



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 16, 2016 – 07:40 AM EDT

PDB ID : 5JZJ
Title : Crystal structure of DCLK1-KD in complex with AMPPN
Authors : Patel, O.; Lucet, I.
Deposited on : 2016-05-17
Resolution : 1.71 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027939

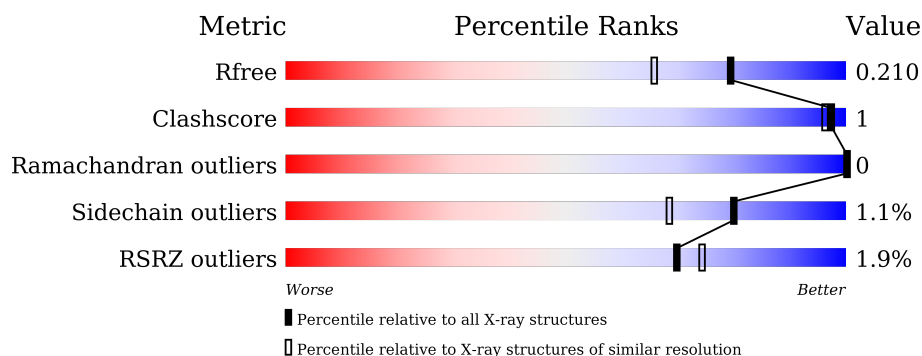
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.71 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}	91344	3998 (1.74-1.70)
Clashscore	102246	4425 (1.74-1.70)
Ramachandran outliers	100387	4360 (1.74-1.70)
Sidechain outliers	100360	4360 (1.74-1.70)
RSRZ outliers	91569	4010 (1.74-1.70)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	297	<div> <div>2%</div> <div> <div></div> <div>88%</div> <div>•</div> <div>9%</div> </div> </div>
1	B	297	<div> <div>2%</div> <div> <div></div> <div>85%</div> <div>•</div> <div>12%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4683 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 Serine/threonine-protein kinase DCLK1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	3	0
			2111	1356	346	398	11			
1	B	261	Total	C	N	O	S	0	4	0
			2018	1302	327	379	10			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	353	MET	-	initiating methionine	UNP O15075
A	354	HIS	-	expression tag	UNP O15075
A	355	HIS	-	expression tag	UNP O15075
A	356	HIS	-	expression tag	UNP O15075
A	357	HIS	-	expression tag	UNP O15075
A	358	HIS	-	expression tag	UNP O15075
A	359	HIS	-	expression tag	UNP O15075
A	360	HIS	-	expression tag	UNP O15075
A	361	HIS	-	expression tag	UNP O15075
A	362	GLY	-	expression tag	UNP O15075
A	363	SER	-	expression tag	UNP O15075
A	364	GLU	-	expression tag	UNP O15075
A	365	ASN	-	expression tag	UNP O15075
A	366	LEU	-	expression tag	UNP O15075
A	367	TYR	-	expression tag	UNP O15075
A	368	PHE	-	expression tag	UNP O15075
A	369	GLN	-	expression tag	UNP O15075
A	370	GLY	-	expression tag	UNP O15075
A	371	SER	-	expression tag	UNP O15075
B	353	MET	-	initiating methionine	UNP O15075
B	354	HIS	-	expression tag	UNP O15075
B	355	HIS	-	expression tag	UNP O15075
B	356	HIS	-	expression tag	UNP O15075
B	357	HIS	-	expression tag	UNP O15075
B	358	HIS	-	expression tag	UNP O15075

