Full wwPDB X-ray Structure Validation Report (i)

Apr 27, 2024 – 02:39 pm BST

PDB ID	:	5J03
Title	:	Crystal Structure of a chimeric Kv7.2 - Kv7.3 proximal C-terminal Domain in Complex with Calmodulin
Authors	:	Strulovich, R.; Hirsch, J.A.
Deposited on	:	2016-03-26
Resolution	:	2.00 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*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

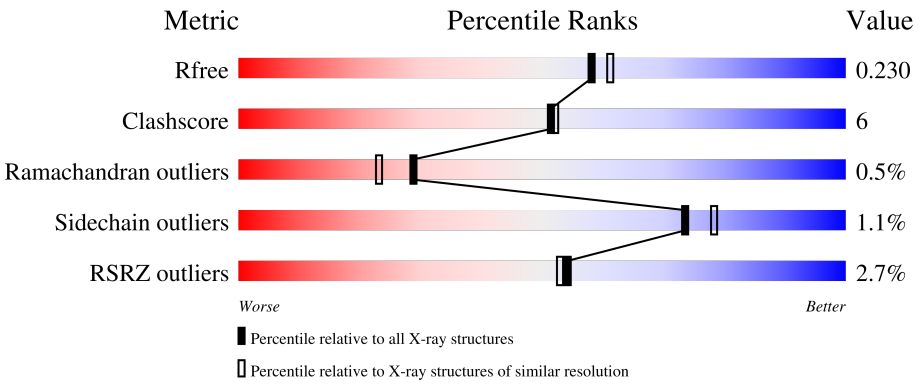
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 of 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	110	5%; 60%; 5%; 34%
2	B	149	%; 89%; 9%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3505 atoms, of which 1634 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 Potassium voltage-gated channel subfamily KQT member 3,Potassium voltage-gated channel subfamily KQT member 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	73	1156	377	571	109	97	2	0	0	0

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	317	MET	-	initiating methionine	UNP O43525
A	318	GLY	-	expression tag	UNP O43525
A	319	SER	-	expression tag	UNP O43525
A	320	HIS	-	expression tag	UNP O43525
A	321	HIS	-	expression tag	UNP O43525
A	322	HIS	-	expression tag	UNP O43525
A	323	HIS	-	expression tag	UNP O43525
A	324	HIS	-	expression tag	UNP O43525
A	325	HIS	-	expression tag	UNP O43525
A	326	HIS	-	expression tag	UNP O43525
A	327	HIS	-	expression tag	UNP O43525
A	328	GLY	-	expression tag	UNP O43525
A	329	SER	-	expression tag	UNP O43525
A	330	ASP	-	expression tag	UNP O43525
A	331	TYR	-	expression tag	UNP O43525
A	332	ASP	-	expression tag	UNP O43525
A	333	ILE	-	expression tag	UNP O43525
A	334	PRO	-	expression tag	UNP O43525
A	335	THR	-	expression tag	UNP O43525
A	336	THR	-	expression tag	UNP O43525
A	337	GLU	-	expression tag	UNP O43525
A	338	ASN	-	expression tag	UNP O43525
A	339	LEU	-	expression tag	UNP O43525
A	340	TYR	-	expression tag	UNP O43525
A	341	PHE	-	expression tag	UNP O43525
A	342	GLN	-	expression tag	UNP O43525

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Chain	Residue	Modelled	Actual	Comment	Reference
A	344	SER	LEU	conflict	UNP O43525

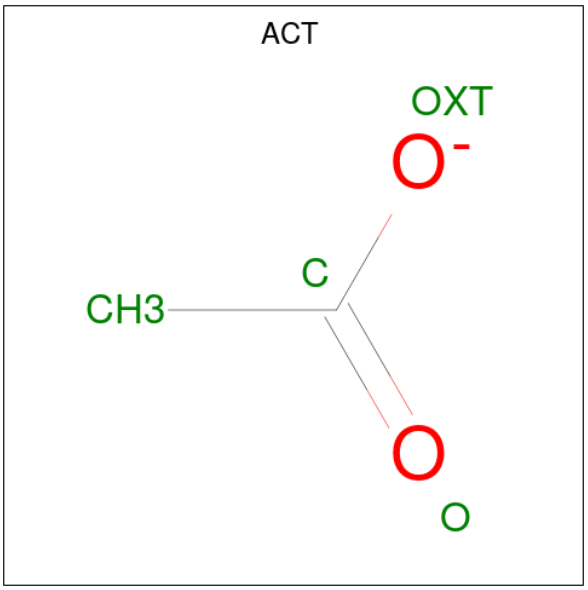
- Molecule 2 is a protein called Calmodulin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	146	2204	702	1060	187	246	9	0	3	0

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
3	B	5	5	5	0	0

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
4	B	1	7	2	3	2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	40	40	40	0	0

