



# Full wwPDB X-ray Structure Validation Report

Oct 9, 2014 – 07:31 AM EDT

PDB ID : 4NTT  
Title : Structure of the catalytic subunit of cAMP-dependent protein kinase bound to ADP and one magnesium ion  
Authors : Bastidas, A.C.; Wu, J.; Taylor, S.S.  
Deposited on : 2013-12-02  
Resolution : 3.50 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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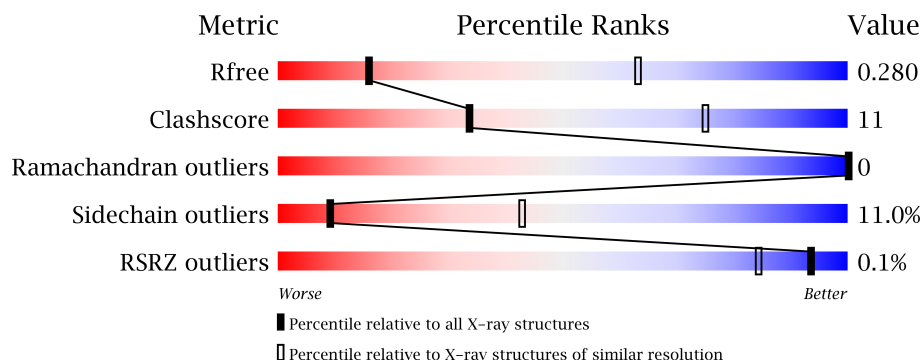
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.16 November 2013  
Xtriage (Phenix) : dev-1439  
EDS : stable23828  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable23828

# 1 Overall quality at a glance

The reported resolution of this entry is 3.50 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))
$R_{free}$	66092	1243 (3.70-3.30)
Clashscore	79885	1039 (3.66-3.34)
Ramachandran outliers	78287	1000 (3.66-3.34)
Sidechain outliers	78261	1000 (3.66-3.34)
RSRZ outliers	66119	1243 (3.70-3.30)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	350	
1	B	350	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	MG	B	402	-	X

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 5262 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

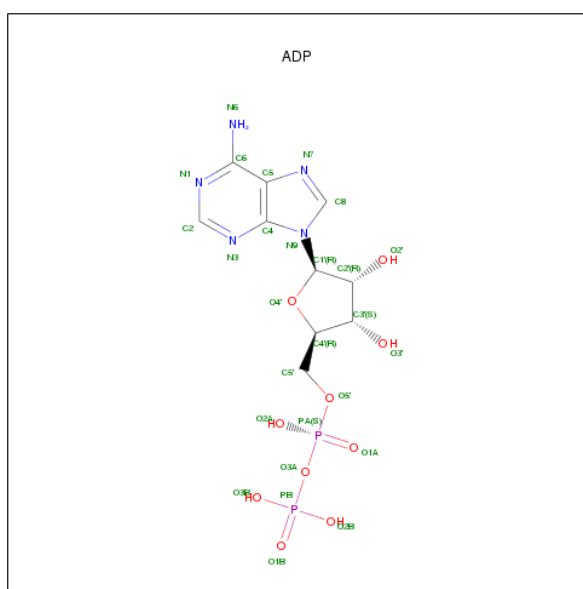
- Molecule 1 is a protein called cAMP-dependent protein kinase catalytic subunit alpha.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	P	S	0	0	0
			2597	1676	439	473	2	7			
1	B	341	Total	C	N	O	P	S	0	0	0
			2609	1688	439	472	2	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	CYS	LYS	EXPRESSION TAG	UNP P05132
B	7	CYS	LYS	ENGINEERED MUTATION	UNP P05132

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



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

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
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).

