



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 28, 2023 – 09:57 AM EDT

PDB ID : 3KXD
Title : Crystal structure of the mthk rck in complex with cadmium
Authors : Dvir, H.; Choe, S.
Deposited on : 2009-12-02
Resolution : 2.20 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtrriage (Phenix) : 1.13
EDS : 2.35
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) : 2.35

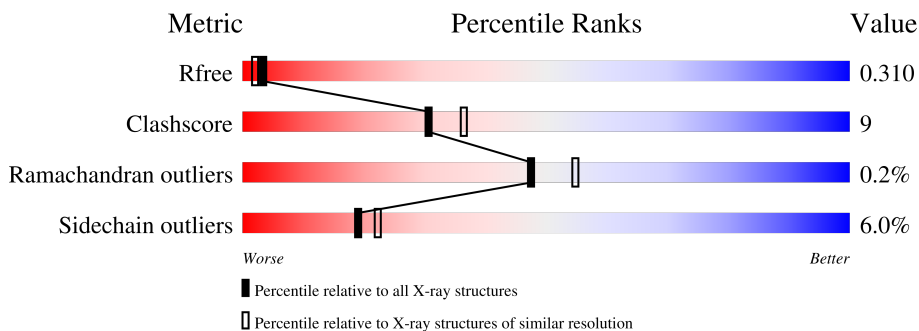
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	224	75% (green), 23% (yellow), 2% (orange), 0% (red), 0% (grey)
1	B	224	73% (green), 23% (yellow), 4% (orange), 0% (red), 0% (grey)

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3494 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 Calcium-gated potassium channel mthK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	224	Total	C	N	O	S	0	0	0
			1706	1067	296	336	7			
1	B	219	Total	C	N	O	S	0	0	0
			1672	1045	296	324	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	337	LEU	-	expression tag	UNP O27564
A	338	VAL	-	expression tag	UNP O27564
A	339	PRO	-	expression tag	UNP O27564
B	337	LEU	-	expression tag	UNP O27564
B	338	VAL	-	expression tag	UNP O27564
B	339	PRO	-	expression tag	UNP O27564

- Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

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

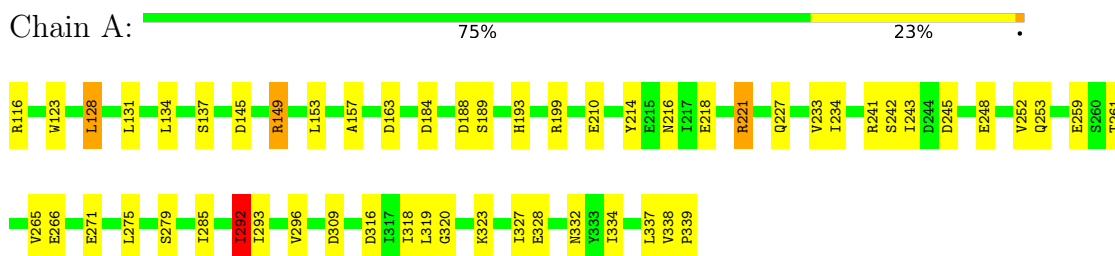
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	67	Total	O	0	0
			67	67		
3	B	45	Total	O	0	0
			45	45		

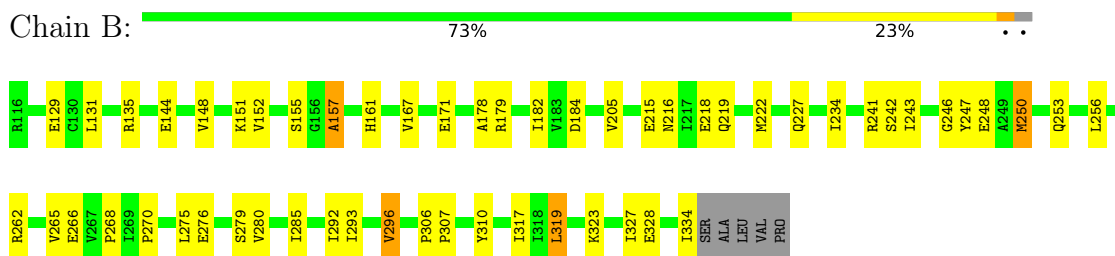
3 Residue-property plots

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. 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. 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: Calcium-gated potassium channel mthK



