



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 02:48 AM UTC

PDB ID : 8T6Q / pdb_00008t6q
EMDB ID : EMD-41077
Title : Cryo-EM structure of dodecameric CaMKII beta holoenzyme T287A T306A T307A
Authors : Chien, C.-T.; Chiu, W.; Khan, S.
Deposited on : 2023-06-16
Resolution : 3.50 Å(reported)

This is a Full wwPDB EM 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/EMValidationReportHelp>
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:

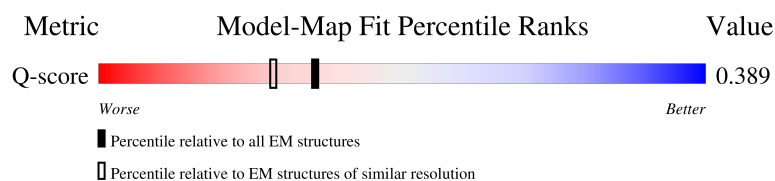
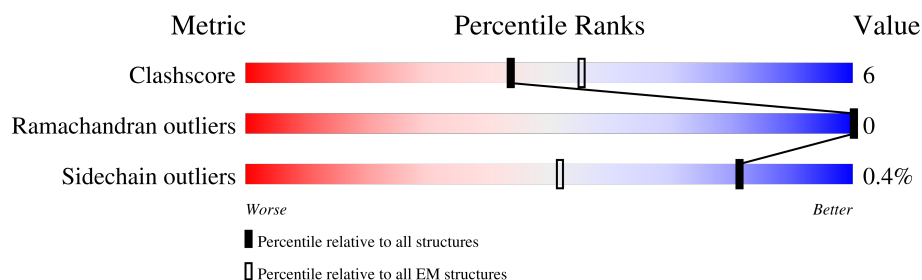
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

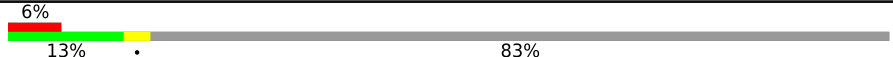
The reported resolution of this entry is 3.50 Å.

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)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13950 (3.00 - 4.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	815	
1	B	815	
1	C	815	
1	D	815	

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Mol	Chain	Length	Quality of chain
1	F	815	13% . 83%
1	G	815	15% . 83%
1	H	815	14% . 83%
1	I	815	15% . 83%
1	J	815	15% . 83%
1	K	815	15% . 83%
1	L	815	13% . 83%
1	N	815	15% . 83%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 25482 atoms, of which 12534 are hydrogens and 0 are deuteriums.

In the tables below, 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 Venus-tagged CaMKII beta holoenzyme mutant.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	135	Total 2124	C 681	H 1045	N 193	O 201	S 4	0	0
1	B	135	Total 2124	C 681	H 1045	N 193	O 201	S 4	0	0
1	C	135	Total 2124	C 681	H 1045	N 193	O 201	S 4	0	0
1	D	135	Total 2124	C 681	H 1045	N 193	O 201	S 4	0	0
1	F	135	Total 2124	C 681	H 1045	N 193	O 201	S 4	0	0
1	G	135	Total 2124	C 681	H 1045	N 193	O 201	S 4	0	0
1	H	135	Total 2123	C 681	H 1044	N 193	O 201	S 4	0	0
1	I	135	Total 2123	C 681	H 1044	N 193	O 201	S 4	0	0
1	J	135	Total 2123	C 681	H 1044	N 193	O 201	S 4	0	0
1	K	135	Total 2123	C 681	H 1044	N 193	O 201	S 4	0	0
1	L	135	Total 2123	C 681	H 1044	N 193	O 201	S 4	0	0
1	N	135	Total 2123	C 681	H 1044	N 193	O 201	S 4	0	0

There are 600 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-272	MET	-	initiating methionine	UNP P42212
A	-271	GLY	-	expression tag	UNP P42212
A	-270	SER	-	expression tag	UNP P42212
A	-269	SER	-	expression tag	UNP P42212
A	-268	HIS	-	expression tag	UNP P42212
A	-267	HIS	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-266	HIS	-	expression tag	UNP P42212
A	-265	HIS	-	expression tag	UNP P42212
A	-264	HIS	-	expression tag	UNP P42212
A	-263	HIS	-	expression tag	UNP P42212
A	-262	SER	-	expression tag	UNP P42212
A	-261	SER	-	expression tag	UNP P42212
A	-260	GLY	-	expression tag	UNP P42212
A	-259	LEU	-	expression tag	UNP P42212
A	-258	VAL	-	expression tag	UNP P42212
A	-257	PRO	-	expression tag	UNP P42212
A	-256	ARG	-	expression tag	UNP P42212
A	-255	GLY	-	expression tag	UNP P42212
A	-254	SER	-	expression tag	UNP P42212
A	-253	HIS	-	expression tag	UNP P42212
A	-252	VAL	-	expression tag	UNP P42212
A	-207	LEU	PHE	conflict	UNP P42212
A	-189	LEU	PHE	conflict	UNP P42212
A	-188	GLY	SER	conflict	UNP P42212
A	-185	LEU	VAL	conflict	UNP P42212
A	-181	ALA	SER	conflict	UNP P42212
A	-100	THR	MET	conflict	UNP P42212
A	-90	ALA	VAL	conflict	UNP P42212
A	-78	GLY	SER	conflict	UNP P42212
A	-50	TYR	THR	conflict	UNP P42212
A	-47	LYS	ALA	conflict	UNP P42212
A	-22	LEU	HIS	conflict	UNP P42212
A	-14	SER	-	linker	UNP P42212
A	-13	GLY	-	linker	UNP P42212
A	-12	LEU	-	linker	UNP P42212
A	-11	ARG	-	linker	UNP P42212
A	-10	SER	-	linker	UNP P42212
A	-9	ARG	-	linker	UNP P42212
A	-8	ALA	-	linker	UNP P42212
A	-7	GLN	-	linker	UNP P42212
A	-6	ALA	-	linker	UNP P42212
A	-5	SER	-	linker	UNP P42212
A	-4	ASN	-	linker	UNP P42212
A	-3	SER	-	linker	UNP P42212
A	-2	ALA	-	linker	UNP P42212
A	-1	VAL	-	linker	UNP P42212
A	0	ASP	-	linker	UNP P42212
A	287	ALA	THR	engineered mutation	UNP P08413