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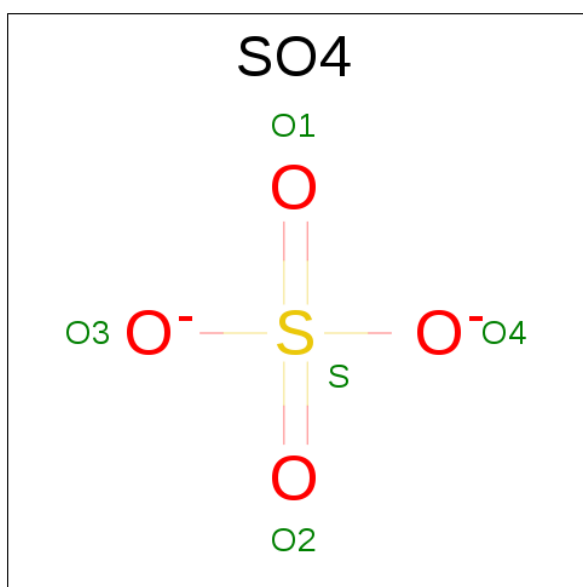
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Chain	Residue	Modelled	Actual	Comment	Reference
B	359	HIS	-	expression tag	UNP O15075
B	360	HIS	-	expression tag	UNP O15075
B	361	HIS	-	expression tag	UNP O15075
B	362	GLY	-	expression tag	UNP O15075
B	363	SER	-	expression tag	UNP O15075
B	364	GLU	-	expression tag	UNP O15075
B	365	ASN	-	expression tag	UNP O15075
B	366	LEU	-	expression tag	UNP O15075
B	367	TYR	-	expression tag	UNP O15075
B	368	PHE	-	expression tag	UNP O15075
B	369	GLN	-	expression tag	UNP O15075
B	370	GLY	-	expression tag	UNP O15075
B	371	SER	-	expression tag	UNP O15075

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

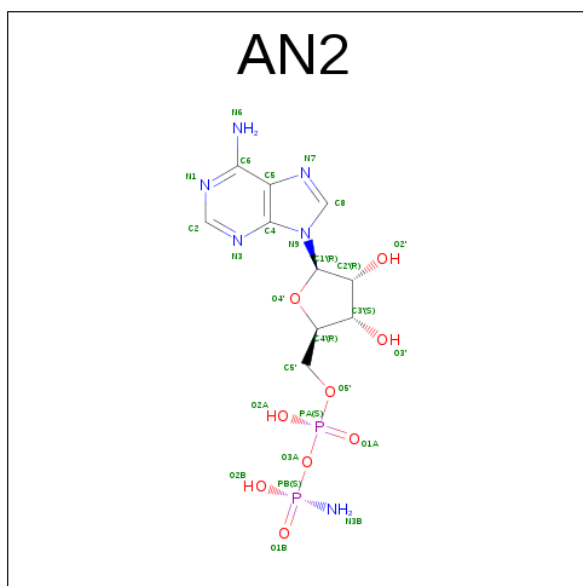
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		

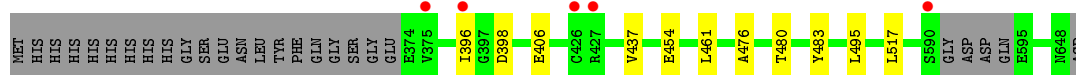
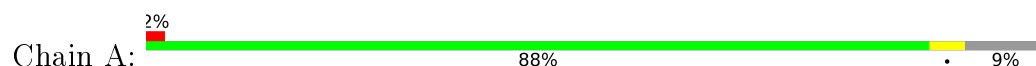
- Molecule 4 is AMP PHOSPHORAMIDATE (three-letter code: AN2) (formula: $C_{10}H_{16}N_6O_9P_2$).



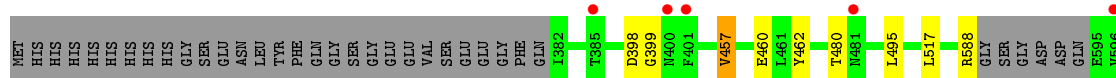
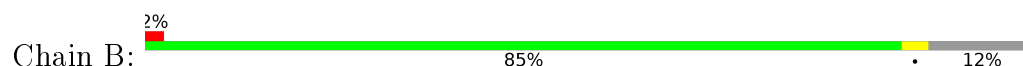
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Serine/threonine-protein kinase DCLK1



- Molecule 1: Serine/threonine-protein kinase DCLK1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.39Å 81.27Å 59.61Å 90.00° 92.10° 90.00°	Depositor
Resolution (Å)	56.35 – 1.71 56.35 – 1.71	Depositor EDS
% Data completeness (in resolution range)	99.6 (56.35-1.71) 99.9 (56.35-1.71)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.88 (at 1.71Å)	Xtriage
Refinement program	BUSTER-TNT	Depositor
R, R_{free}	0.177 , 0.208 0.178 , 0.210	Depositor DCC
R_{free} test set	2932 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	18.1	Xtriage
Anisotropy	0.392	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4683	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MG, SO4, AN2