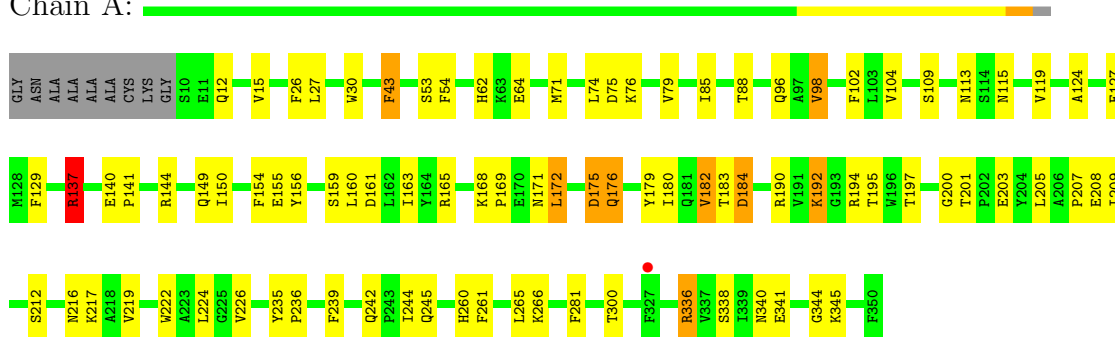
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		

### 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 errors 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.

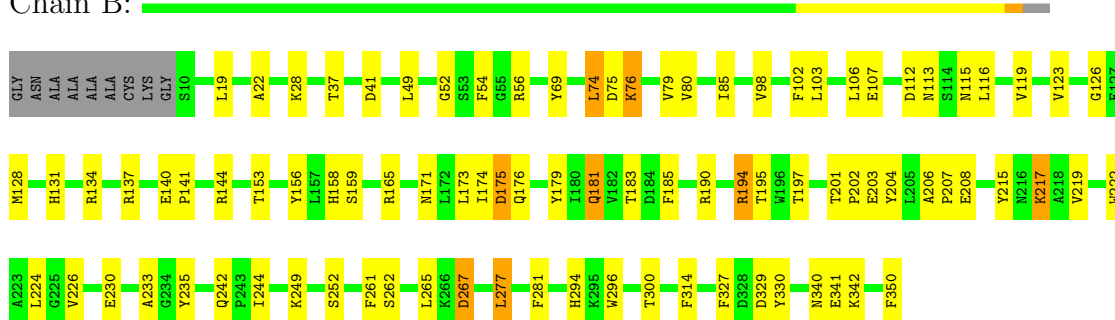
- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha

Chain A:



- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	50.48Å 143.15Å 62.57Å 90.00° 103.61° 90.00°	Depositor
Resolution (Å)	71.58 – 3.50 32.15 – 3.50	Depositor EDS
% Data completeness (in resolution range)	97.1 (71.58-3.50) 97.3 (32.15-3.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.22 (at 3.47Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.227 , 0.284 0.223 , 0.280	Depositor DCC
$R_{free}$ test set	516 reflections (5.10%)	DCC
Wilson B-factor (Å <sup>2</sup> )	95.3	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 42.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 10630 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5262	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

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

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	0.42	1/2641 (0.0%)	0.54	0/3590
1	B	0.41	0/2655	0.55	0/3610
All	All	0.42	1/5296 (0.0%)	0.55	0/7200

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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	137	ARG	CZ-NH2	6.73	1.41	1.33