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Chain	Residue	Modelled	Actual	Comment	Reference
A	306	ALA	THR	engineered mutation	UNP P08413
A	307	ALA	THR	engineered mutation	UNP P08413
B	-272	MET	-	initiating methionine	UNP P42212
B	-271	GLY	-	expression tag	UNP P42212
B	-270	SER	-	expression tag	UNP P42212
B	-269	SER	-	expression tag	UNP P42212
B	-268	HIS	-	expression tag	UNP P42212
B	-267	HIS	-	expression tag	UNP P42212
B	-266	HIS	-	expression tag	UNP P42212
B	-265	HIS	-	expression tag	UNP P42212
B	-264	HIS	-	expression tag	UNP P42212
B	-263	HIS	-	expression tag	UNP P42212
B	-262	SER	-	expression tag	UNP P42212
B	-261	SER	-	expression tag	UNP P42212
B	-260	GLY	-	expression tag	UNP P42212
B	-259	LEU	-	expression tag	UNP P42212
B	-258	VAL	-	expression tag	UNP P42212
B	-257	PRO	-	expression tag	UNP P42212
B	-256	ARG	-	expression tag	UNP P42212
B	-255	GLY	-	expression tag	UNP P42212
B	-254	SER	-	expression tag	UNP P42212
B	-253	HIS	-	expression tag	UNP P42212
B	-252	VAL	-	expression tag	UNP P42212
B	-207	LEU	PHE	conflict	UNP P42212
B	-189	LEU	PHE	conflict	UNP P42212
B	-188	GLY	SER	conflict	UNP P42212
B	-185	LEU	VAL	conflict	UNP P42212
B	-181	ALA	SER	conflict	UNP P42212
B	-100	THR	MET	conflict	UNP P42212
B	-90	ALA	VAL	conflict	UNP P42212
B	-78	GLY	SER	conflict	UNP P42212
B	-50	TYR	THR	conflict	UNP P42212
B	-47	LYS	ALA	conflict	UNP P42212
B	-22	LEU	HIS	conflict	UNP P42212
B	-14	SER	-	linker	UNP P42212
B	-13	GLY	-	linker	UNP P42212
B	-12	LEU	-	linker	UNP P42212
B	-11	ARG	-	linker	UNP P42212
B	-10	SER	-	linker	UNP P42212
B	-9	ARG	-	linker	UNP P42212
B	-8	ALA	-	linker	UNP P42212
B	-7	GLN	-	linker	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	ALA	-	linker	UNP P42212
B	-5	SER	-	linker	UNP P42212
B	-4	ASN	-	linker	UNP P42212
B	-3	SER	-	linker	UNP P42212
B	-2	ALA	-	linker	UNP P42212
B	-1	VAL	-	linker	UNP P42212
B	0	ASP	-	linker	UNP P42212
B	287	ALA	THR	engineered mutation	UNP P08413
B	306	ALA	THR	engineered mutation	UNP P08413
B	307	ALA	THR	engineered mutation	UNP P08413
C	-272	MET	-	initiating methionine	UNP P42212
C	-271	GLY	-	expression tag	UNP P42212
C	-270	SER	-	expression tag	UNP P42212
C	-269	SER	-	expression tag	UNP P42212
C	-268	HIS	-	expression tag	UNP P42212
C	-267	HIS	-	expression tag	UNP P42212
C	-266	HIS	-	expression tag	UNP P42212
C	-265	HIS	-	expression tag	UNP P42212
C	-264	HIS	-	expression tag	UNP P42212
C	-263	HIS	-	expression tag	UNP P42212
C	-262	SER	-	expression tag	UNP P42212
C	-261	SER	-	expression tag	UNP P42212
C	-260	GLY	-	expression tag	UNP P42212
C	-259	LEU	-	expression tag	UNP P42212
C	-258	VAL	-	expression tag	UNP P42212
C	-257	PRO	-	expression tag	UNP P42212
C	-256	ARG	-	expression tag	UNP P42212
C	-255	GLY	-	expression tag	UNP P42212
C	-254	SER	-	expression tag	UNP P42212
C	-253	HIS	-	expression tag	UNP P42212
C	-252	VAL	-	expression tag	UNP P42212
C	-207	LEU	PHE	conflict	UNP P42212
C	-189	LEU	PHE	conflict	UNP P42212
C	-188	GLY	SER	conflict	UNP P42212
C	-185	LEU	VAL	conflict	UNP P42212
C	-181	ALA	SER	conflict	UNP P42212
C	-100	THR	MET	conflict	UNP P42212
C	-90	ALA	VAL	conflict	UNP P42212
C	-78	GLY	SER	conflict	UNP P42212
C	-50	TYR	THR	conflict	UNP P42212
C	-47	LYS	ALA	conflict	UNP P42212
C	-22	LEU	HIS	conflict	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-14	SER	-	linker	UNP P42212
C	-13	GLY	-	linker	UNP P42212
C	-12	LEU	-	linker	UNP P42212
C	-11	ARG	-	linker	UNP P42212
C	-10	SER	-	linker	UNP P42212
C	-9	ARG	-	linker	UNP P42212
C	-8	ALA	-	linker	UNP P42212
C	-7	GLN	-	linker	UNP P42212
C	-6	ALA	-	linker	UNP P42212
C	-5	SER	-	linker	UNP P42212
C	-4	ASN	-	linker	UNP P42212
C	-3	SER	-	linker	UNP P42212
C	-2	ALA	-	linker	UNP P42212
C	-1	VAL	-	linker	UNP P42212
C	0	ASP	-	linker	UNP P42212
C	287	ALA	THR	engineered mutation	UNP P08413
C	306	ALA	THR	engineered mutation	UNP P08413
C	307	ALA	THR	engineered mutation	UNP P08413
D	-272	MET	-	initiating methionine	UNP P42212
D	-271	GLY	-	expression tag	UNP P42212
D	-270	SER	-	expression tag	UNP P42212
D	-269	SER	-	expression tag	UNP P42212
D	-268	HIS	-	expression tag	UNP P42212
D	-267	HIS	-	expression tag	UNP P42212
D	-266	HIS	-	expression tag	UNP P42212
D	-265	HIS	-	expression tag	UNP P42212
D	-264	HIS	-	expression tag	UNP P42212
D	-263	HIS	-	expression tag	UNP P42212
D	-262	SER	-	expression tag	UNP P42212
D	-261	SER	-	expression tag	UNP P42212
D	-260	GLY	-	expression tag	UNP P42212
D	-259	LEU	-	expression tag	UNP P42212
D	-258	VAL	-	expression tag	UNP P42212
D	-257	PRO	-	expression tag	UNP P42212
D	-256	ARG	-	expression tag	UNP P42212
D	-255	GLY	-	expression tag	UNP P42212
D	-254	SER	-	expression tag	UNP P42212
D	-253	HIS	-	expression tag	UNP P42212
D	-252	VAL	-	expression tag	UNP P42212
D	-207	LEU	PHE	conflict	UNP P42212
D	-189	LEU	PHE	conflict	UNP P42212
D	-188	GLY	SER	conflict	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-185	LEU	VAL	conflict	UNP P42212
D	-181	ALA	SER	conflict	UNP P42212
D	-100	THR	MET	conflict	UNP P42212
D	-90	ALA	VAL	conflict	UNP P42212
D	-78	GLY	SER	conflict	UNP P42212
D	-50	TYR	THR	conflict	UNP P42212
D	-47	LYS	ALA	conflict	UNP P42212
D	-22	LEU	HIS	conflict	UNP P42212
D	-14	SER	-	linker	UNP P42212
D	-13	GLY	-	linker	UNP P42212
D	-12	LEU	-	linker	UNP P42212
D	-11	ARG	-	linker	UNP P42212
D	-10	SER	-	linker	UNP P42212
D	-9	ARG	-	linker	UNP P42212
D	-8	ALA	-	linker	UNP P42212
D	-7	GLN	-	linker	UNP P42212
D	-6	ALA	-	linker	UNP P42212
D	-5	SER	-	linker	UNP P42212
D	-4	ASN	-	linker	UNP P42212
D	-3	SER	-	linker	UNP P42212
D	-2	ALA	-	linker	UNP P42212
D	-1	VAL	-	linker	UNP P42212
D	0	ASP	-	linker	UNP P42212
D	287	ALA	THR	engineered mutation	UNP P08413
D	306	ALA	THR	engineered mutation	UNP P08413
D	307	ALA	THR	engineered mutation	UNP P08413
F	-272	MET	-	initiating methionine	UNP P42212
F	-271	GLY	-	expression tag	UNP P42212
F	-270	SER	-	expression tag	UNP P42212
F	-269	SER	-	expression tag	UNP P42212
F	-268	HIS	-	expression tag	UNP P42212
F	-267	HIS	-	expression tag	UNP P42212
F	-266	HIS	-	expression tag	UNP P42212
F	-265	HIS	-	expression tag	UNP P42212
F	-264	HIS	-	expression tag	UNP P42212
F	-263	HIS	-	expression tag	UNP P42212
F	-262	SER	-	expression tag	UNP P42212
F	-261	SER	-	expression tag	UNP P42212
F	-260	GLY	-	expression tag	UNP P42212
F	-259	LEU	-	expression tag	UNP P42212
F	-258	VAL	-	expression tag	UNP P42212
F	-257	PRO	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-256	ARG	-	expression tag	UNP P42212
F	-255	GLY	-	expression tag	UNP P42212
F	-254	SER	-	expression tag	UNP P42212
F	-253	HIS	-	expression tag	UNP P42212
F	-252	VAL	-	expression tag	UNP P42212
F	-207	LEU	PHE	conflict	UNP P42212
F	-189	LEU	PHE	conflict	UNP P42212
F	-188	GLY	SER	conflict	UNP P42212
F	-185	LEU	VAL	conflict	UNP P42212
F	-181	ALA	SER	conflict	UNP P42212
F	-100	THR	MET	conflict	UNP P42212
F	-90	ALA	VAL	conflict	UNP P42212
F	-78	GLY	SER	conflict	UNP P42212
F	-50	TYR	THR	conflict	UNP P42212
F	-47	LYS	ALA	conflict	UNP P42212
F	-22	LEU	HIS	conflict	UNP P42212
F	-14	SER	-	linker	UNP P42212
F	-13	GLY	-	linker	UNP P42212
F	-12	LEU	-	linker	UNP P42212
F	-11	ARG	-	linker	UNP P42212
F	-10	SER	-	linker	UNP P42212
F	-9	ARG	-	linker	UNP P42212
F	-8	ALA	-	linker	UNP P42212
F	-7	GLN	-	linker	UNP P42212
F	-6	ALA	-	linker	UNP P42212
F	-5	SER	-	linker	UNP P42212
F	-4	ASN	-	linker	UNP P42212
F	-3	SER	-	linker	UNP P42212
F	-2	ALA	-	linker	UNP P42212
F	-1	VAL	-	linker	UNP P42212
F	0	ASP	-	linker	UNP P42212
F	287	ALA	THR	engineered mutation	UNP P08413
F	306	ALA	THR	engineered mutation	UNP P08413
F	307	ALA	THR	engineered mutation	UNP P08413
G	-272	MET	-	initiating methionine	UNP P42212
G	-271	GLY	-	expression tag	UNP P42212
G	-270	SER	-	expression tag	UNP P42212
G	-269	SER	-	expression tag	UNP P42212
G	-268	HIS	-	expression tag	UNP P42212
G	-267	HIS	-	expression tag	UNP P42212
G	-266	HIS	-	expression tag	UNP P42212
G	-265	HIS	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-264	HIS	-	expression tag	UNP P42212
G	-263	HIS	-	expression tag	UNP P42212
G	-262	SER	-	expression tag	UNP P42212
G	-261	SER	-	expression tag	UNP P42212
G	-260	GLY	-	expression tag	UNP P42212
G	-259	LEU	-	expression tag	UNP P42212
G	-258	VAL	-	expression tag	UNP P42212
G	-257	PRO	-	expression tag	UNP P42212
G	-256	ARG	-	expression tag	UNP P42212
G	-255	GLY	-	expression tag	UNP P42212
G	-254	SER	-	expression tag	UNP P42212
G	-253	HIS	-	expression tag	UNP P42212
G	-252	VAL	-	expression tag	UNP P42212
G	-207	LEU	PHE	conflict	UNP P42212
G	-189	LEU	PHE	conflict	UNP P42212
G	-188	GLY	SER	conflict	UNP P42212
G	-185	LEU	VAL	conflict	UNP P42212
G	-181	ALA	SER	conflict	UNP P42212
G	-100	THR	MET	conflict	UNP P42212
G	-90	ALA	VAL	conflict	UNP P42212
G	-78	GLY	SER	conflict	UNP P42212
G	-50	TYR	THR	conflict	UNP P42212
G	-47	LYS	ALA	conflict	UNP P42212
G	-22	LEU	HIS	conflict	UNP P42212
G	-14	SER	-	linker	UNP P42212
G	-13	GLY	-	linker	UNP P42212
G	-12	LEU	-	linker	UNP P42212
G	-11	ARG	-	linker	UNP P42212
G	-10	SER	-	linker	UNP P42212
G	-9	ARG	-	linker	UNP P42212
G	-8	ALA	-	linker	UNP P42212
G	-7	GLN	-	linker	UNP P42212
G	-6	ALA	-	linker	UNP P42212
G	-5	SER	-	linker	UNP P42212
G	-4	ASN	-	linker	UNP P42212
G	-3	SER	-	linker	UNP P42212
G	-2	ALA	-	linker	UNP P42212
G	-1	VAL	-	linker	UNP P42212
G	0	ASP	-	linker	UNP P42212
G	287	ALA	THR	engineered mutation	UNP P08413
G	306	ALA	THR	engineered mutation	UNP P08413
G	307	ALA	THR	engineered mutation	UNP P08413

