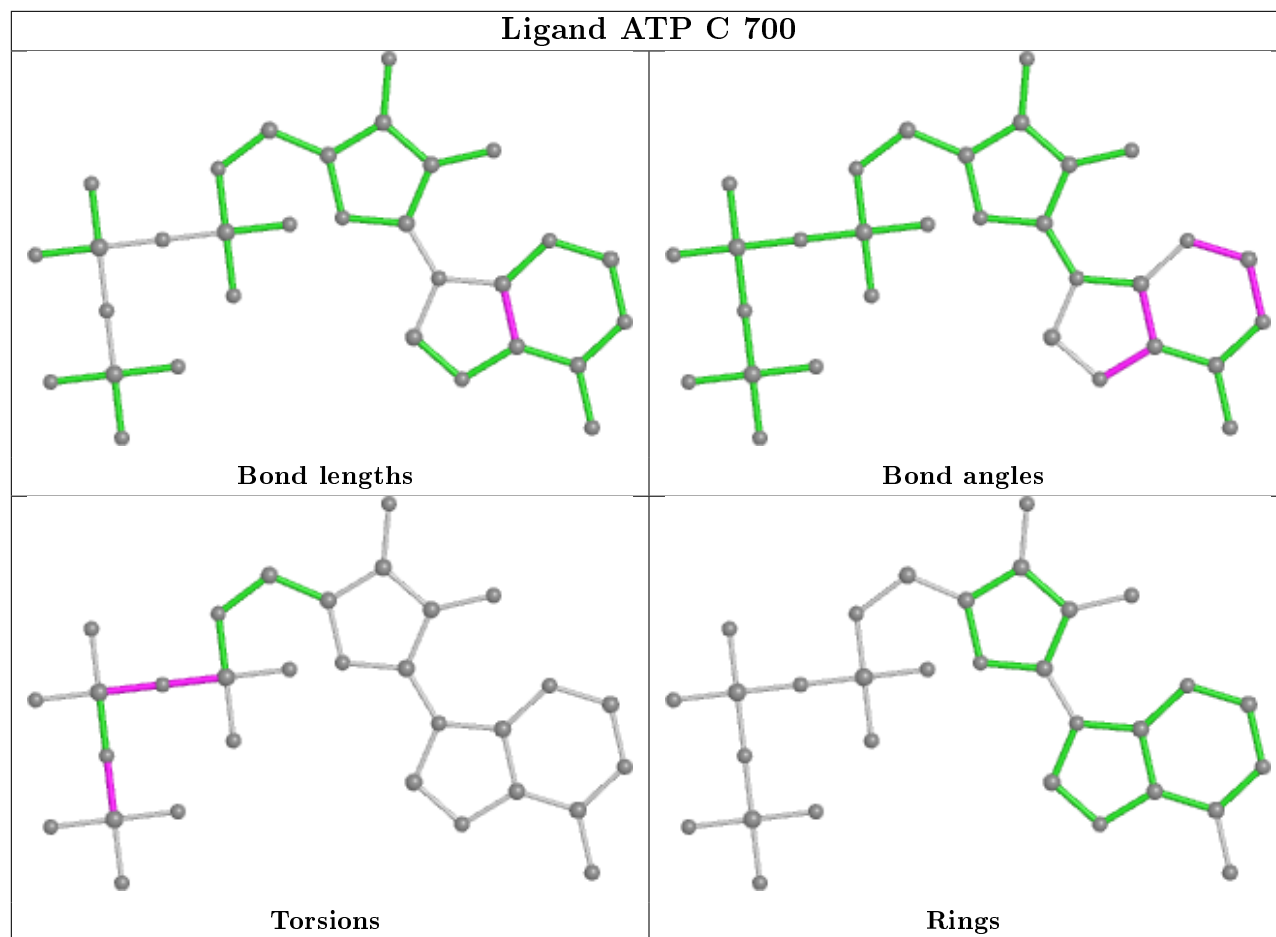
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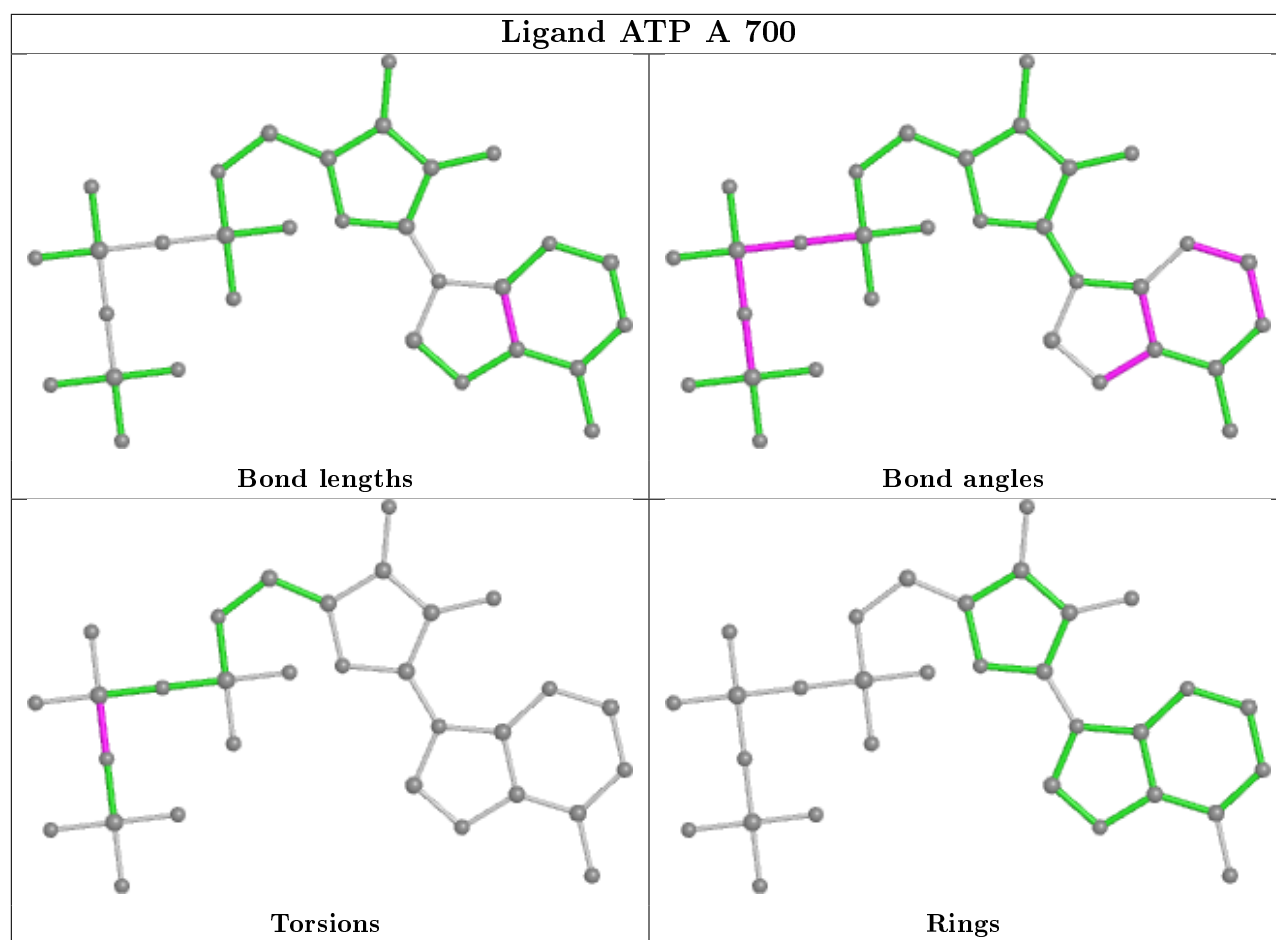
Mol	Chain	Res	Type	Atoms
3	C	700	ATP	PA-O3A-PB-O1B
3	C	700	ATP	PA-O3A-PB-O2B
3	A	700	ATP	PG-O3B-PB-O1B
3	A	700	ATP	PG-O3B-PB-O2B

