



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2018 – 05:54 PM EST

PDB ID : 6CCF  
Title : Crystal Structure of the Human CAMKK1A in complex with Hesperadin  
Authors : Santiago, A.S.; Counago, R.M.; dos Reis, C.V.; Ramos, P.Z.; Silva, P.N.B.; Drewry, D.; Elkins, J.M.; Massirer, K.B.; Arruda, P.; Edwards, A.M.; Structural Genomics Consortium (SGC)  
Deposited on : 2018-02-07  
Resolution : 2.10 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

<http://wwpdb.org/validation/2016/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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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-20030736

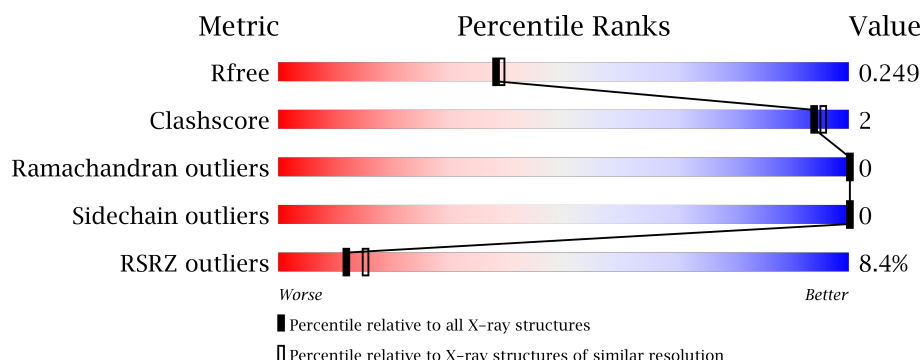
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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))
$R_{free}$	100719	4243 (2.10-2.10)
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (2.10-2.10)

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. 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	290	<div> <div>9%</div> <div>87%</div> <div>10%</div> </div>
1	B	290	<div> <div>6%</div> <div>86%</div> <div>10%</div> </div>



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.

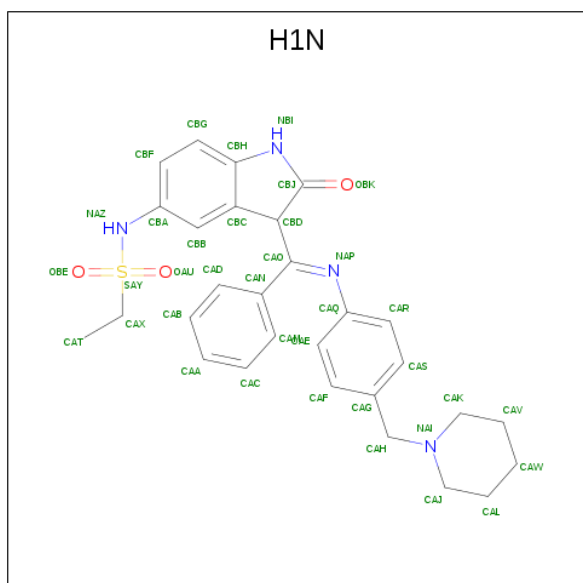
- Molecule 1 is a protein called Calcium/calmodulin-dependent protein kinase kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	260	Total 2007	C 1290	N 329	O 380	S 8	0	0	0
1	B	262	Total 2016	C 1300	N 334	O 374	S 8	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

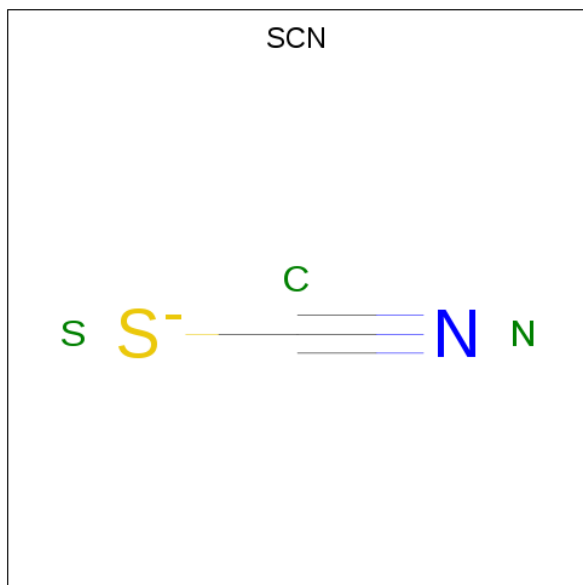
Chain	Residue	Modelled	Actual	Comment	Reference
A	122	SER	-	expression tag	UNP Q8N5S9
A	123	MET	-	expression tag	UNP Q8N5S9
B	122	SER	-	expression tag	UNP Q8N5S9
B	123	MET	-	expression tag	UNP Q8N5S9