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Chain	Residue	Modelled	Actual	Comment	Reference
H	-272	MET	-	initiating methionine	UNP P42212
H	-271	GLY	-	expression tag	UNP P42212
H	-270	SER	-	expression tag	UNP P42212
H	-269	SER	-	expression tag	UNP P42212
H	-268	HIS	-	expression tag	UNP P42212
H	-267	HIS	-	expression tag	UNP P42212
H	-266	HIS	-	expression tag	UNP P42212
H	-265	HIS	-	expression tag	UNP P42212
H	-264	HIS	-	expression tag	UNP P42212
H	-263	HIS	-	expression tag	UNP P42212
H	-262	SER	-	expression tag	UNP P42212
H	-261	SER	-	expression tag	UNP P42212
H	-260	GLY	-	expression tag	UNP P42212
H	-259	LEU	-	expression tag	UNP P42212
H	-258	VAL	-	expression tag	UNP P42212
H	-257	PRO	-	expression tag	UNP P42212
H	-256	ARG	-	expression tag	UNP P42212
H	-255	GLY	-	expression tag	UNP P42212
H	-254	SER	-	expression tag	UNP P42212
H	-253	HIS	-	expression tag	UNP P42212
H	-252	VAL	-	expression tag	UNP P42212
H	-207	LEU	PHE	conflict	UNP P42212
H	-189	LEU	PHE	conflict	UNP P42212
H	-188	GLY	SER	conflict	UNP P42212
H	-185	LEU	VAL	conflict	UNP P42212
H	-181	ALA	SER	conflict	UNP P42212
H	-100	THR	MET	conflict	UNP P42212
H	-90	ALA	VAL	conflict	UNP P42212
H	-78	GLY	SER	conflict	UNP P42212
H	-50	TYR	THR	conflict	UNP P42212
H	-47	LYS	ALA	conflict	UNP P42212
H	-22	LEU	HIS	conflict	UNP P42212
H	-14	SER	-	linker	UNP P42212
H	-13	GLY	-	linker	UNP P42212
H	-12	LEU	-	linker	UNP P42212
H	-11	ARG	-	linker	UNP P42212
H	-10	SER	-	linker	UNP P42212
H	-9	ARG	-	linker	UNP P42212
H	-8	ALA	-	linker	UNP P42212
H	-7	GLN	-	linker	UNP P42212
H	-6	ALA	-	linker	UNP P42212
H	-5	SER	-	linker	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
H	-4	ASN	-	linker	UNP P42212
H	-3	SER	-	linker	UNP P42212
H	-2	ALA	-	linker	UNP P42212
H	-1	VAL	-	linker	UNP P42212
H	0	ASP	-	linker	UNP P42212
H	287	ALA	THR	engineered mutation	UNP P08413
H	306	ALA	THR	engineered mutation	UNP P08413
H	307	ALA	THR	engineered mutation	UNP P08413
I	-272	MET	-	initiating methionine	UNP P42212
I	-271	GLY	-	expression tag	UNP P42212
I	-270	SER	-	expression tag	UNP P42212
I	-269	SER	-	expression tag	UNP P42212
I	-268	HIS	-	expression tag	UNP P42212
I	-267	HIS	-	expression tag	UNP P42212
I	-266	HIS	-	expression tag	UNP P42212
I	-265	HIS	-	expression tag	UNP P42212
I	-264	HIS	-	expression tag	UNP P42212
I	-263	HIS	-	expression tag	UNP P42212
I	-262	SER	-	expression tag	UNP P42212
I	-261	SER	-	expression tag	UNP P42212
I	-260	GLY	-	expression tag	UNP P42212
I	-259	LEU	-	expression tag	UNP P42212
I	-258	VAL	-	expression tag	UNP P42212
I	-257	PRO	-	expression tag	UNP P42212
I	-256	ARG	-	expression tag	UNP P42212
I	-255	GLY	-	expression tag	UNP P42212
I	-254	SER	-	expression tag	UNP P42212
I	-253	HIS	-	expression tag	UNP P42212
I	-252	VAL	-	expression tag	UNP P42212
I	-207	LEU	PHE	conflict	UNP P42212
I	-189	LEU	PHE	conflict	UNP P42212
I	-188	GLY	SER	conflict	UNP P42212
I	-185	LEU	VAL	conflict	UNP P42212
I	-181	ALA	SER	conflict	UNP P42212
I	-100	THR	MET	conflict	UNP P42212
I	-90	ALA	VAL	conflict	UNP P42212
I	-78	GLY	SER	conflict	UNP P42212
I	-50	TYR	THR	conflict	UNP P42212
I	-47	LYS	ALA	conflict	UNP P42212
I	-22	LEU	HIS	conflict	UNP P42212
I	-14	SER	-	linker	UNP P42212
I	-13	GLY	-	linker	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
I	-12	LEU	-	linker	UNP P42212
I	-11	ARG	-	linker	UNP P42212
I	-10	SER	-	linker	UNP P42212
I	-9	ARG	-	linker	UNP P42212
I	-8	ALA	-	linker	UNP P42212
I	-7	GLN	-	linker	UNP P42212
I	-6	ALA	-	linker	UNP P42212
I	-5	SER	-	linker	UNP P42212
I	-4	ASN	-	linker	UNP P42212
I	-3	SER	-	linker	UNP P42212
I	-2	ALA	-	linker	UNP P42212
I	-1	VAL	-	linker	UNP P42212
I	0	ASP	-	linker	UNP P42212
I	287	ALA	THR	engineered mutation	UNP P08413
I	306	ALA	THR	engineered mutation	UNP P08413
I	307	ALA	THR	engineered mutation	UNP P08413
J	-272	MET	-	initiating methionine	UNP P42212
J	-271	GLY	-	expression tag	UNP P42212
J	-270	SER	-	expression tag	UNP P42212
J	-269	SER	-	expression tag	UNP P42212
J	-268	HIS	-	expression tag	UNP P42212
J	-267	HIS	-	expression tag	UNP P42212
J	-266	HIS	-	expression tag	UNP P42212
J	-265	HIS	-	expression tag	UNP P42212
J	-264	HIS	-	expression tag	UNP P42212
J	-263	HIS	-	expression tag	UNP P42212
J	-262	SER	-	expression tag	UNP P42212
J	-261	SER	-	expression tag	UNP P42212
J	-260	GLY	-	expression tag	UNP P42212
J	-259	LEU	-	expression tag	UNP P42212
J	-258	VAL	-	expression tag	UNP P42212
J	-257	PRO	-	expression tag	UNP P42212
J	-256	ARG	-	expression tag	UNP P42212
J	-255	GLY	-	expression tag	UNP P42212
J	-254	SER	-	expression tag	UNP P42212
J	-253	HIS	-	expression tag	UNP P42212
J	-252	VAL	-	expression tag	UNP P42212
J	-207	LEU	PHE	conflict	UNP P42212
J	-189	LEU	PHE	conflict	UNP P42212
J	-188	GLY	SER	conflict	UNP P42212
J	-185	LEU	VAL	conflict	UNP P42212
J	-181	ALA	SER	conflict	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
J	-100	THR	MET	conflict	UNP P42212
J	-90	ALA	VAL	conflict	UNP P42212
J	-78	GLY	SER	conflict	UNP P42212
J	-50	TYR	THR	conflict	UNP P42212
J	-47	LYS	ALA	conflict	UNP P42212
J	-22	LEU	HIS	conflict	UNP P42212
J	-14	SER	-	linker	UNP P42212
J	-13	GLY	-	linker	UNP P42212
J	-12	LEU	-	linker	UNP P42212
J	-11	ARG	-	linker	UNP P42212
J	-10	SER	-	linker	UNP P42212
J	-9	ARG	-	linker	UNP P42212
J	-8	ALA	-	linker	UNP P42212
J	-7	GLN	-	linker	UNP P42212
J	-6	ALA	-	linker	UNP P42212
J	-5	SER	-	linker	UNP P42212
J	-4	ASN	-	linker	UNP P42212
J	-3	SER	-	linker	UNP P42212
J	-2	ALA	-	linker	UNP P42212
J	-1	VAL	-	linker	UNP P42212
J	0	ASP	-	linker	UNP P42212
J	287	ALA	THR	engineered mutation	UNP P08413
J	306	ALA	THR	engineered mutation	UNP P08413
J	307	ALA	THR	engineered mutation	UNP P08413
K	-272	MET	-	initiating methionine	UNP P42212
K	-271	GLY	-	expression tag	UNP P42212
K	-270	SER	-	expression tag	UNP P42212
K	-269	SER	-	expression tag	UNP P42212
K	-268	HIS	-	expression tag	UNP P42212
K	-267	HIS	-	expression tag	UNP P42212
K	-266	HIS	-	expression tag	UNP P42212
K	-265	HIS	-	expression tag	UNP P42212
K	-264	HIS	-	expression tag	UNP P42212
K	-263	HIS	-	expression tag	UNP P42212
K	-262	SER	-	expression tag	UNP P42212
K	-261	SER	-	expression tag	UNP P42212
K	-260	GLY	-	expression tag	UNP P42212
K	-259	LEU	-	expression tag	UNP P42212
K	-258	VAL	-	expression tag	UNP P42212
K	-257	PRO	-	expression tag	UNP P42212
K	-256	ARG	-	expression tag	UNP P42212
K	-255	GLY	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
K	-254	SER	-	expression tag	UNP P42212
K	-253	HIS	-	expression tag	UNP P42212
K	-252	VAL	-	expression tag	UNP P42212
K	-207	LEU	PHE	conflict	UNP P42212
K	-189	LEU	PHE	conflict	UNP P42212
K	-188	GLY	SER	conflict	UNP P42212
K	-185	LEU	VAL	conflict	UNP P42212
K	-181	ALA	SER	conflict	UNP P42212
K	-100	THR	MET	conflict	UNP P42212
K	-90	ALA	VAL	conflict	UNP P42212
K	-78	GLY	SER	conflict	UNP P42212
K	-50	TYR	THR	conflict	UNP P42212
K	-47	LYS	ALA	conflict	UNP P42212
K	-22	LEU	HIS	conflict	UNP P42212
K	-14	SER	-	linker	UNP P42212
K	-13	GLY	-	linker	UNP P42212
K	-12	LEU	-	linker	UNP P42212
K	-11	ARG	-	linker	UNP P42212
K	-10	SER	-	linker	UNP P42212
K	-9	ARG	-	linker	UNP P42212
K	-8	ALA	-	linker	UNP P42212
K	-7	GLN	-	linker	UNP P42212
K	-6	ALA	-	linker	UNP P42212
K	-5	SER	-	linker	UNP P42212
K	-4	ASN	-	linker	UNP P42212
K	-3	SER	-	linker	UNP P42212
K	-2	ALA	-	linker	UNP P42212
K	-1	VAL	-	linker	UNP P42212
K	0	ASP	-	linker	UNP P42212
K	287	ALA	THR	engineered mutation	UNP P08413
K	306	ALA	THR	engineered mutation	UNP P08413
K	307	ALA	THR	engineered mutation	UNP P08413
L	-272	MET	-	initiating methionine	UNP P42212
L	-271	GLY	-	expression tag	UNP P42212
L	-270	SER	-	expression tag	UNP P42212
L	-269	SER	-	expression tag	UNP P42212
L	-268	HIS	-	expression tag	UNP P42212
L	-267	HIS	-	expression tag	UNP P42212
L	-266	HIS	-	expression tag	UNP P42212
L	-265	HIS	-	expression tag	UNP P42212
L	-264	HIS	-	expression tag	UNP P42212
L	-263	HIS	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
L	-262	SER	-	expression tag	UNP P42212
L	-261	SER	-	expression tag	UNP P42212
L	-260	GLY	-	expression tag	UNP P42212
L	-259	LEU	-	expression tag	UNP P42212
L	-258	VAL	-	expression tag	UNP P42212
L	-257	PRO	-	expression tag	UNP P42212
L	-256	ARG	-	expression tag	UNP P42212
L	-255	GLY	-	expression tag	UNP P42212
L	-254	SER	-	expression tag	UNP P42212
L	-253	HIS	-	expression tag	UNP P42212
L	-252	VAL	-	expression tag	UNP P42212
L	-207	LEU	PHE	conflict	UNP P42212
L	-189	LEU	PHE	conflict	UNP P42212
L	-188	GLY	SER	conflict	UNP P42212
L	-185	LEU	VAL	conflict	UNP P42212
L	-181	ALA	SER	conflict	UNP P42212
L	-100	THR	MET	conflict	UNP P42212
L	-90	ALA	VAL	conflict	UNP P42212
L	-78	GLY	SER	conflict	UNP P42212
L	-50	TYR	THR	conflict	UNP P42212
L	-47	LYS	ALA	conflict	UNP P42212
L	-22	LEU	HIS	conflict	UNP P42212
L	-14	SER	-	linker	UNP P42212
L	-13	GLY	-	linker	UNP P42212
L	-12	LEU	-	linker	UNP P42212
L	-11	ARG	-	linker	UNP P42212
L	-10	SER	-	linker	UNP P42212
L	-9	ARG	-	linker	UNP P42212
L	-8	ALA	-	linker	UNP P42212
L	-7	GLN	-	linker	UNP P42212
L	-6	ALA	-	linker	UNP P42212
L	-5	SER	-	linker	UNP P42212
L	-4	ASN	-	linker	UNP P42212
L	-3	SER	-	linker	UNP P42212
L	-2	ALA	-	linker	UNP P42212
L	-1	VAL	-	linker	UNP P42212
L	0	ASP	-	linker	UNP P42212
L	287	ALA	THR	engineered mutation	UNP P08413
L	306	ALA	THR	engineered mutation	UNP P08413
L	307	ALA	THR	engineered mutation	UNP P08413
N	-272	MET	-	initiating methionine	UNP P42212
N	-271	GLY	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
N	-270	SER	-	expression tag	UNP P42212
N	-269	SER	-	expression tag	UNP P42212
N	-268	HIS	-	expression tag	UNP P42212
N	-267	HIS	-	expression tag	UNP P42212
N	-266	HIS	-	expression tag	UNP P42212
N	-265	HIS	-	expression tag	UNP P42212
N	-264	HIS	-	expression tag	UNP P42212
N	-263	HIS	-	expression tag	UNP P42212
N	-262	SER	-	expression tag	UNP P42212
N	-261	SER	-	expression tag	UNP P42212
N	-260	GLY	-	expression tag	UNP P42212
N	-259	LEU	-	expression tag	UNP P42212
N	-258	VAL	-	expression tag	UNP P42212
N	-257	PRO	-	expression tag	UNP P42212
N	-256	ARG	-	expression tag	UNP P42212
N	-255	GLY	-	expression tag	UNP P42212
N	-254	SER	-	expression tag	UNP P42212
N	-253	HIS	-	expression tag	UNP P42212
N	-252	VAL	-	expression tag	UNP P42212
N	-207	LEU	PHE	conflict	UNP P42212
N	-189	LEU	PHE	conflict	UNP P42212
N	-188	GLY	SER	conflict	UNP P42212
N	-185	LEU	VAL	conflict	UNP P42212
N	-181	ALA	SER	conflict	UNP P42212
N	-100	THR	MET	conflict	UNP P42212
N	-90	ALA	VAL	conflict	UNP P42212
N	-78	GLY	SER	conflict	UNP P42212
N	-50	TYR	THR	conflict	UNP P42212
N	-47	LYS	ALA	conflict	UNP P42212
N	-22	LEU	HIS	conflict	UNP P42212
N	-14	SER	-	linker	UNP P42212
N	-13	GLY	-	linker	UNP P42212
N	-12	LEU	-	linker	UNP P42212
N	-11	ARG	-	linker	UNP P42212
N	-10	SER	-	linker	UNP P42212
N	-9	ARG	-	linker	UNP P42212
N	-8	ALA	-	linker	UNP P42212
N	-7	GLN	-	linker	UNP P42212
N	-6	ALA	-	linker	UNP P42212
N	-5	SER	-	linker	UNP P42212
N	-4	ASN	-	linker	UNP P42212
N	-3	SER	-	linker	UNP P42212