- Molecule 1: Calcium-gated potassium channel mthK



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.02Å 38.29Å 96.63Å 90.00° 95.10° 90.00°	Depositor
Resolution (Å)	57.83 – 2.20 57.79 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.3 (57.83-2.20) 99.3 (57.79-2.20)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.04 (at 2.20Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.184 , 0.254 0.253 , 0.310	Depositor DCC
R_{free} test set	1108 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	35.7	Xtrriage
Anisotropy	0.043	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3494	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.04% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CD

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	1.06	1/1728 (0.1%)	1.07	10/2340 (0.4%)
1	B	0.94	2/1693 (0.1%)	0.85	1/2290 (0.0%)
All	All	1.00	3/3421 (0.1%)	0.97	11/4630 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	247	TYR	CD1-CE1	-5.84	1.30	1.39
1	A	199	ARG	CB-CG	5.38	1.67	1.52
1	B	247	TYR	CD2-CE2	-5.30	1.31	1.39

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	ARG	NE-CZ-NH1	8.15	124.37	120.30
1	A	149	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	A	292	ILE	CB-CA-C	-6.33	98.95	111.60
1	B	184	ASP	CB-CG-OD2	-6.31	112.62	118.30
1	A	199	ARG	NE-CZ-NH2	-6.26	117.17	120.30
1	A	128	LEU	CB-CG-CD1	5.81	120.88	111.00
1	A	221	ARG	NE-CZ-NH2	-5.73	117.44	120.30
1	A	163	ASP	CB-CG-OD1	5.72	123.45	118.30
1	A	184	ASP	CB-CG-OD2	-5.22	113.61	118.30
1	A	316	ASP	CB-CG-OD1	5.18	122.97	118.30
1	A	145	ASP	CB-CG-OD1	5.12	122.91	118.30

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	1706	0	1691	32	0
1	B	1672	0	1665	38	0
2	A	3	0	0	0	0
2	B	1	0	0	0	0
3	A	67	0	0	3	0
3	B	45	0	0	0	0
All	All	3494	0	3356	62	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 9.