- Molecule 2 is N-[2-OXO-3-((E)-PHENYL{[4-(PIPERIDIN-1-YLMETHYL)PHENYL]IMINO}METHYL)-2,6-DIHYDRO-1H-INDOL-5-YL]ETHANESULFONAMIDE (three-letter code: H1N) (formula:  $C_{29}H_{32}N_4O_3S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			37	29	4	3	1		
2	B	1	Total	C	N	O	S	0	0
			37	29	4	3	1		

- Molecule 3 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	S	0	0
			3	1	1	1		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



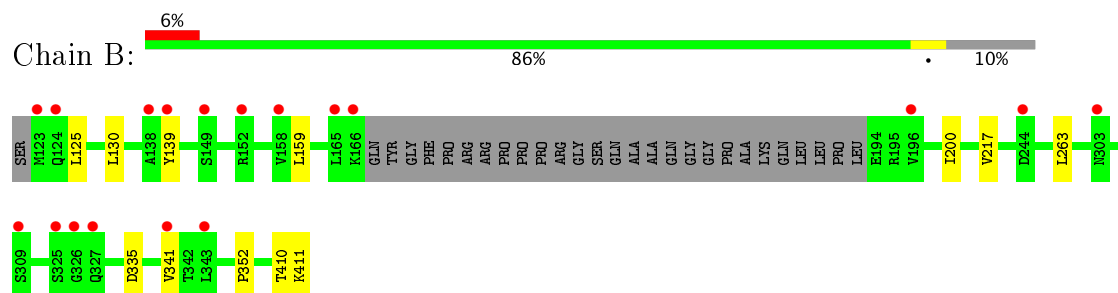
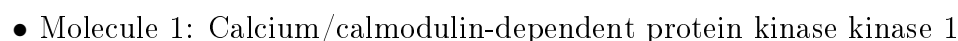
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	126	Total	O	0	0
			126	126		
5	B	145	Total	O	0	0
			145	145		



- Molecule 1: Calcium/calmodulin-dependent protein kinase kinase 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	49.22Å 83.17Å 80.12Å 90.00° 97.27° 90.00°	Depositor
Resolution (Å)	19.87 – 2.10 19.87 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.4 (19.87-2.10) 99.5 (19.87-2.10)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.10 (at 2.09Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, $R_{free}$	0.210 , 0.244 0.217 , 0.249	Depositor DCC
$R_{free}$ test set	1885 reflections (5.33%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.5	Xtriage
Anisotropy	0.185	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 39.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4375	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.40% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SCN, EDO, H1N

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	0.43	0/2049	0.64	0/2782
1	B	0.44	0/2058	0.64	0/2795
All	All	0.44	0/4107	0.64	0/5577

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	2007	0	1950	5	0
1	B	2016	0	1970	6	0
2	A	37	0	31	1	0
2	B	37	0	31	1	0
3	A	3	0	0	0	0
4	B	4	0	6	0	0
5	A	126	0	0	0	0
5	B	145	0	0	0	0
All	All	4375	0	3988	13	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 (13) 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:B:263:LEU:HD21	1:B:335:ASP:HB3	1.88	0.55
1:A:296:VAL:HG13	1:A:311:ALA:HB1	1.93	0.49
1:B:200:ILE:CG2	1:B:217:VAL:HG21	2.43	0.49
2:A:501:H1N:HAR	2:A:501:H1N:CAN	2.45	0.46
1:B:139:TYR:HB2	1:B:159:LEU:HD22	1.98	0.45
2:B:501:H1N:HAR	2:B:501:H1N:CAN	2.46	0.45
1:B:410:THR:O	1:B:411:LYS:CB	2.66	0.43
1:A:270:LYS:HA	1:A:300:PHE:CZ	2.54	0.43
1:B:341:VAL:HG13	1:B:352:PRO:HD2	2.00	0.42
1:A:329:PHE:CE1	1:A:334:LEU:HD11	2.54	0.42
1:B:125:LEU:CD1	1:B:130:LEU:HD11	2.51	0.41
1:A:139:TYR:HB2	1:A:159:LEU:HD22	2.02	0.40
1:A:296:VAL:HG13	1:A:311:ALA:CB	2.51	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	256/290 (88%)	251 (98%)	5 (2%)	0	100	100
1	B	258/290 (89%)	249 (96%)	9 (4%)	0	100	100
All	All	514/580 (89%)	500 (97%)	14 (3%)	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	212/253 (84%)	212 (100%)	0	100	100
1	B	211/253 (83%)	211 (100%)	0	100	100
All	All	423/506 (84%)	423 (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 ⓘ