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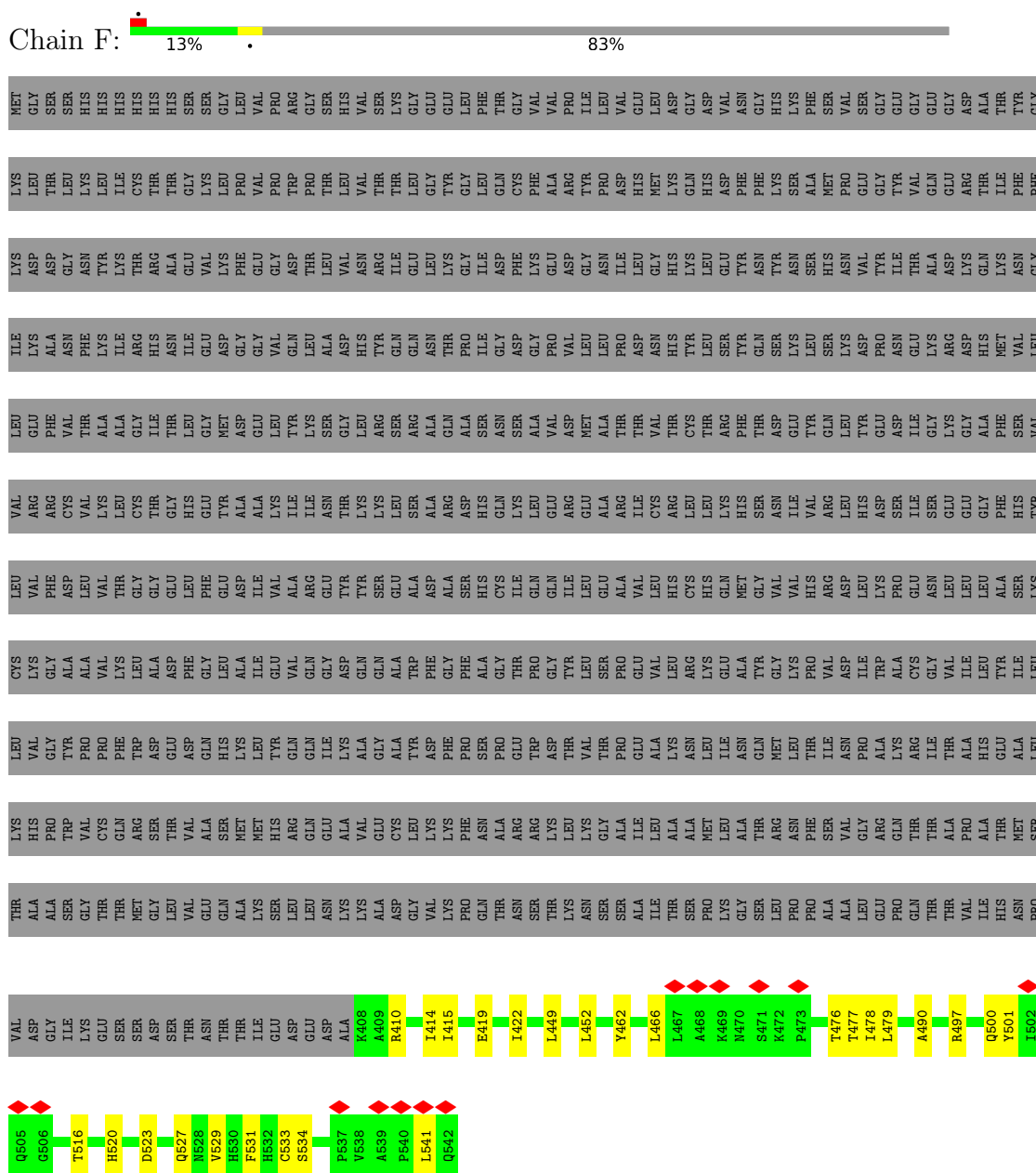
Chain	Residue	Modelled	Actual	Comment	Reference
N	-2	ALA	-	linker	UNP P42212
N	-1	VAL	-	linker	UNP P42212
N	0	ASP	-	linker	UNP P42212
N	287	ALA	THR	engineered mutation	UNP P08413
N	306	ALA	THR	engineered mutation	UNP P08413
N	307	ALA	THR	engineered mutation	UNP P08413

Chain B: 15% 83%

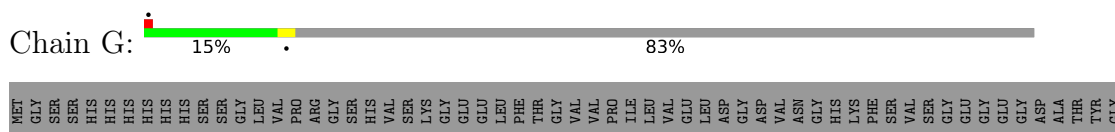




- Molecule 1: Venus-tagged CaMKII beta holoenzyme mutant



- Molecule 1: Venus-tagged CaMKII beta holoenzyme mutant

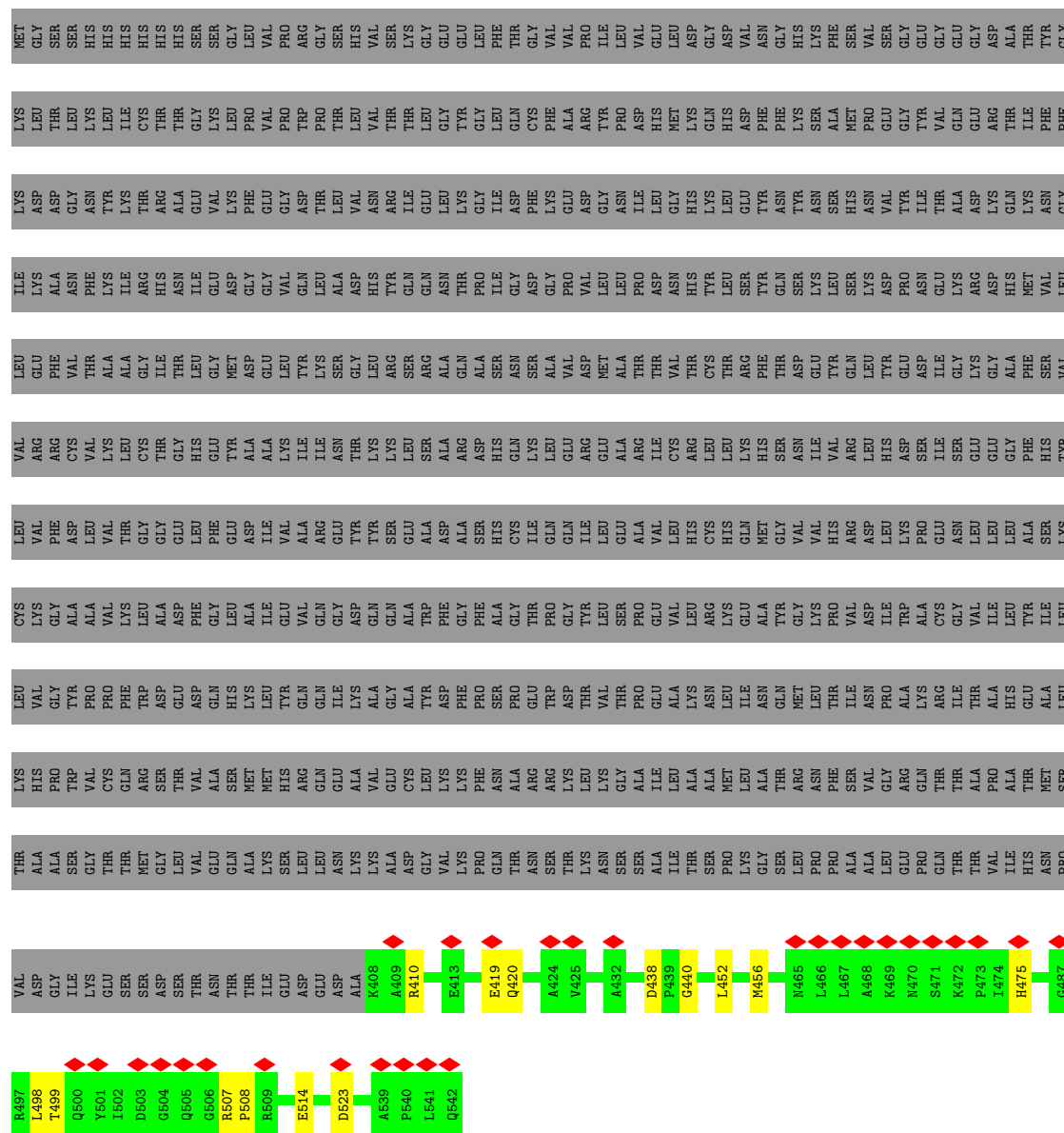






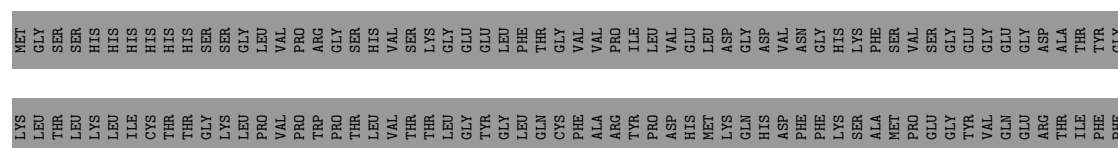
- Molecule 1: Venus-tagged CaMKII beta holoenzyme mutant

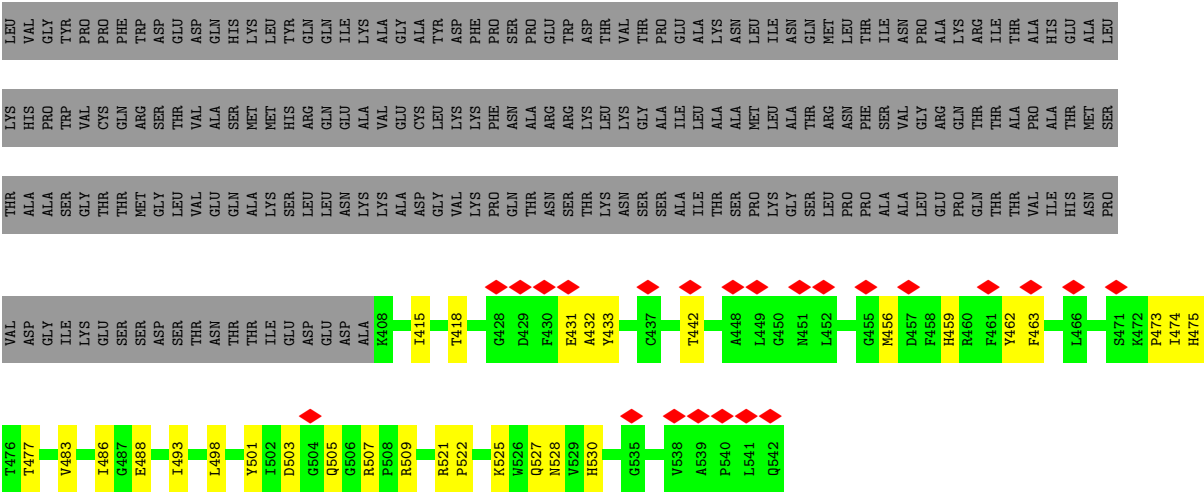
Chain J: 15% 83%



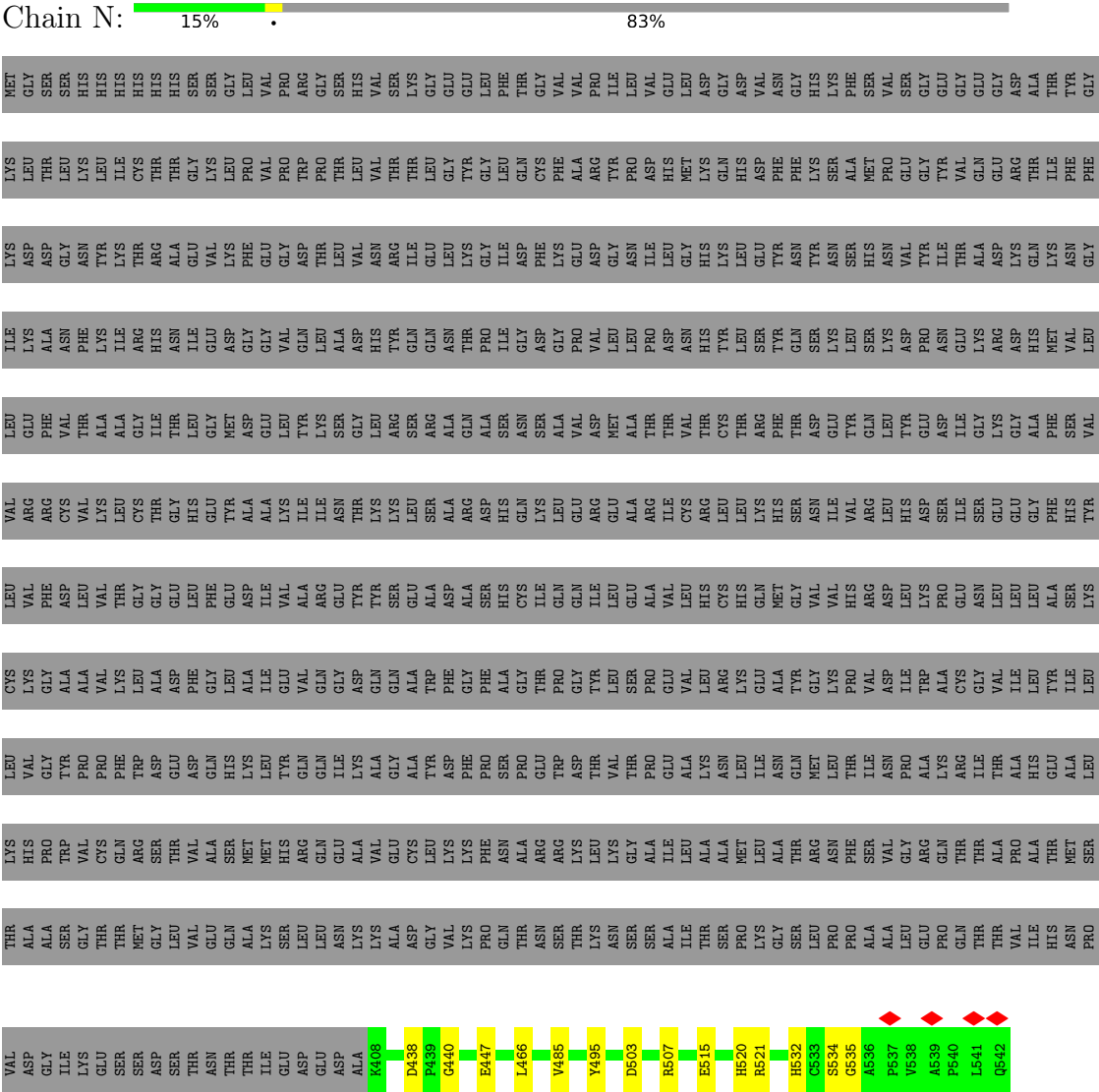
- Molecule 1: Venus-tagged CaMKII beta holoenzyme mutant