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	182	VAL	Peptide

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit,

and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2597	0	2327	58	0
1	B	2609	0	2347	52	0
2	A	27	0	12	2	0
2	B	27	0	12	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
All	All	5262	0	4698	110	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (110) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:171:ASN:C	1:A:183:THR:HG22	1.70	1.10
1:A:171:ASN:HB3	1:A:183:THR:HG21	1.40	1.02
1:B:126:GLY:O	1:B:173:LEU:HA	1.67	0.94
1:A:12:GLN:O	1:A:15:VAL:HG22	1.74	0.87
1:A:171:ASN:HB3	1:A:183:THR:CG2	2.05	0.87
1:B:76:LYS:HG2	1:B:115:ASN:HA	1.64	0.79
1:A:163:ILE:HG21	1:A:217:LYS:HA	1.64	0.79
1:A:171:ASN:O	1:A:183:THR:HG22	1.83	0.77
1:B:277:LEU:HD23	1:B:277:LEU:H	1.51	0.74
1:A:236:PRO:HG2	1:A:239:PHE:HB2	1.70	0.73
1:A:75:ASP:HA	1:A:115:ASN:HB3	1.71	0.72
1:B:75:ASP:HA	1:B:115:ASN:HB3	1.71	0.71
1:A:183:THR:OG1	1:A:184:ASP:N	2.19	0.71
1:A:124:ALA:HB3	1:A:176:GLN:HG2	1.74	0.69
1:B:102:PHE:O	1:B:181:GLN:HA	1.92	0.69
1:A:175:ASP:HB2	1:A:179:TYR:H	1.56	0.69
1:B:194:ARG:HH11	1:B:194:ARG:HG3	1.60	0.66
1:B:204:TYR:OH	1:B:230:GLU:OE2	2.09	0.65
1:A:175:ASP:HB2	1:A:179:TYR:N	2.11	0.64
1:B:103:LEU:HD22	1:B:185:PHE:HZ	1.63	0.63
1:A:137:ARG:HH22	1:A:260:HIS:CG	2.16	0.62
1:B:144:ARG:HD3	1:B:300:THR:HG23	1.80	0.62
1:B:156:TYR:O	1:B:159:SER:HB3	2.00	0.61
1:A:76:LYS:HG2	1:A:115:ASN:HA	1.83	0.60
1:A:171:ASN:C	1:A:183:THR:CG2	2.60	0.60
1:B:208:GLU:HB3	1:B:215:TYR:HB3	1.84	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:209:ILE:HD11	1:A:219:VAL:HG11	1.84	0.60
1:A:172:LEU:N	1:A:183:THR:HG22	2.15	0.59
1:B:165:ARG:HH12	1:B:197:TPO:HB	1.68	0.59
1:B:80:VAL:HG13	1:B:85:ILE:HD11	1.85	0.59
1:B:126:GLY:HA2	1:B:174:ILE:H	1.67	0.58
1:A:43:PHE:HB3	1:A:62:HIS:HA	1.85	0.57
1:B:54:PHE:HB2	1:B:74:LEU:HB2	1.87	0.56
1:A:137:ARG:NH1	1:A:260:HIS:CE1	2.75	0.55
1:A:161:ASP:HB2	1:A:190:ARG:HE	1.72	0.55
1:B:171:ASN:O	1:B:183:THR:HG22	2.06	0.55
1:B:113:ASN:HA	1:B:341:GLU:HG3	1.89	0.55
1:A:137:ARG:NH2	1:A:260:HIS:CD2	2.75	0.54
1:A:222:TRP:O	1:A:226:VAL:HG23	2.08	0.54
1:B:52:GLY:H	1:B:56:ARG:HA	1.73	0.54
1:B:267:ASP:OD1	1:B:294:HIS:NE2	2.40	0.54
1:B:144:ARG:HD3	1:B:300:THR:CG2	2.38	0.53
1:A:98:VAL:HG22	1:A:156:TYR:HE2	1.73	0.53
1:A:144:ARG:HD3	1:A:300:THR:HG23	1.91	0.53
1:A:163:ILE:CG2	1:A:217:LYS:HA	2.35	0.52
1:B:76:LYS:NZ	1:B:112:ASP:O	2.41	0.52
1:A:182:VAL:HG13	1:A:183:THR:HG23	1.92	0.52
1:B:75:ASP:O	1:B:79:VAL:HG23	2.10	0.51
1:B:54:PHE:HB2	1:B:74:LEU:CB	2.41	0.51
1:A:172:LEU:H	1:A:172:LEU:HD23	1.75	0.51
1:A:53:SER:OG	2:A:401:ADP:PB	2.68	0.51
1:B:69:TYR:CE2	1:B:107:GLU:HG3	2.46	0.51
1:A:102:PHE:O	1:A:182:VAL:HG23	2.11	0.51
1:A:137:ARG:NH2	1:A:260:HIS:CG	2.78	0.51