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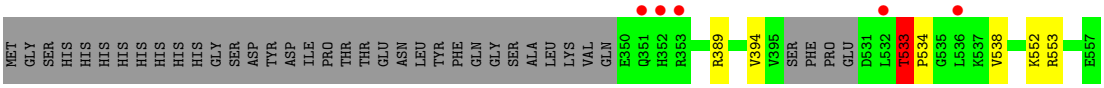
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	B	93	93	93	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Potassium voltage-gated channel subfamily KQT member 3,Potassium voltage-gated channel subfamily KQT member 2

Chain A:



• Molecule 2: Calmodulin

Chain B:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	47.66Å 47.66Å 182.22Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.26 – 2.00 40.26 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.26-2.00) 99.9 (40.26-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.01 (at 2.00Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.179 , 0.228 0.180 , 0.230	Depositor DCC
R_{free} test set	919 reflections (5.39%)	wwPDB-VP
Wilson B-factor (Å^2)	28.7	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.41 , 52.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.043 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3505	wwPDB-VP
Average B, all atoms (Å^2)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical 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.

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: CA, ACT

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	0.61	0/598	0.73	1/806 (0.1%)
2	B	0.65	0/1163	0.72	1/1564 (0.1%)
All	All	0.64	0/1761	0.72	2/2370 (0.1%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	533	THR	C-N-CD	-5.54	108.42	120.60
2	B	106	LEU	CB-CG-CD2	-5.09	102.35	111.00

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	585	571	569	10	0
2	B	1144	1060	1058	14	0
3	B	5	0	0	0	0
4	B	4	3	3	0	0
5	A	40	0	0	2	0
5	B	93	0	0	3	0
All	All	1871	1634	1630	19	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 6.

All (19) 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)
2:B:23[B]:ASP:OD1	5:B:301:HOH:O	1.79	1.00
2:B:78:LYS:H	2:B:78:LYS:HD2	1.46	0.80
1:A:534:PRO:O	5:A:601:HOH:O	1.99	0.79
1:A:553:ARG:NH1	2:B:55:GLU:OE2	2.16	0.77
2:B:138:ASN:OD1	2:B:140:GLU:HG2	1.97	0.64
1:A:533:THR:CB	1:A:534:PRO:HA	2.32	0.60
2:B:78:LYS:H	2:B:78:LYS:CD	2.14	0.57
2:B:23[B]:ASP:OD1	2:B:23[B]:ASP:N	2.38	0.57
1:A:389:ARG:NH2	5:A:602:HOH:O	2.28	0.53
2:B:3:ASP:N	5:B:305:HOH:O	2.42	0.52
1:A:533:THR:CB	1:A:534:PRO:CA	2.91	0.49
1:A:538:VAL:HG11	2:B:16:ALA:HB2	1.96	0.48
1:A:394:VAL:HG12	2:B:41:GLY:HA3	1.94	0.48
2:B:122:VAL:O	2:B:126:ILE:HG12	2.15	0.47
1:A:552:LYS:NZ	2:B:80:THR:O	2.48	0.47
2:B:78:LYS:HD2	2:B:78:LYS:N	2.22	0.46
1:A:538:VAL:CG1	2:B:16:ALA:HB2	2.45	0.46
2:B:140:GLU:HG2	5:B:312:HOH:O	2.17	0.42
1:A:394:VAL:HG12	1:A:394:VAL:O	2.19	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	Percentiles
1	A	69/110 (63%)	63 (91%)	5 (7%)	1 (1%)	11 5
2	B	147/149 (99%)	146 (99%)	1 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	216/259 (83%)	209 (97%)	6 (3%)	1 (0%)	29 23

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	533	THR

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	56/98 (57%)	56 (100%)	0	100 100
2	B	121/127 (95%)	119 (98%)	2 (2%)	60 65
All	All	177/225 (79%)	175 (99%)	2 (1%)	73 78

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	131	ILE
2	B	145	MET

Sometimes 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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 5 are monoatomic - leaving 1 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	$\#\|Z\| > 2$
4	ACT	B	206	3	3,3,3	0.52	0	3,3,3	1.48	1 (33%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	206	ACT	OXT-C-CH3	2.01	123.48	115.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	73/110 (66%)	0.18	5 (6%)	17	16	24, 38, 97, 118	0
2	B	146/149 (97%)	-0.32	1 (0%)	87	87	24, 36, 58, 104	0
All	All	219/259 (84%)	-0.15	6 (2%)	54	53	24, 36, 81, 118	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	532	LEU	3.9
2	B	3	ASP	3.5
1	A	351	GLN	3.3
1	A	353	ARG	2.6
1	A	536	LEU	2.4
1	A	352	HIS	2.1

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 monosaccharides 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, 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	B-factors(Å^2)	Q<0.9
3	CA	B	201	1/1	0.97	0.12	35,35,35,35	0
3	CA	B	204	1/1	0.98	0.10	24,24,24,24	1
4	ACT	B	206	4/4	0.98	0.12	27,30,36,36	0
3	CA	B	202	1/1	0.99	0.10	34,34,34,34	0
3	CA	B	205	1/1	0.99	0.05	30,30,30,30	1
3	CA	B	203	1/1	0.99	0.09	29,29,29,29	0

6.5 Other polymers (i)

There are no such residues in this entry.