Chain K: 15% 83%





● Molecule 1: Venus-tagged CaMKII beta holoenzyme mutant



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	332360	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.068	Depositor
Minimum map value	-0.680	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.020	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	275.2, 275.2, 275.2	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.86, 0.86, 0.86	Depositor

5 Model quality

5.1 Standard geometry

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	0.12	0/1107	0.30	0/1501
1	B	0.12	0/1107	0.30	0/1501
1	C	0.12	0/1107	0.26	0/1501
1	D	0.12	0/1107	0.30	0/1501
1	F	0.14	0/1107	0.34	0/1501
1	G	0.12	0/1107	0.27	0/1501
1	H	0.20	0/1107	0.47	2/1501 (0.1%)
1	I	0.11	0/1107	0.27	0/1501
1	J	0.11	0/1107	0.28	0/1501
1	K	0.13	0/1107	0.30	0/1501
1	L	0.14	0/1107	0.37	0/1501
1	N	0.12	0/1107	0.28	0/1501
All	All	0.13	0/13284	0.32	2/18012 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	508	PRO	CA-N-CD	-10.67	97.06	112.00
1	H	508	PRO	N-CD-CG	-5.87	94.40	103.20

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	1079	1045	1043	21	0
1	B	1079	1045	1043	15	0
1	C	1079	1045	1043	21	0
1	D	1079	1045	1043	15	0
1	F	1079	1045	1043	18	0
1	G	1079	1045	1043	9	0
1	H	1079	1044	1043	15	0
1	I	1079	1044	1043	8	0
1	J	1079	1044	1043	10	0
1	K	1079	1044	1043	12	0
1	L	1079	1044	1043	25	0
1	N	1079	1044	1043	11	0
All	All	12948	12534	12516	165	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 (165) 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:C:497:ARG:NE	1:C:515:GLU:OE2	2.15	0.79
1:L:488:GLU:OE1	1:L:521:ARG:NH1	2.16	0.78
1:B:488:GLU:N	1:B:488:GLU:OE1	2.18	0.77
1:L:527:GLN:N	1:L:527:GLN:OE1	2.18	0.76
1:F:533:CYS:SG	1:F:534:SER:N	2.61	0.74
1:C:437:CYS:O	1:C:527:GLN:NE2	2.21	0.74
1:B:423:GLU:OE2	1:B:427:ASN:ND2	2.20	0.73
1:H:508:PRO:O	1:H:509:ARG:NE	2.21	0.73
1:L:483:VAL:HG22	1:L:493:ILE:HD13	1.71	0.72
1:H:488:GLU:O	1:H:521:ARG:NH2	2.23	0.71
1:B:457:ASP:OD2	1:C:475:HIS:NE2	2.24	0.69
1:J:456:MET:N	1:J:456:MET:SD	2.66	0.69
1:H:499:THR:OG1	1:H:511:SER:OG	2.05	0.69
1:H:533:CYS:SG	1:H:534:SER:N	2.66	0.69
1:K:408:LYS:O	1:K:412:GLN:NE2	2.26	0.69
1:H:508:PRO:O	1:H:508:PRO:HD2	1.92	0.68
1:B:532:HIS:NE2	1:I:534:SER:OG	2.26	0.67
1:I:503:ASP:OD2	1:I:507:ARG:NH2	2.28	0.66
1:I:514:GLU:OE1	1:I:514:GLU:N	2.31	0.64
1:A:470:ASN:ND2	1:A:472:LYS:O	2.32	0.63
1:L:503:ASP:OD1	1:L:507:ARG:N	2.32	0.62
1:D:534:SER:OG	1:N:532:HIS:NE2	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:520:HIS:CG	1:H:486:ILE:HD12	2.34	0.62
1:B:534:SER:OG	1:I:532:HIS:NE2	2.32	0.62
1:I:523:ASP:N	1:I:523:ASP:OD1	2.33	0.61
1:J:498:LEU:HD12	1:J:499:THR:N	2.15	0.61
1:L:433:TYR:OH	1:L:456:MET:HE3	2.00	0.60
1:N:447:GLU:OE1	1:N:447:GLU:N	2.31	0.60
1:K:488:GLU:N	1:K:488:GLU:OE1	2.35	0.60
1:G:514:GLU:N	1:G:514:GLU:OE1	2.35	0.59
1:K:423:GLU:OE2	1:K:427:ASN:ND2	2.36	0.59
1:A:503:ASP:OD1	1:A:507:ARG:N	2.35	0.59
1:A:538:VAL:HG12	1:A:540:PRO:HD2	1.85	0.58
1:A:512:GLN:NE2	1:A:514:GLU:OE2	2.35	0.58
1:J:514:GLU:N	1:J:514:GLU:OE1	2.36	0.58
1:C:534:SER:OG	1:H:532:HIS:NE2	2.29	0.57
1:G:438:ASP:OD1	1:G:440:GLY:N	2.34	0.57
1:B:523:ASP:O	1:B:523:ASP:OD1	2.22	0.57
1:L:473:PRO:O	1:L:474:ILE:HD13	2.04	0.57
1:A:509:ARG:O	1:C:462:TYR:OH	2.22	0.56
1:A:506:GLY:O	1:C:469:LYS:NZ	2.36	0.56
1:G:503:ASP:OD1	1:G:507:ARG:N	2.36	0.56
1:H:514:GLU:N	1:H:514:GLU:OE1	2.39	0.56
1:C:514:GLU:N	1:C:514:GLU:OE1	2.39	0.56
1:J:452:LEU:O	1:J:452:LEU:HD23	2.06	0.56
1:K:534:SER:OG	1:K:535:GLY:N	2.37	0.56
1:K:523:ASP:O	1:K:523:ASP:OD1	2.25	0.55
1:L:483:VAL:CG2	1:L:493:ILE:HD13	2.37	0.55
1:D:410:ARG:NH1	1:D:488:GLU:O	2.39	0.55
1:K:488:GLU:O	1:K:521:ARG:NH2	2.39	0.55
1:L:474:ILE:HD12	1:L:501:TYR:CB	2.37	0.55
1:D:523:ASP:OD2	1:D:523:ASP:O	2.24	0.55
1:D:488:GLU:N	1:D:488:GLU:OE1	2.39	0.55
1:F:520:HIS:O	1:F:527:GLN:N	2.39	0.55
1:I:466:LEU:O	1:I:469:LYS:NZ	2.40	0.55
1:F:490:ALA:HB2	1:L:486:ILE:HG21	1.89	0.54
1:L:501:TYR:O	1:L:509:ARG:N	2.40	0.54
1:B:429:ASP:OD2	1:B:431:GLU:N	2.40	0.54
1:B:466:LEU:C	1:B:466:LEU:HD23	2.32	0.54
1:D:470:ASN:OD1	1:D:472:LYS:N	2.40	0.54
1:F:478:ILE:C	1:F:479:LEU:HD12	2.33	0.54
1:G:495:TYR:OH	1:G:515:GLU:OE1	2.24	0.54
1:D:466:LEU:HD23	1:D:467:LEU:N	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:503:ASP:OD1	1:H:506:GLY:N	2.42	0.52
1:A:470:ASN:O	1:A:472:LYS:HG3	2.10	0.52
1:K:442:THR:HB	1:K:452:LEU:HD21	1.90	0.52
1:B:514:GLU:N	1:B:514:GLU:OE2	2.43	0.52
1:C:430:PHE:CE2	1:C:434:ALA:HB2	2.44	0.52
1:F:523:ASP:OD1	1:F:523:ASP:C	2.53	0.52
1:G:523:ASP:OD1	1:G:523:ASP:C	2.54	0.51
1:F:516:THR:HG21	1:L:530:HIS:CD2	2.46	0.51
1:L:415:ILE:O	1:L:418:THR:OG1	2.28	0.51
1:D:514:GLU:N	1:D:514:GLU:OE1	2.44	0.51
1:J:523:ASP:CG	1:J:523:ASP:O	2.54	0.51
1:L:474:ILE:HD12	1:L:501:TYR:HB3	1.92	0.50
1:K:457:ASP:OD2	1:K:457:ASP:C	2.54	0.50
1:F:501:TYR:OH	1:F:541:LEU:O	2.24	0.50
1:L:477:THR:N	1:L:498:LEU:O	2.44	0.50
1:A:502:ILE:HG21	1:C:466:LEU:HD13	1.94	0.50
1:B:464:GLU:N	1:B:464:GLU:OE2	2.45	0.50
1:H:477:THR:N	1:H:498:LEU:O	2.39	0.49
1:C:528:ASN:ND2	1:C:530:HIS:O	2.39	0.49
1:F:415:ILE:HG22	1:F:419:GLU:OE2	2.12	0.49
1:H:488:GLU:OE1	1:H:488:GLU:N	2.46	0.49
1:B:469:LYS:O	1:B:472:LYS:NZ	2.45	0.49
1:J:438:ASP:OD1	1:J:440:GLY:N	2.41	0.48
1:A:437:CYS:SG	1:A:528:ASN:ND2	2.84	0.48
1:J:507:ARG:NH1	1:J:508:PRO:O	2.46	0.48
1:F:477:THR:OG1	1:F:500:GLN:NE2	2.47	0.48
1:F:529:VAL:HG21	1:L:486:ILE:HD11	1.96	0.48
1:N:466:LEU:C	1:N:466:LEU:HD23	2.39	0.48
1:N:503:ASP:OD2	1:N:507:ARG:NH2	2.47	0.47
1:A:538:VAL:HG12	1:A:540:PRO:CD	2.44	0.47
1:B:429:ASP:OD2	1:B:429:ASP:C	2.56	0.47
1:H:496:ILE:HG23	1:H:514:GLU:HG3	1.95	0.47
1:C:438:ASP:OD1	1:C:438:ASP:C	2.58	0.47
1:K:523:ASP:O	1:K:525:LYS:N	2.48	0.47
1:D:486:ILE:HG23	1:N:520:HIS:CD2	2.50	0.47
1:L:442:THR:O	1:L:530:HIS:ND1	2.48	0.46
1:I:419:GLU:OE1	1:I:420:GLN:N	2.48	0.46
1:N:495:TYR:OH	1:N:515:GLU:OE1	2.20	0.46
1:N:534:SER:OG	1:N:535:GLY:N	2.48	0.46
1:C:419:GLU:OE1	1:C:420:GLN:N	2.48	0.46
1:H:499:THR:O	1:H:510:THR:HG23	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:463:PHE:CD2	1:C:467:LEU:HD12	2.51	0.46
1:A:488:GLU:OE1	1:A:488:GLU:N	2.49	0.46
1:A:416:LYS:HE3	1:A:416:LYS:HA	1.98	0.46
1:C:436:ILE:N	1:C:436:ILE:HD13	2.30	0.46
1:C:520:HIS:O	1:C:527:GLN:N	2.44	0.45
1:C:453:VAL:HG23	1:C:453:VAL:O	2.15	0.45
1:L:505:GLN:OE1	1:L:505:GLN:N	2.44	0.45
1:F:479:LEU:HD12	1:F:479:LEU:N	2.31	0.45
1:J:410:ARG:NH2	1:J:490:ALA:O	2.50	0.45
1:N:438:ASP:OD1	1:N:440:GLY:N	2.43	0.45
1:F:462:TYR:O	1:F:466:LEU:HD23	2.17	0.45
1:G:503:ASP:OD2	1:G:507:ARG:NH1	2.50	0.45
1:D:523:ASP:OD2	1:D:523:ASP:C	2.61	0.44
1:L:528:ASN:ND2	1:L:530:HIS:O	2.45	0.44
1:D:466:LEU:HD23	1:D:466:LEU:C	2.41	0.44
1:D:466:LEU:HD23	1:D:467:LEU:HA	1.99	0.44
1:N:521:ARG:O	1:N:521:ARG:HG2	2.17	0.44
1:L:525:LYS:CG	1:L:525:LYS:O	2.65	0.44
1:A:444:PHE:CD1	1:J:496:ILE:HD11	2.53	0.44
1:L:431:GLU:OE1	1:L:432:ALA:N	2.51	0.44
1:A:484:HIS:ND1	1:A:492:CYS:SG	2.87	0.44
1:J:419:GLU:OE1	1:J:420:GLN:N	2.50	0.44
1:N:485:VAL:O	1:N:485:VAL:HG13	2.18	0.44
1:F:410:ARG:O	1:F:414:ILE:HG13	2.17	0.44
1:B:523:ASP:O	1:B:525:LYS:N	2.51	0.43
1:N:521:ARG:O	1:N:521:ARG:CG	2.66	0.43
1:C:430:PHE:CE2	1:C:460:ARG:HG2	2.54	0.43
1:F:449:LEU:HD12	1:F:449:LEU:N	2.33	0.43
1:K:523:ASP:OD1	1:K:523:ASP:C	2.61	0.43
1:D:466:LEU:HD23	1:D:467:LEU:CA	2.48	0.43
1:A:470:ASN:O	1:A:472:LYS:HE2	2.18	0.43
1:A:533:CYS:SG	1:A:534:SER:N	2.91	0.43
1:F:531:PHE:CD1	1:F:531:PHE:C	2.97	0.42
1:K:496:ILE:HG22	1:K:497:ARG:N	2.34	0.42
1:D:496:ILE:HG22	1:D:497:ARG:N	2.35	0.42
1:F:452:LEU:C	1:F:452:LEU:HD23	2.44	0.42
1:B:523:ASP:OD1	1:B:523:ASP:C	2.63	0.42
1:C:486:ILE:HG21	1:H:490:ALA:HB2	2.02	0.42
1:F:422:ILE:HG23	1:F:497:ARG:NH1	2.35	0.42
1:G:508:PRO:O	1:G:509:ARG:NH2	2.52	0.42
1:C:485:VAL:O	1:C:485:VAL:HG13	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:473:PRO:O	1:D:474:ILE:HG23	2.20	0.42
1:I:457:ASP:OD1	1:I:460:ARG:NH2	2.53	0.42
1:C:497:ARG:CZ	1:C:515:GLU:OE2	2.68	0.42
1:A:413:GLU:O	1:A:417:THR:HG23	2.18	0.42
1:L:521:ARG:O	1:L:522:PRO:C	2.62	0.42
1:H:480:ASN:OD1	1:H:480:ASN:C	2.63	0.41
1:K:466:LEU:HD23	1:K:466:LEU:C	2.46	0.41
1:A:472:LYS:HB3	1:A:473:PRO:CD	2.50	0.41
1:A:460:ARG:O	1:A:464:GLU:N	2.45	0.41
1:D:429:ASP:OD1	1:D:430:PHE:N	2.54	0.41
1:A:531:PHE:CG	1:A:532:HIS:N	2.89	0.41
1:G:423:GLU:OE1	1:G:424:ALA:N	2.54	0.41
1:L:474:ILE:HG22	1:L:475:HIS:N	2.34	0.41
1:B:460:ARG:HH11	1:B:460:ARG:HG2	1.85	0.41
1:L:525:LYS:O	1:L:525:LYS:HG2	2.19	0.41
1:A:472:LYS:HE3	1:A:502:ILE:HB	2.03	0.41
1:L:459:HIS:HA	1:L:462:TYR:CE2	2.56	0.41
1:L:459:HIS:O	1:L:463:PHE:HB2	2.21	0.40
1:F:414:ILE:HG13	1:F:414:ILE:H	1.74	0.40
1:G:496:ILE:HG22	1:G:497:ARG:N	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	133/815 (16%)	129 (97%)	4 (3%)	0	100	100
1	B	133/815 (16%)	131 (98%)	2 (2%)	0	100	100
1	C	133/815 (16%)	127 (96%)	6 (4%)	0	100	100
1	D	133/815 (16%)	129 (97%)	4 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	133/815 (16%)	129 (97%)	4 (3%)	0	100	100
1	G	133/815 (16%)	133 (100%)	0	0	100	100
1	H	133/815 (16%)	132 (99%)	1 (1%)	0	100	100
1	I	133/815 (16%)	128 (96%)	5 (4%)	0	100	100
1	J	133/815 (16%)	133 (100%)	0	0	100	100
1	K	133/815 (16%)	125 (94%)	8 (6%)	0	100	100
1	L	133/815 (16%)	126 (95%)	7 (5%)	0	100	100
1	N	133/815 (16%)	128 (96%)	5 (4%)	0	100	100
All	All	1596/9780 (16%)	1550 (97%)	46 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	115/689 (17%)	114 (99%)	1 (1%)	70	76
1	B	115/689 (17%)	115 (100%)	0	100	100
1	C	115/689 (17%)	114 (99%)	1 (1%)	70	76
1	D	115/689 (17%)	115 (100%)	0	100	100
1	F	115/689 (17%)	114 (99%)	1 (1%)	70	76
1	G	115/689 (17%)	114 (99%)	1 (1%)	70	76
1	H	115/689 (17%)	115 (100%)	0	100	100
1	I	115/689 (17%)	114 (99%)	1 (1%)	70	76
1	J	115/689 (17%)	114 (99%)	1 (1%)	70	76
1	K	115/689 (17%)	115 (100%)	0	100	100
1	L	115/689 (17%)	115 (100%)	0	100	100
1	N	115/689 (17%)	115 (100%)	0	100	100
All	All	1380/8268 (17%)	1374 (100%)	6 (0%)	81	81