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 [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	519/589 (88%)	0.21	22 (4%) 36 39	25, 49, 97, 147	0
1	C	524/589 (88%)	0.26	23 (4%) 34 37	22, 40, 94, 133	0
2	B	52/63 (82%)	0.93	10 (19%) 1 1	41, 80, 114, 122	0
2	D	54/63 (85%)	0.70	6 (11%) 5 5	34, 76, 101, 113	0
All	All	1149/1304 (88%)	0.29	61 (5%) 26 28	22, 47, 99, 147	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	432	VAL	5.7
1	C	463	GLU	4.9
1	C	462	LYS	4.9
1	C	11	GLY	4.6
2	B	17	TRP	4.2
1	C	433	SER	3.9
1	C	460	ALA	3.8
1	A	556	ASP	3.8
1	C	461	PRO	3.8
1	C	522	LYS	3.6
1	A	460	ALA	3.6
1	C	521	GLU	3.5
1	A	463	GLU	3.5
1	A	564	ARG	3.5
1	A	480	VAL	3.4
2	D	55	HIS	3.4
2	B	11	PRO	3.3
2	D	12	THR	3.3
2	D	15	GLU	3.3
1	A	547	VAL	3.3
1	A	432	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
2	B	15	GLU	3.2
1	A	467	VAL	3.1
1	C	431	SER	3.1
1	C	545	ALA	3.1
1	A	566	VAL	3.1
1	A	565	GLU	3.0
1	A	464	GLY	3.0
1	C	544	GLY	2.9
1	C	543	LYS	2.9
1	A	462	LYS	2.8
2	D	54	ARG	2.8
2	B	12	THR	2.8
2	D	2	ASN	2.8
1	C	464	GLY	2.7
2	B	13	LEU	2.7
2	B	6	VAL	2.7
1	C	106	ASP	2.7
1	A	468	LEU	2.6
1	A	466	ASP	2.6
1	A	543	LYS	2.6
1	C	541	VAL	2.5
1	C	459	ASP	2.5
1	A	477	HIS	2.5
1	A	12	GLU	2.5
1	C	540	GLY	2.4
1	C	523	ARG	2.4
1	C	466	ASP	2.4
1	C	539	ARG	2.4
1	C	434	GLU	2.3
1	A	106	ASP	2.2
2	B	31	GLN	2.1
1	A	459	ASP	2.1
2	B	7	SER	2.1
2	D	11	PRO	2.1
1	C	404	PHE	2.0
2	B	14	PRO	2.0
2	B	28	LYS	2.0
1	A	222	LEU	2.0
1	A	366	TYR	2.0
1	A	542	LYS	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