1:A:336:ARG:HH22	1:A:338:SEP:HB3	1.75	0.50
1:A:124:ALA:CB	1:A:176:GLN:HG2	2.41	0.50
1:A:168:LYS:O	1:A:171:ASN:HB2	2.12	0.49
1:A:26:PHE:CG	1:A:160:LEU:HD13	2.48	0.48
1:A:71:MET:HG3	1:A:119:VAL:HG22	1.95	0.48
1:B:265:LEU:HD23	1:B:296:TRP:CE2	2.49	0.48
1:B:249:LYS:O	1:B:252:SER:O	2.32	0.48
1:B:233:ALA:HB2	1:B:261:PHE:CZ	2.48	0.48
1:B:175:ASP:HB2	1:B:179:TYR:N	2.29	0.47
1:B:49:LEU:O	1:B:327:PHE:HD1	1.96	0.47
1:A:161:ASP:OD2	1:A:192:LYS:HE3	2.14	0.47
1:B:206:ALA:HB3	1:B:219:VAL:HG22	1.97	0.47
1:B:69:TYR:HE2	1:B:107:GLU:HG3	1.80	0.47
1:A:129:PHE:HB2	1:A:169:PRO:HB2	1.98	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:149:GLN:OE1	1:A:180:ILE:HB	2.16	0.46
1:A:27:LEU:HA	1:A:30:TRP:HB3	1.97	0.45
1:A:341:GLU:OE2	1:A:344:GLY:HA3	2.16	0.45
1:B:329:ASP:O	1:B:330:TYR:HB2	2.17	0.45
1:A:244:ILE:H	1:A:244:ILE:HD12	1.81	0.45
1:B:107:GLU:HB2	1:B:119:VAL:HG12	1.99	0.45
1:B:140:GLU:N	1:B:141:PRO:HD2	2.32	0.45
1:B:242:GLN:HB3	1:B:244:ILE:HD12	1.98	0.44
1:A:140:GLU:N	1:A:141:PRO:HD2	2.31	0.44
1:A:242:GLN:HB2	1:A:245:GLN:NE2	2.33	0.44
1:B:165:ARG:NH1	1:B:197:TPO:O1P	2.51	0.44
1:B:233:ALA:HB2	1:B:261:PHE:HZ	1.82	0.44
1:A:208:GLU:CD	1:A:208:GLU:H	2.21	0.44
1:B:131:HIS:O	1:B:134:ARG:HB3	2.18	0.44
1:B:175:ASP:HB2	1:B:179:TYR:H	1.82	0.43
1:B:222:TRP:O	1:B:226:VAL:HG23	2.18	0.43
1:A:165:ARG:NH1	1:A:197:TPO:O1P	2.50	0.42
1:B:116:LEU:HD22	1:B:350:PHE:CD2	2.54	0.42
1:B:140:GLU:OE1	1:B:262:SER:HB3	2.19	0.42
1:A:194:ARG:HG2	1:A:216:ASN:HB3	2.01	0.42
1:B:98:VAL:HG22	1:B:156:TYR:HE2	1.84	0.42
1:B:19:LEU:HA	1:B:22:ALA:HB3	2.02	0.42
1:A:54:PHE:CE1	1:A:79:VAL:HA	2.55	0.42
1:A:154:PHE:O	1:A:155:GLU:C	2.58	0.41
1:B:207:PRO:O	1:B:208:GLU:C	2.58	0.41
1:B:76:LYS:HE2	1:B:342:LYS:O	2.20	0.41
1:A:182:VAL:CG1	1:A:183:THR:HG23	2.49	0.41
1:A:172:LEU:HD12	1:A:180:ILE:HD13	2.03	0.41
1:B:201:THR:HG23	1:B:202:PRO:HD2	2.03	0.41
1:A:200:GLY:HA3	1:A:205:LEU:HD21	2.02	0.41
1:A:62:HIS:NE2	1:A:64:GLU:HB3	2.36	0.41
1:B:98:VAL:CG2	1:B:156:TYR:HE2	2.33	0.41
1:B:144:ARG:CD	1:B:300:THR:HG23	2.48	0.41
1:A:184:ASP:HB3	2:A:401:ADP:O3A	2.21	0.41
1:B:131:HIS:NE2	1:B:314:PHE:CE1	2.88	0.41
1:A:207:PRO:HD3	1:A:222:TRP:CE2	2.56	0.41
1:A:171:ASN:CB	1:A:183:THR:CG2	2.88	0.40
1:A:261:PHE:HB3	1:A:266:LYS:HG3	2.04	0.40
1:A:150:ILE:HD11	1:A:172:LEU:HD13	2.03	0.40
1:B:140:GLU:OE1	1:B:262:SER:CB	2.69	0.40
1:B:158:HIS:ND1	1:B:217:LYS:HD2	2.37	0.40
1:A:113:ASN:OD1	1:A:341:GLU:HA	2.21	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	337/350 (96%)	298 (88%)	39 (12%)	0	100	100
1	B	337/350 (96%)	302 (90%)	35 (10%)	0	100	100
All	All	674/700 (96%)	600 (89%)	74 (11%)	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	231/303 (76%)	204 (88%)	27 (12%)	8	38
1	B	234/303 (77%)	210 (90%)	24 (10%)	10	47
All	All	465/606 (77%)	414 (89%)	51 (11%)	9	43