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	530	HIS
1	C	514	GLU
1	F	476	THR
1	G	421	LEU
1	I	534	SER
1	J	475	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	427	ASN
1	A	470	ASN
1	A	482	HIS
1	A	528	ASN
1	B	420	GLN
1	B	484	HIS
1	B	520	HIS
1	B	528	ASN
1	C	484	HIS
1	C	530	HIS
1	C	532	HIS
1	D	420	GLN
1	D	484	HIS
1	F	451	ASN
1	H	459	HIS
1	H	475	HIS
1	I	420	GLN
1	I	451	ASN
1	K	470	ASN
1	K	500	GLN
1	N	484	HIS
1	N	512	GLN
1	N	520	HIS

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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.

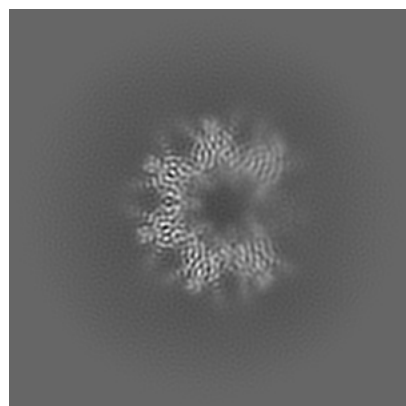
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-41077. These allow visual inspection of the internal detail of the map and identification of artifacts.

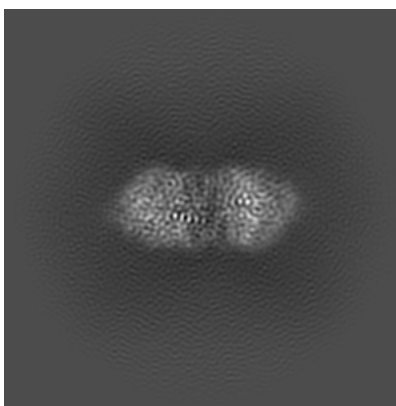
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

6.1.1 Primary map



X

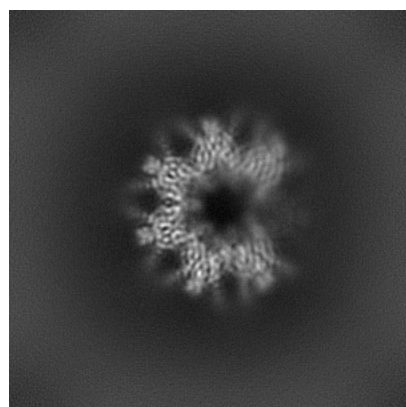


Y

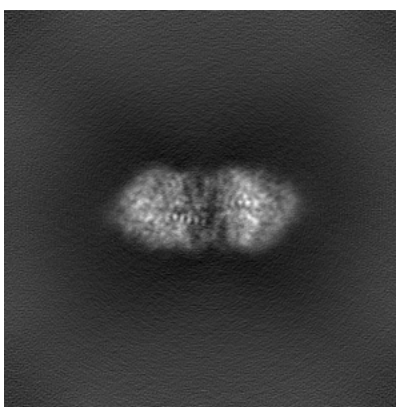


Z

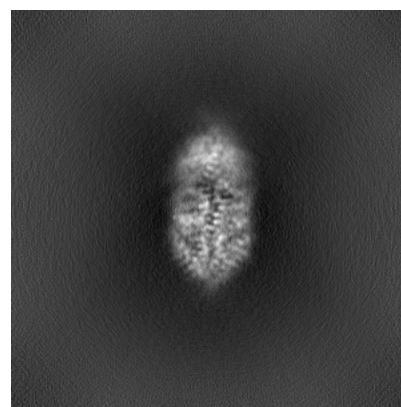
6.1.2 Raw map



X



Y

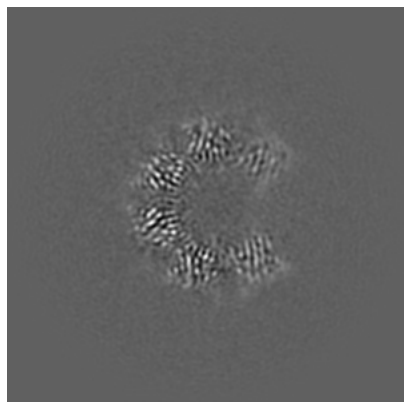


Z

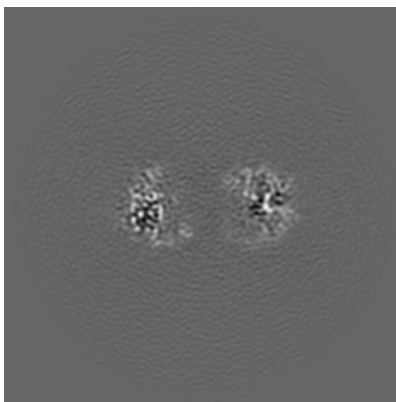
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

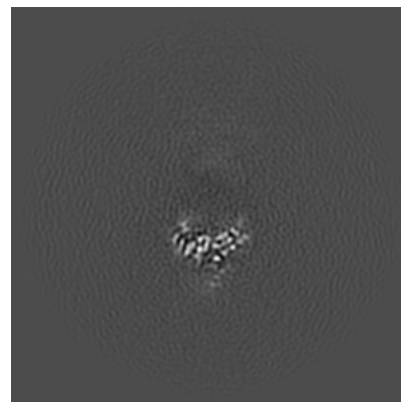
6.2.1 Primary map



X Index: 160

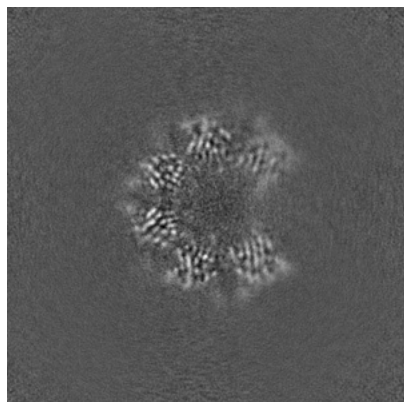


Y Index: 160

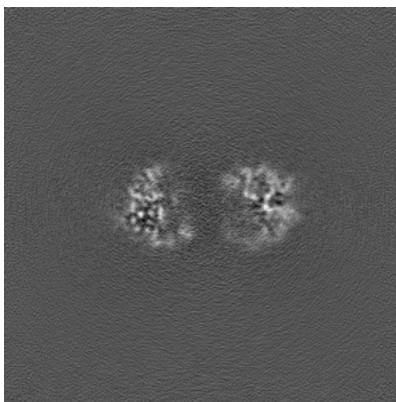


Z Index: 160

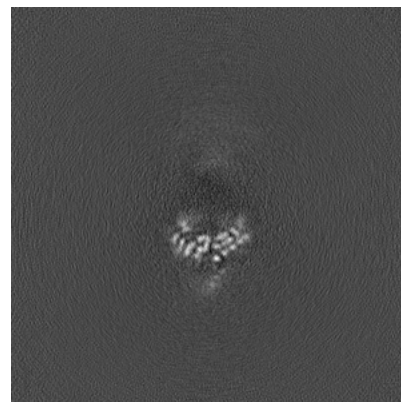
6.2.2 Raw map



X Index: 160



Y Index: 160

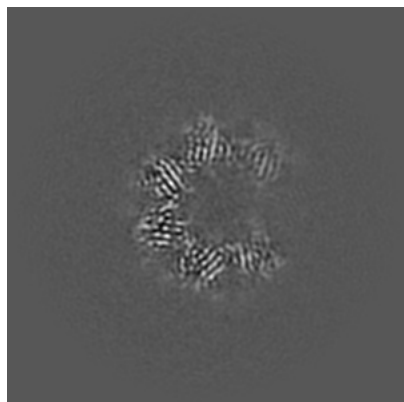


Z Index: 160

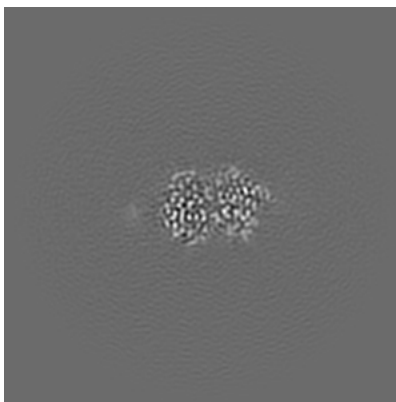
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