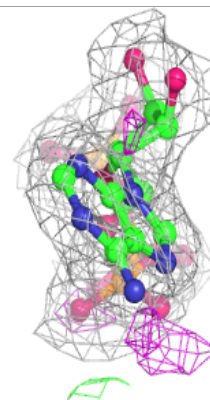
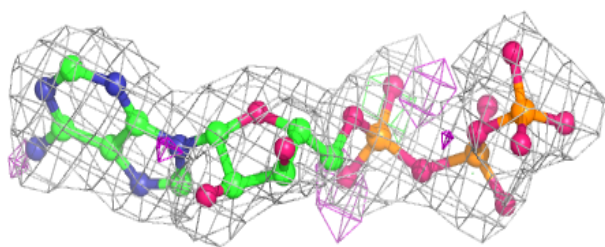
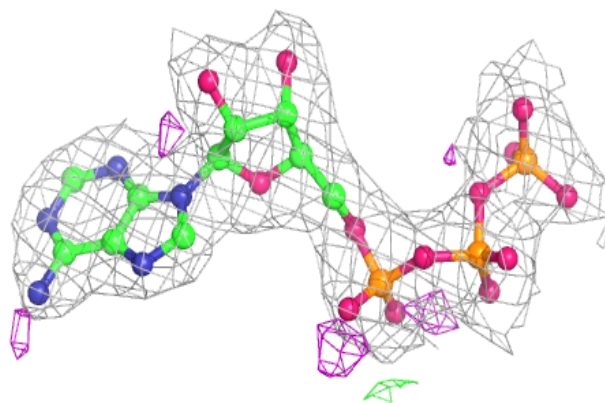
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ATP	C	700	31/31	0.97	0.17	19,31,37,42	0
4	MG	C	701	1/1	0.98	0.16	37,37,37,37	0
3	ATP	A	700	31/31	0.98	0.17	26,38,42,45	0
4	MG	A	701	1/1	0.99	0.07	30,30,30,30	0

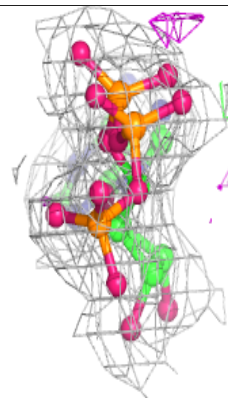
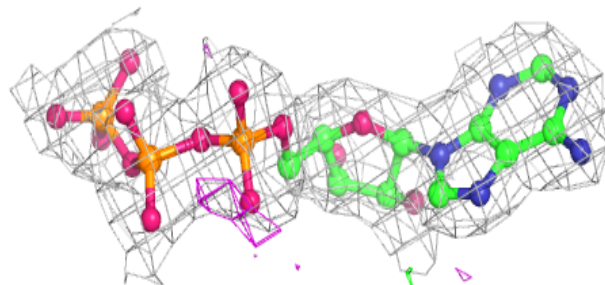
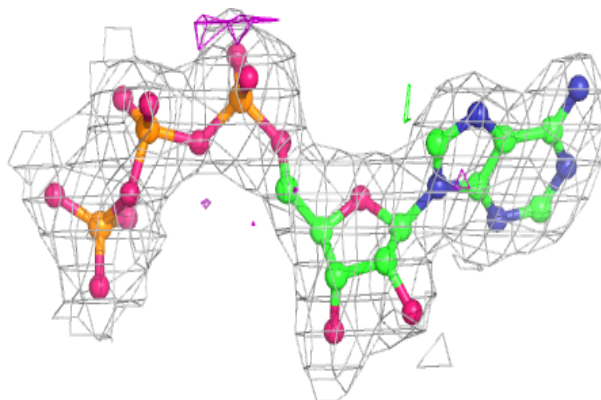
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around ATP C 700:**

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

**Electron density around ATP A 700:**

$2mF_o-DF_c$  (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.