

Full wwPDB X-ray Structure Validation Report (i)

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PDB ID : 3OSM

> Title : Structure of the Kinase Associated Domain-1 (KA1) from Kcc4p

Authors Moravcevic, K.; Lemmon, M.A.

2010-09-09 Deposited on

1.70 Å(reported) Resolution

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 A user guide is available at

https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity 4.02b-467

> 1.8.5 (274361), CSD as541be (2020) Mogul

Xtriage (Phenix) 1.13 EDS 2.11

Percentile statistics 20191225.v01 (using entries in the PDB archive December 25th 2019)

> Refmac 5.8.0158

7.0.044 (Gargrove) CCP4

Ideal geometry (proteins) Engh & Huber (2001) Ideal geometry (DNA, RNA) Parkinson et al. (1996)

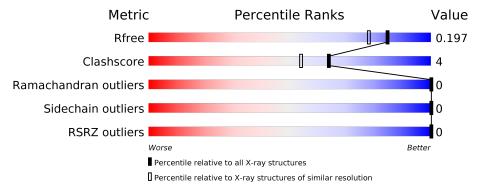
Validation Pipeline (wwPDB-VP) 2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
R_{free}	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 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	A	128	86%	7%	7%

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	\mathbf{Type}	Chain	m Res	Chirality	$\mathbf{Geometry}$	Clashes	Electron density
2	SRT	A	1039	X	-	-	-



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 1056 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.

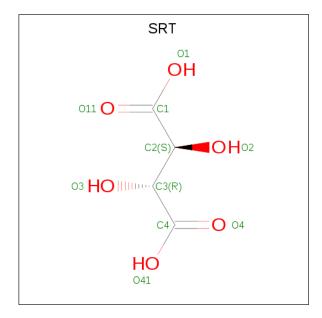
• Molecule 1 is a protein called serine/threonine-protein kinase KCC4.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	119	Total	С	N	О	S	0	0	0
1	Λ	119	920	602	155	161	2	0	U	

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	910	MET	_	expression tag	UNP P25389
A	911	HIS	-	expression tag	UNP P25389
A	912	HIS	_	expression tag	UNP P25389
A	913	HIS	-	expression tag	UNP P25389
A	914	HIS	-	expression tag	UNP P25389
A	915	HIS	-	expression tag	UNP P25389
A	916	HIS	_	expression tag	UNP P25389

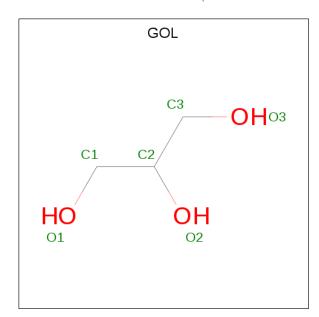
• Molecule 2 is S,R MESO-TARTARIC ACID (three-letter code: SRT) (formula: C₄H₆O₆).





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
2	A	1	Total 10	C 4	O 6	0	0

• Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total C 6 3	O 3	0	0

• Molecule 4 is water.

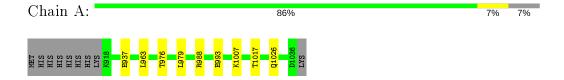
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	120	Total O 120 120	0	0



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: serine/threonine-protein kinase KCC4





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants	26.24Å 33.23Å 38.15Å	Depositor
a, b, c, α , β , γ	104.15° 90.70° 107.00°	Depositor
Resolution (Å)	36.84 - 1.70	Depositor
Resolution (A)	30.69 - 1.70	EDS
% Data completeness	96.5 (36.84-1.70)	Depositor
(in resolution range)	96.3 (30.69-1.70)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$< I/\sigma(I) > 1$	2.30 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
D D.	0.187 , 0.197	Depositor
R, R_{free}	0.187 , 0.197	DCC
R_{free} test set	614 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.305	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.37, 46.3	EDS
L-test for twinning ²	$ < L > = 0.50, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	1056	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 14.56% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $< L^2>$ 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: GOL, SRT

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).

Mal	N/IOI ('hain		Bond lengths		Bond angles	
MIOI	Chain	RMSZ	# Z >5	RMSZ	# Z > 5	
1	A	0.54	0/934	0.60	0/1259	

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)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
1	A	920	0	970	8	0
2	A	10	0	4	2	0
3	A	6	0	8	0	0
4	A	120	0	0	2	0
All	All	1056	0	982	8	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 4.

All (8) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.



Atom-1	Atom-2	$egin{array}{l} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{array}$	$egin{array}{c} ext{Clash} \ ext{overlap } (ext{Å}) \end{array}$
1:A:976:THR:HG22	1:A:993:GLU:HG2	1.77	0.65
1:A:988:ARG:HH22	2:A:1039:SRT:H3	1.69	0.58
1:A:976:THR:HG23	4:A:83:HOH:O	2.06	0.54
1:A:979:LEU:HD13	1:A:1017:THR:HG22	1.93	0.51
1:A:988:ARG:NH2	2:A:1039:SRT:H3	2.33	0.43
1:A:937:GLU:HG2	1:A:1007:LYS:HD3	2.01	0.42
1:A:1026:GLN:NE2	4:A:97:HOH:O	2.46	0.42
1:A:963:LEU:HD23	1:A:979:LEU:HD21	2.02	0.42

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	Perce	ntiles
1	A	117/128 (91%)	116 (99%)	1 (1%)	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	99/117 (85%)	99 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no



such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

2 ligands 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	T 0	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Dag	Link	\mathbf{B}	Bond lengths			ond ang	gles
	Chain	nes	nes	Res	res Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2							
3	GOL	A	1038	-	5,5,5	0.28	0	5,5,5	0.43	0								
2	SRT	A	1039	-	3,9,9	0.30	0	6,12,12	2.13	2 (33%)								

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	GOL	A	1038	-	1	2/4/4/4	-
2	SRT	A	1039	-	2/2/4/4	4/4/12/12	-

There are no bond length outliers.



All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
2	A	1039	SRT	O2-C2-C3	-3.91	95.85	108.90
2	A	1039	SRT	C1-C2-C3	2.35	118.16	113.11

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	1039	SRT	C2
2	A	1039	SRT	С3

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1038	GOL	O1-C1-C2-O2
3	A	1038	GOL	O1-C1-C2-C3
2	A	1039	SRT	O2-C2-C3-O3
2	A	1039	SRT	C1-C2-C3-O3
2	A	1039	SRT	C1-C2-C3-C4
2	A	1039	SRT	O2-C2-C3-C4

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1039	SRT	2	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, 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	$\langle { m RSRZ} \rangle$	$\#\text{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	119/128 (92%)	-0.19	0 100 100	12, 16, 24, 28	0

There are no RSRZ outliers to report.

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^{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	${f B-factors}({f \AA}^2)$	Q<0.9
3	GOL	A	1038	6/6	0.86	0.19	24,28,31,31	0
2	SRT	A	1039	10/10	0.88	0.15	21,25,27,28	0

6.5 Other polymers (i)

There are no such residues in this entry.