All (62) 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:265:VAL:HG21	1:B:327:ILE:HD12	1.35	1.05
1:A:265:VAL:HG21	1:A:327:ILE:HD12	1.49	0.90
1:B:285:ILE:HG21	1:B:293:ILE:HD11	1.56	0.86
1:B:265:VAL:HG21	1:B:327:ILE:CD1	2.09	0.81
1:B:265:VAL:CG2	1:B:327:ILE:HD12	2.11	0.80
1:B:262:ARG:HH11	1:B:262:ARG:HG3	1.46	0.78
1:A:116:ARG:HA	3:A:60:HOH:O	1.89	0.73
1:A:265:VAL:CG2	1:A:327:ILE:HD12	2.23	0.68
1:B:167:VAL:O	1:B:171:GLU:HG3	1.93	0.67
1:A:252:VAL:HG21	1:B:256:LEU:HD11	1.79	0.65
1:A:221:ARG:HG2	3:A:44:HOH:O	1.95	0.64
1:B:144:GLU:O	1:B:161:HIS:HE1	1.81	0.63
1:A:285:ILE:HG21	1:A:293:ILE:HD11	1.81	0.63
1:B:218:GLU:O	1:B:222:MET:HG3	1.99	0.63
1:B:265:VAL:CG2	1:B:327:ILE:CD1	2.75	0.62
1:B:131:LEU:HD11	1:B:157:ALA:HB2	1.81	0.62
1:A:275:LEU:HD12	1:A:334:ILE:HG22	1.81	0.62
1:A:296:VAL:HG22	1:A:318:ILE:HD13	1.83	0.61
1:B:262:ARG:HG3	1:B:262:ARG:NH1	2.14	0.59
1:B:151:LYS:O	1:B:155:SER:HB2	2.03	0.58
1:B:148:VAL:O	1:B:152:VAL:HG23	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:243:ILE:HG22	1:B:243:ILE:O	2.08	0.54
1:B:279:SER:HA	1:B:310:TYR:O	2.09	0.53
1:B:144:GLU:O	1:B:161:HIS:CE1	2.61	0.53
1:B:241:ARG:NH1	1:B:248:GLU:OE1	2.41	0.52
1:A:189:SER:O	1:A:193:HIS:CD2	2.62	0.52
1:A:188:ASP:CG	1:A:216:ASN:ND2	2.64	0.51
1:B:148:VAL:O	1:B:148:VAL:HG23	2.11	0.51
1:A:242:SER:HB3	1:B:227:GLN:OE1	2.11	0.51
1:B:266:GLU:HG2	1:B:319:LEU:HD23	1.93	0.50
1:B:262:ARG:NH2	1:B:323:LYS:HE2	2.27	0.50
1:A:233:VAL:HA	1:B:129:GLU:HG2	1.94	0.49
1:B:266:GLU:HG2	1:B:319:LEU:CD2	2.44	0.47
1:B:275:LEU:HD12	1:B:334:ILE:HD11	1.96	0.47
1:A:218:GLU:OE2	1:A:221:ARG:HD2	2.15	0.46
1:A:243:ILE:HD11	1:B:182:ILE:HD11	1.98	0.46
1:A:227:GLN:OE1	1:B:242:SER:HB3	2.16	0.46
1:A:241:ARG:NH1	1:A:248:GLU:OE1	2.49	0.45
1:B:178:ALA:O	1:B:205:VAL:HG23	2.17	0.44
1:A:292:ILE:O	1:A:320:GLY:HA2	2.18	0.44
1:A:134:LEU:O	1:A:137:SER:HB3	2.17	0.44
1:A:189:SER:O	1:A:193:HIS:HD2	2.00	0.44
1:B:280:VAL:HG23	1:B:310:TYR:HB3	1.99	0.43
1:B:268:PRO:O	1:B:270:PRO:HD3	2.18	0.43
1:A:323:LYS:O	1:A:327:ILE:HG12	2.19	0.43
1:A:234:ILE:HG21	1:B:234:ILE:HG21	2.00	0.42
1:A:332:ASN:N	1:A:332:ASN:HD22	2.17	0.42
1:A:319:LEU:N	1:A:319:LEU:HD12	2.35	0.42
1:B:296:VAL:HA	1:B:317:ILE:O	2.20	0.41
1:A:218:GLU:HG3	3:A:48:HOH:O	2.20	0.41
1:A:245:ASP:C	1:A:245:ASP:OD1	2.58	0.41
1:B:262:ARG:NH1	1:B:262:ARG:CG	2.82	0.41
1:A:266:GLU:HG3	1:B:248:GLU:OE2	2.21	0.41
1:A:131:LEU:HD11	1:A:157:ALA:HB2	2.01	0.41
1:A:243:ILE:CG2	1:B:179:ARG:HD3	2.51	0.41
1:A:271:GLU:HA	1:A:271:GLU:OE1	2.20	0.41
1:A:338:VAL:O	1:A:339:PRO:O	2.39	0.41
1:B:306:PRO:HA	1:B:307:PRO:HD3	1.94	0.41
1:A:123:TRP:CH2	1:A:131:LEU:HD12	2.56	0.41
1:A:188:ASP:CG	1:A:216:ASN:HD21	2.23	0.40
1:B:246:GLY:O	1:B:250:MET:HB2	2.21	0.40
1:B:216:ASN:O	1:B:219:GLN:HB2	2.22	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	222/224 (99%)	217 (98%)	5 (2%)	0	100	100
1	B	217/224 (97%)	205 (94%)	11 (5%)	1 (0%)	29	31
All	All	439/448 (98%)	422 (96%)	16 (4%)	1 (0%)	47	55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	157	ALA

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	184/192 (96%)	171 (93%)	13 (7%)	14	16
1	B	180/192 (94%)	171 (95%)	9 (5%)	24	30
All	All	364/384 (95%)	342 (94%)	22 (6%)	19	22

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	128	LEU
1	A	149	ARG

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Mol	Chain	Res	Type
1	A	153	LEU
1	A	210	GLU
1	A	214	TYR
1	A	253	GLN
1	A	259	GLU
1	A	261	THR
1	A	279	SER
1	A	292	ILE
1	A	309	ASP
1	A	328	GLU
1	A	337	LEU
1	B	135	ARG
1	B	215	GLU
1	B	250	MET
1	B	253	GLN
1	B	276	GLU
1	B	292	ILE
1	B	296	VAL
1	B	319	LEU
1	B	328	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	193	HIS
1	A	332	ASN
1	B	158	ASN
1	B	161	HIS
1	B	253	GLN
1	B	332	ASN

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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.