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	0.50	0/2162	0.59	0/2939
1	B	0.50	0/2071	0.59	0/2824
All	All	0.50	0/4233	0.59	0/5763

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	2111	0	2100	5	0
1	B	2018	0	1998	5	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	5	0	0	0	0
4	A	27	0	14	0	0
4	B	27	0	14	1	0
5	A	271	0	0	0	0
5	B	222	0	0	1	0
All	All	4683	0	4126	10	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 1.

The worst 5 of 10 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:396:ILE:HG21	1:A:406:GLU:HG3	1.88	0.54
1:A:476:ALA:O	1:A:480:THR:HG22	2.09	0.51
1:B:457:VAL:HG12	1:B:460:GLU:H	1.79	0.47
1:B:495:LEU:HD13	1:B:517:LEU:HD21	1.97	0.46
1:A:495:LEU:HD13	1:A:517:LEU:HD21	1.98	0.45

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	270/297 (91%)	265 (98%)	5 (2%)	0	100	100
1	B	261/297 (88%)	257 (98%)	4 (2%)	0	100	100
All	All	531/594 (89%)	522 (98%)	9 (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	230/259 (89%)	227 (99%)	3 (1%)	76	61
1	B	218/259 (84%)	214 (98%)	4 (2%)	66	48
All	All	448/518 (86%)	441 (98%)	7 (2%)	80	53

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	398[A]	ASP
1	B	480	THR
1	B	398[B]	ASP
1	A	398[B]	ASP
1	B	457	VAL

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 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	SO4	A	702	-	4,4,4	0.34	0	6,6,6	0.14	0
4	AN2	A	703	2	24,29,29	1.80	6 (25%)	22,45,45	2.76	2 (9%)
4	AN2	B	902	2	24,29,29	1.98	8 (33%)	22,45,45	3.33	3 (13%)

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	SO4	A	702	-	-	0/0/0/0	0/0/0/0
4	AN2	A	703	2	-	0/9/32/32	0/3/3/3
4	AN2	B	902	2	-	0/9/32/32	0/3/3/3

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	902	AN2	PB-O2B	-3.55	1.47	1.56
4	A	703	AN2	PB-O2B	-3.20	1.48	1.56
4	B	902	AN2	C5-C4	-2.51	1.34	1.40
4	A	703	AN2	C5-C4	-2.23	1.35	1.40
4	B	902	AN2	O3'-C3'	2.07	1.47	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	902	AN2	N3-C2-N1	-14.35	117.60	128.87
4	A	703	AN2	N3-C2-N1	-12.04	119.41	128.87
4	B	902	AN2	O2A-PA-O3A	2.28	115.03	105.27
4	A	703	AN2	O2B-PB-O1B	3.43	118.56	110.16
4	B	902	AN2	O2B-PB-O1B	4.20	120.45	110.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	902	AN2	1	0

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, 95th 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 ²)	Q<0.9
1	A	271/297 (91%)	-0.12	5 (1%) 71 76	10, 21, 44, 63	0
1	B	261/297 (87%)	-0.10	5 (1%) 70 74	12, 23, 43, 52	0
All	All	532/594 (89%)	-0.11	10 (1%) 70 74	10, 23, 44, 63	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	375	VAL	3.4
1	B	401	PHE	3.1
1	B	481	ASN	2.7
1	B	400	ASN	2.7
1	A	427	ARG	2.6

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. 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, 95th 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	LLDF	B-factors(\AA^2)	Q<0.9
4	AN2	B	902	27/27	0.98	0.07	-0.75	14,16,21,27	0
4	AN2	A	703	27/27	0.98	0.08	-0.91	15,17,20,26	0
3	SO4	A	702	5/5	1.00	0.06	-3.76	13,13,15,16	0
2	MG	B	901	1/1	1.00	0.05	-	16,16,16,16	0
2	MG	A	701	1/1	0.99	0.03	-	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.