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 ⓘ

4 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	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	H1N	A	501	-	38,41,41	1.35	1 (2%)	44,58,58	1.34	3 (6%)
3	SCN	A	502	-	1,2,2	0.05	0	0,1,1	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	H1N	B	501	-	38,41,41	1.19	2 (5%)	44,58,58	1.27	3 (6%)
4	EDO	B	502	-	3,3,3	0.49	0	2,2,2	0.34	0

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	H1N	A	501	-	-	0/24/57/57	0/5/5/5
3	SCN	A	502	-	-	0/0/0/0	0/0/0/0
2	H1N	B	501	-	-	0/24/57/57	0/5/5/5
4	EDO	B	502	-	-	0/1/1/1	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	H1N	CBH-NBI	2.16	1.39	1.35
2	B	501	H1N	CAO-NAP	6.51	1.38	1.28
2	A	501	H1N	CAO-NAP	7.59	1.40	1.28

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	H1N	CAX-SAY-NAZ	-2.54	97.82	105.26
2	B	501	H1N	CAX-SAY-NAZ	-2.20	98.82	105.26
2	B	501	H1N	CBH-NBI-CBJ	2.77	111.53	106.81
2	A	501	H1N	CBH-NBI-CBJ	2.90	111.75	106.81
2	B	501	H1N	CAQ-NAP-CAO	5.93	133.52	121.98
2	A	501	H1N	CAQ-NAP-CAO	6.73	135.10	121.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	H1N	1	0
2	B	501	H1N	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 ⓘ

### 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	260/290 (89%)	0.59	26 (10%) 8 10	20, 35, 62, 79	0
1	B	262/290 (90%)	0.51	18 (6%) 18 22	20, 32, 59, 99	0
All	All	522/580 (90%)	0.55	44 (8%) 12 15	20, 33, 60, 99	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	326	GLY	12.3
1	A	326	GLY	8.2
1	A	139	TYR	4.9
1	B	138	ALA	4.8
1	A	149	SER	4.5
1	A	164	LEU	4.2
1	B	149	SER	3.9
1	B	123	MET	3.7
1	A	196	VAL	3.6
1	A	138	ALA	3.6
1	A	137	GLY	3.5
1	A	325	SER	3.3
1	B	325	SER	3.3
1	B	327	GLN	3.2
1	B	343	LEU	3.2
1	A	195	ARG	3.1
1	B	165	LEU	3.1
1	B	139	TYR	3.0
1	A	148	GLU	3.0
1	A	343	LEU	2.9
1	A	197	TYR	2.8
1	B	196	VAL	2.7
1	B	244	ASP	2.7
1	A	310	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	198	GLN	2.6
1	A	340	GLY	2.6
1	A	162	LYS	2.5
1	A	194	GLU	2.4
1	B	124	GLN	2.4
1	B	166	LYS	2.4
1	B	309	SER	2.4
1	A	336	VAL	2.4
1	A	308	SER	2.4
1	B	341	VAL	2.4
1	A	140	GLY	2.3
1	A	201	ALA	2.3
1	A	199	GLU	2.3
1	A	341	VAL	2.3
1	B	303	ASN	2.2
1	B	158	VAL	2.2
1	A	276	ILE	2.1
1	A	132	SER	2.1
1	A	327	GLN	2.0
1	B	152	ARG	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. 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	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	EDO	B	502	4/4	0.96	0.12	0.85	40,40,41,41	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	H1N	A	501	37/37	0.86	0.17	0.29	33,38,56,57	0
2	H1N	B	501	37/37	0.92	0.15	0.26	26,29,62,65	0
3	SCN	A	502	3/3	0.87	0.12	-	58,58,59,68	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.