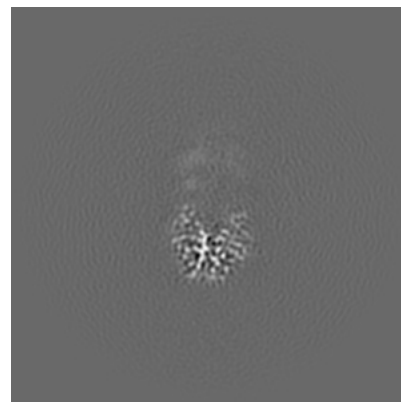
6.3.1 Primary map



X Index: 156

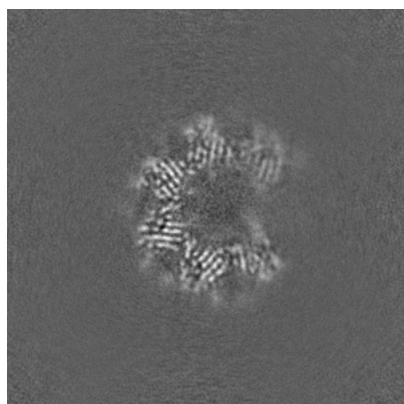


Y Index: 125

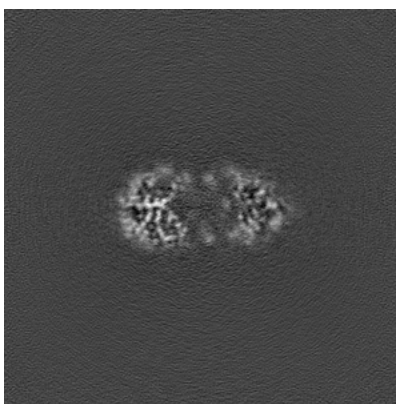


Z Index: 142

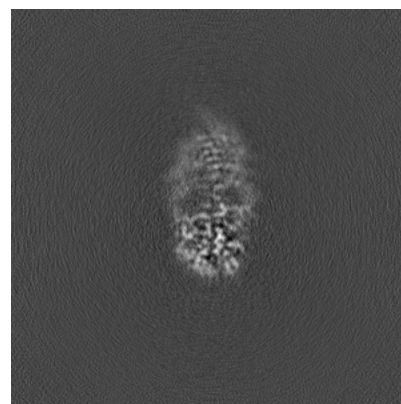
6.3.2 Raw map



X Index: 155



Y Index: 150

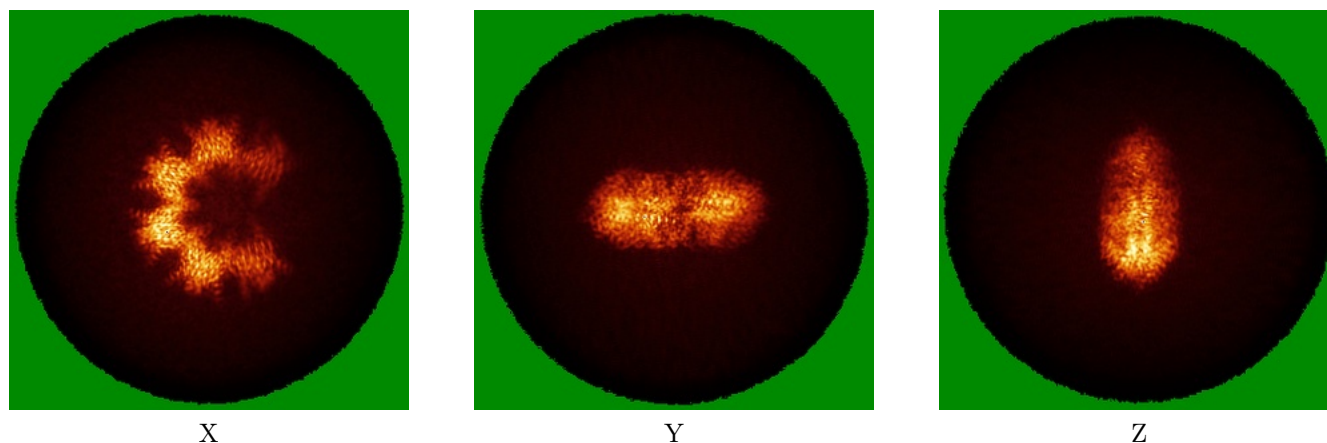


Z Index: 194

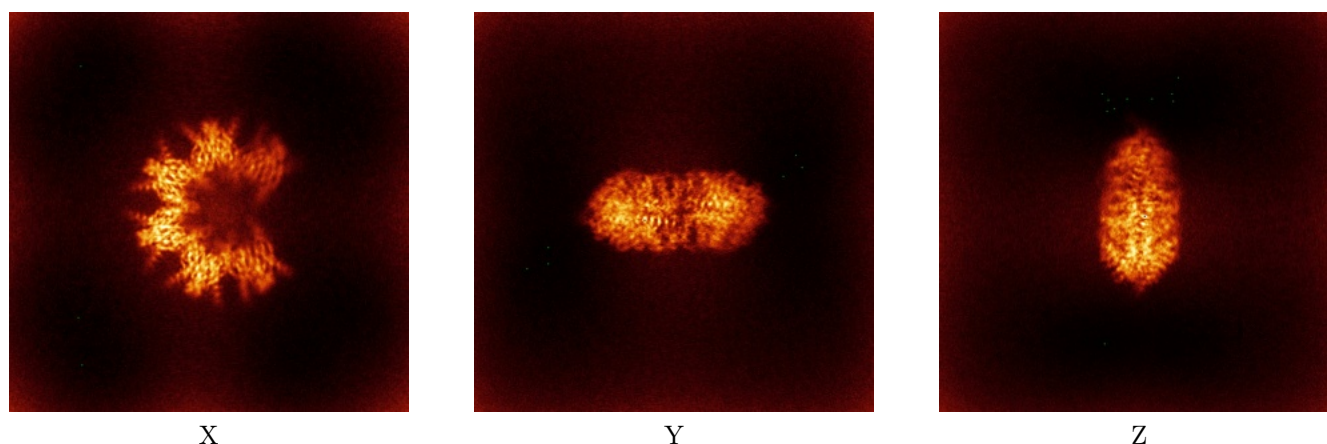
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

6.4.1 Primary map



6.4.2 Raw map



The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

This section was not generated.

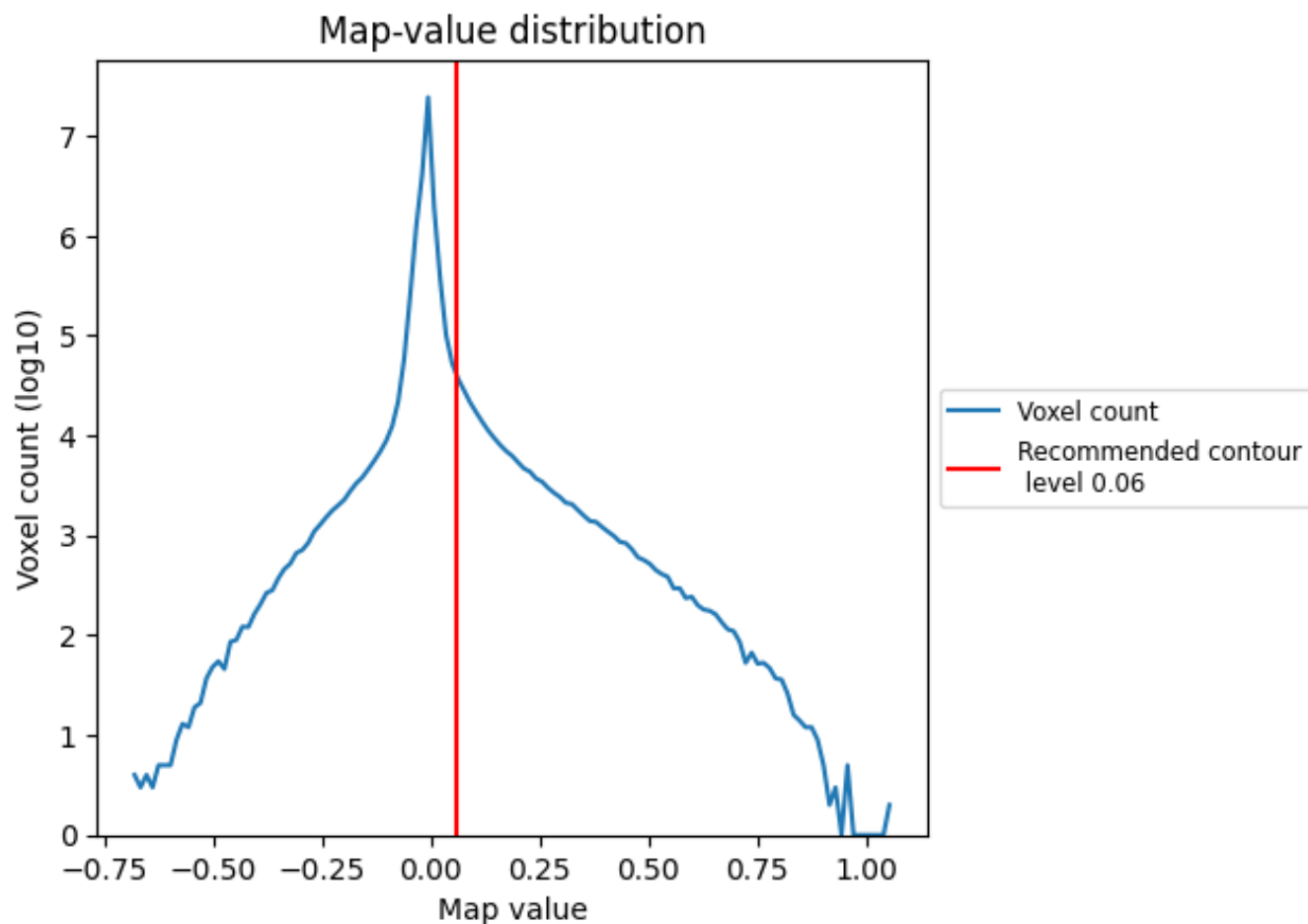
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