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	PHE
1	A	74	LEU
1	A	85	ILE
1	A	88	THR
1	A	96	GLN
1	A	98	VAL
1	A	104	VAL

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Mol	Chain	Res	Type
1	A	109	SER
1	A	127	GLU
1	A	137	ARG
1	A	159	SER
1	A	172	LEU
1	A	175	ASP
1	A	176	GLN
1	A	184	ASP
1	A	192	LYS
1	A	195	THR
1	A	201	THR
1	A	203	GLU
1	A	212	SER
1	A	224	LEU
1	A	235	TYR
1	A	265	LEU
1	A	281	PHE
1	A	336	ARG
1	A	340	ASN
1	A	345	LYS
1	B	28	LYS
1	B	37	THR
1	B	41	ASP
1	B	74	LEU
1	B	76	LYS
1	B	106	LEU
1	B	123	VAL
1	B	128	MET
1	B	137	ARG
1	B	153	THR
1	B	175	ASP
1	B	176	GLN
1	B	181	GLN
1	B	190	ARG
1	B	194	ARG
1	B	195	THR
1	B	203	GLU
1	B	217	LYS
1	B	224	LEU
1	B	235	TYR
1	B	267	ASP
1	B	277	LEU

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Mol	Chain	Res	Type
1	B	281	PHE
1	B	340	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	307	GLN
1	B	181	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

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
1	TPO	A	197	1	10,10,11	5.71	2 (20%)	12,14,16	1.22	1 (8%)
1	SEP	A	338	1	9,9,10	6.23	2 (22%)	10,12,14	1.33	1 (10%)
1	TPO	B	197	1	10,10,11	5.46	2 (20%)	12,14,16	1.34	1 (8%)
1	SEP	B	338	1	9,9,10	6.28	3 (33%)	10,12,14	1.37	1 (10%)

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
1	TPO	A	197	1	-	0/9/11/13	0/0/0/0
1	SEP	A	338	1	-	0/6/8/10	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TPO	B	197	1	-	0/9/11/13	0/0/0/0
1	SEP	B	338	1	-	0/6/8/10	0/0/0/0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	338	SEP	O-C	18.22	1.23	1.11
1	A	338	SEP	O-C	18.14	1.23	1.11
1	A	197	TPO	O-C	17.80	1.23	1.11
1	B	197	TPO	O-C	16.95	1.23	1.11
1	B	338	SEP	P-O1P	3.13	1.61	1.51
1	A	338	SEP	P-O1P	3.09	1.61	1.51
1	B	197	TPO	OG1-CB	-2.90	1.41	1.45
1	A	197	TPO	OG1-CB	-2.58	1.42	1.45
1	B	338	SEP	CA-C	2.29	1.54	1.49

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	338	SEP	P-OG-CB	-3.64	109.32	118.63
1	A	338	SEP	P-OG-CB	-3.50	109.69	118.63
1	A	197	TPO	O3P-P-O2P	2.24	115.78	107.38
1	B	197	TPO	O3P-P-O2P	2.17	115.51	107.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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	$\# Z  > 2$
2	ADP	A	401	3	29,29,29	1.08	3 (10%)	45,45,45	2.14	7 (15%)
2	ADP	B	401	3	29,29,29	1.13	2 (6%)	45,45,45	2.03	7 (15%)

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	A	401	3	-	0/16/32/32	0/3/3/3
2	ADP	B	401	3	-	0/16/32/32	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	ADP	C5-C4	3.25	1.47	1.40
2	A	401	ADP	C5-C4	3.20	1.47	1.40
2	B	401	ADP	C4-N9	-2.36	1.34	1.37
2	A	401	ADP	C4-N9	-2.31	1.34	1.37
2	A	401	ADP	O4'-C1'	2.13	1.43	1.41

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	ADP	C5-C4-N3	-7.86	118.32	125.98
2	B	401	ADP	C5-C4-N3	-7.60	118.57	125.98
2	A	401	ADP	N3-C2-N1	-6.26	123.38	128.89
2	B	401	ADP	N3-C2-N1	-6.17	123.47	128.89
2	B	401	ADP	N3-C4-N9	5.83	135.39	125.39
2	A	401	ADP	N3-C4-N9	5.80	135.34	125.39
2	A	401	ADP	C3'-C2'-C1'	3.57	106.51	100.92
2	A	401	ADP	PA-O3A-PB	-3.28	122.85	131.93
2	A	401	ADP	C4-C5-N7	-2.88	106.63	109.41
2	A	401	ADP	C2-N3-C4	2.80	121.33	113.27
2	B	401	ADP	C4-C5-N7	-2.77	106.73	109.41
2	B	401	ADP	C2-N3-C4	2.69	121.01	113.27
2	B	401	ADP	C3'-C2'-C1'	2.50	104.84	100.92
2	B	401	ADP	PA-O3A-PB	-2.48	125.05	131.93

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

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	341/350 (97%)	-0.17	1 (0%) 91 76	82, 101, 129, 146	0
1	B	341/350 (97%)	-0.13	0 100 100	78, 96, 128, 147	0
All	All	682/700 (97%)	-0.15	1 (0%) 93 84	78, 99, 129, 147	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	327	PHE	2.2

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	TPO	B	197	11/12	0.14	-0.90	90,92,94,94	0
1	SEP	B	338	10/11	0.07	-1.81	134,138,146,147	0
1	TPO	A	197	11/12	0.11	-1.99	91,93,94,94	0
1	SEP	A	338	10/11	0.09	-2.32	130,134,141,143	0

### 6.3 Carbohydrates ⓘ

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	MG	B	402	1/1	0.54	4.36	78,78,78,78	0
3	MG	A	402	1/1	0.39	1.40	62,62,62,62	0
2	ADP	A	401	27/27	0.32	1.33	144,156,189,198	0
2	ADP	B	401	27/27	0.25	0.05	115,122,133,139	0

## 6.5 Other polymers

There are no such residues in this entry.