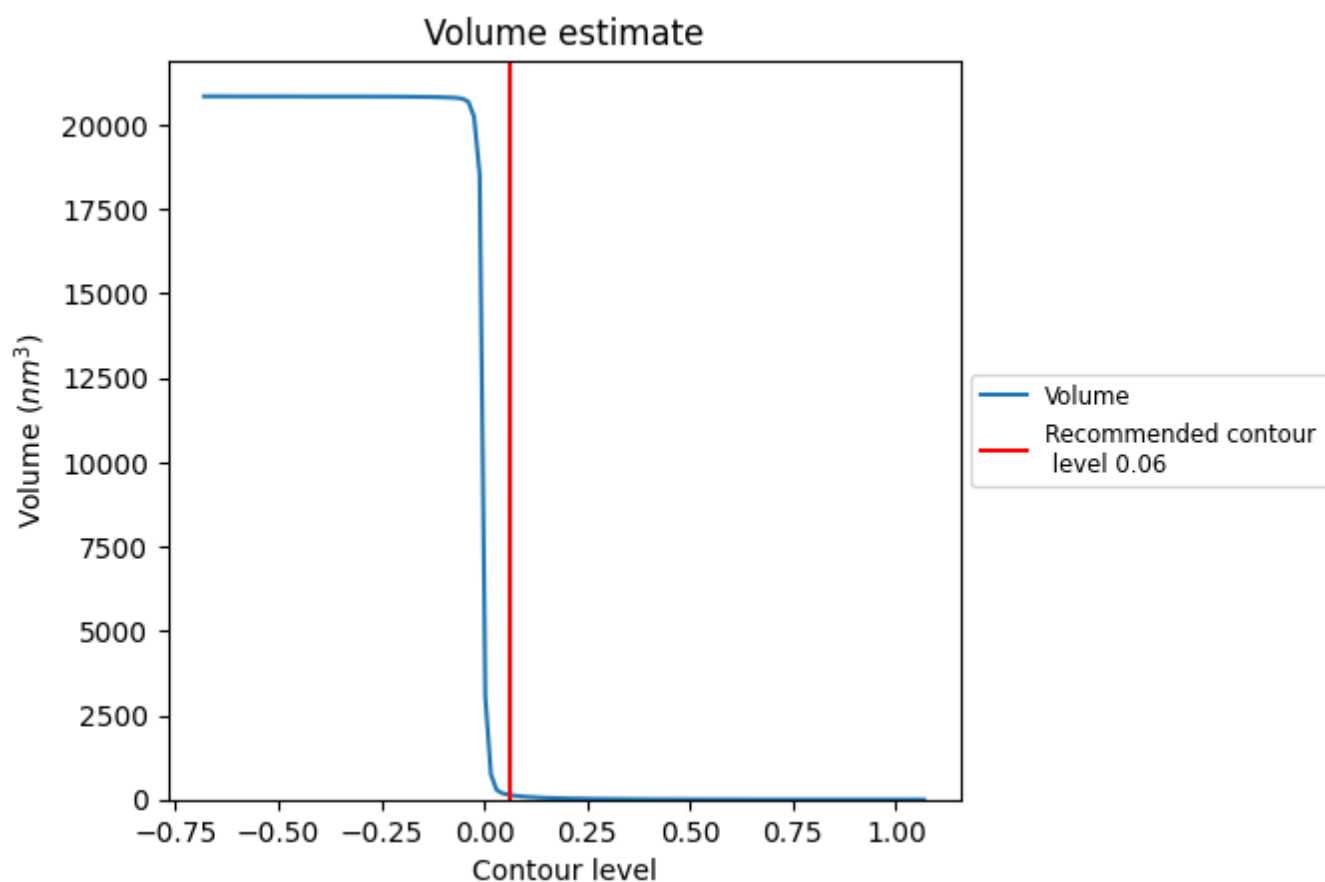
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

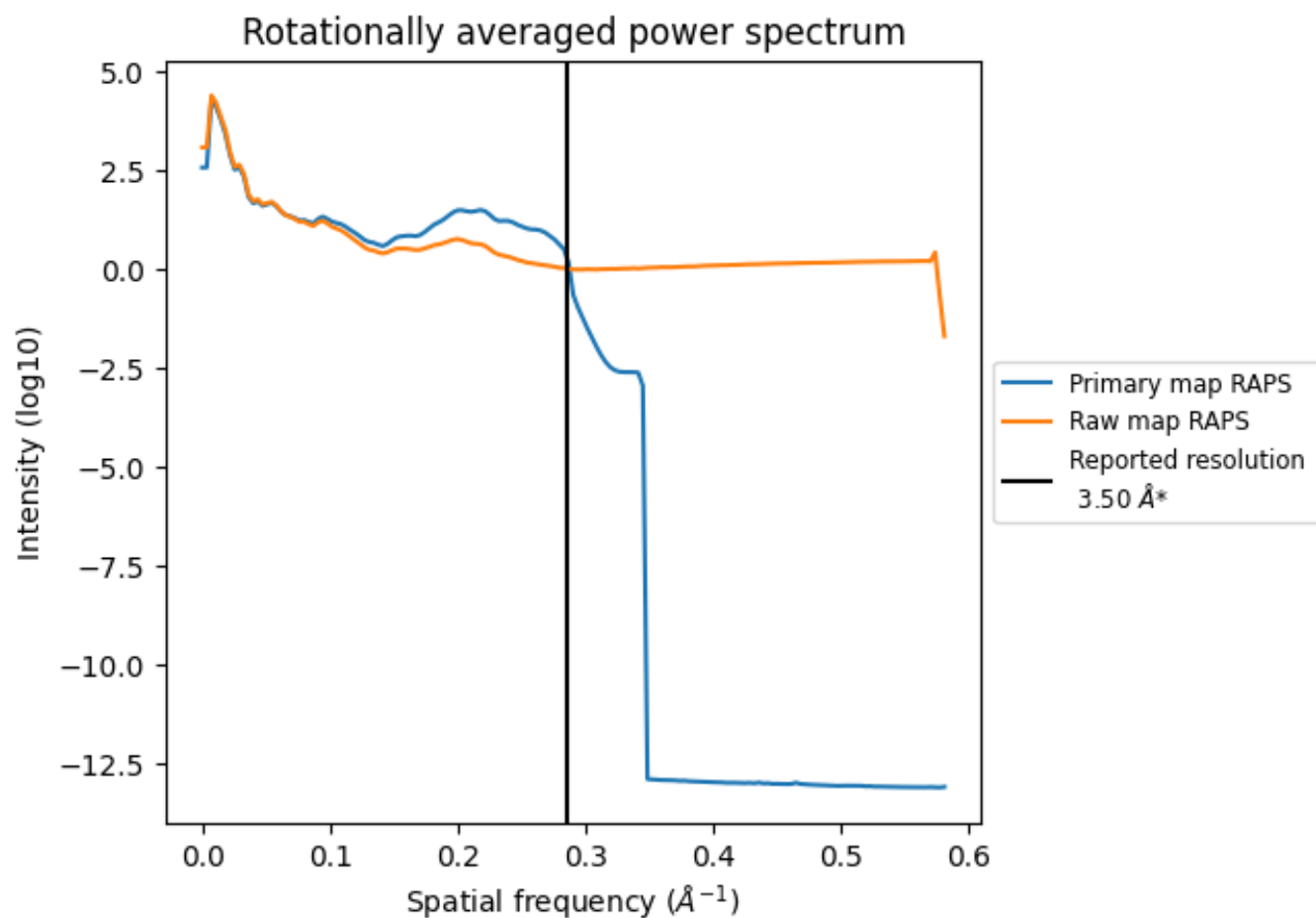
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 141 nm³; this corresponds to an approximate mass of 128 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

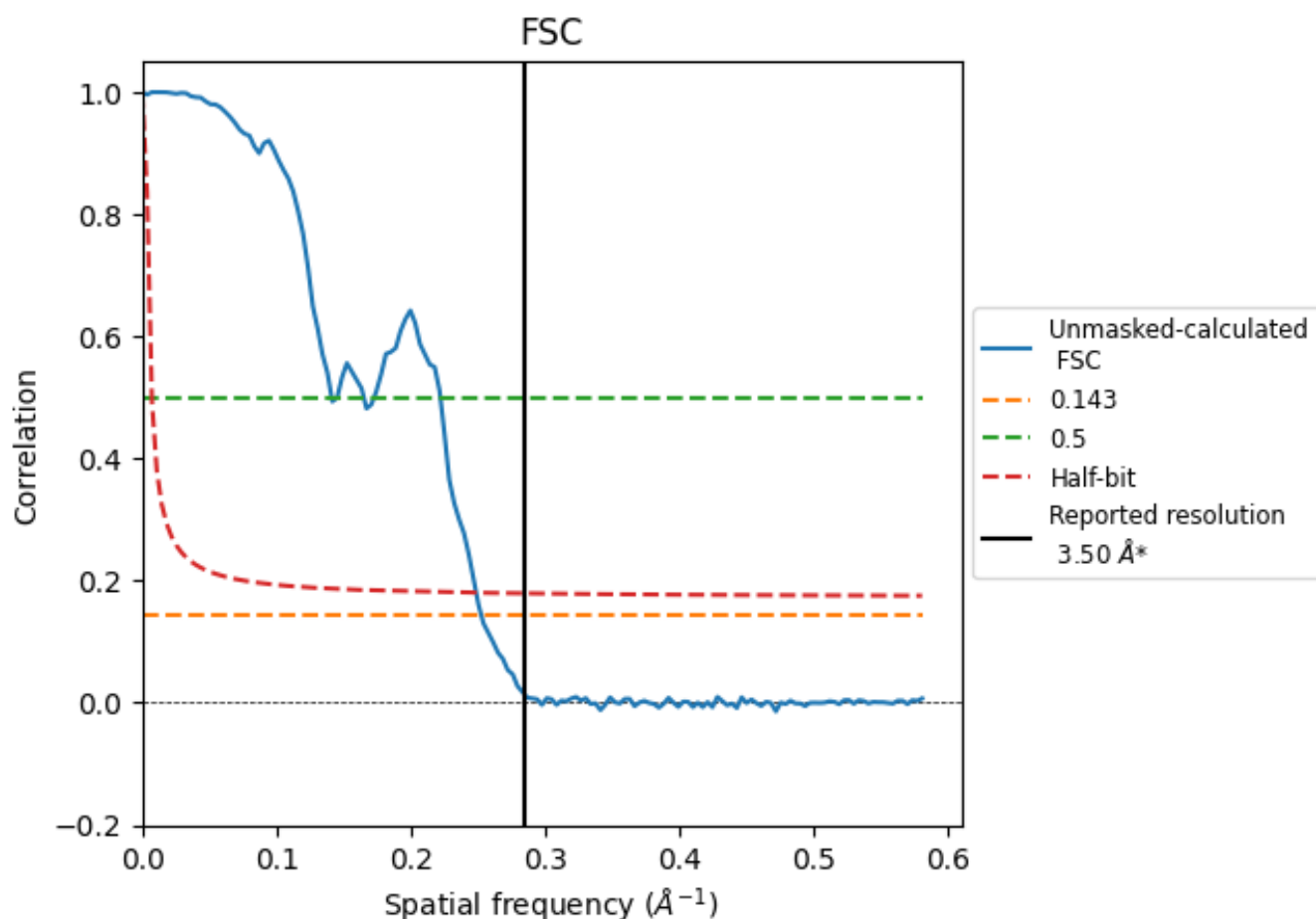


*Reported resolution corresponds to spatial frequency of 0.286 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.286 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.96	7.08	4.02

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.96 differs from the reported value 3.5 by more than 10 %

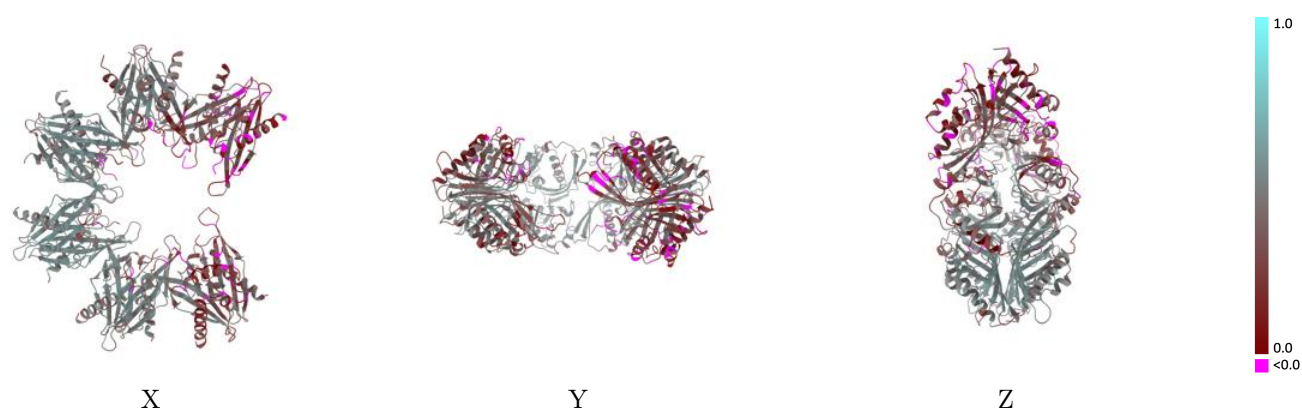
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-41077 and PDB model 8T6Q. Per-residue inclusion information can be found in section 3 on page 20.

9.1 Map-model overlay [i](#)

This section was not generated.

9.2 Q-score mapped to coordinate model [i](#)

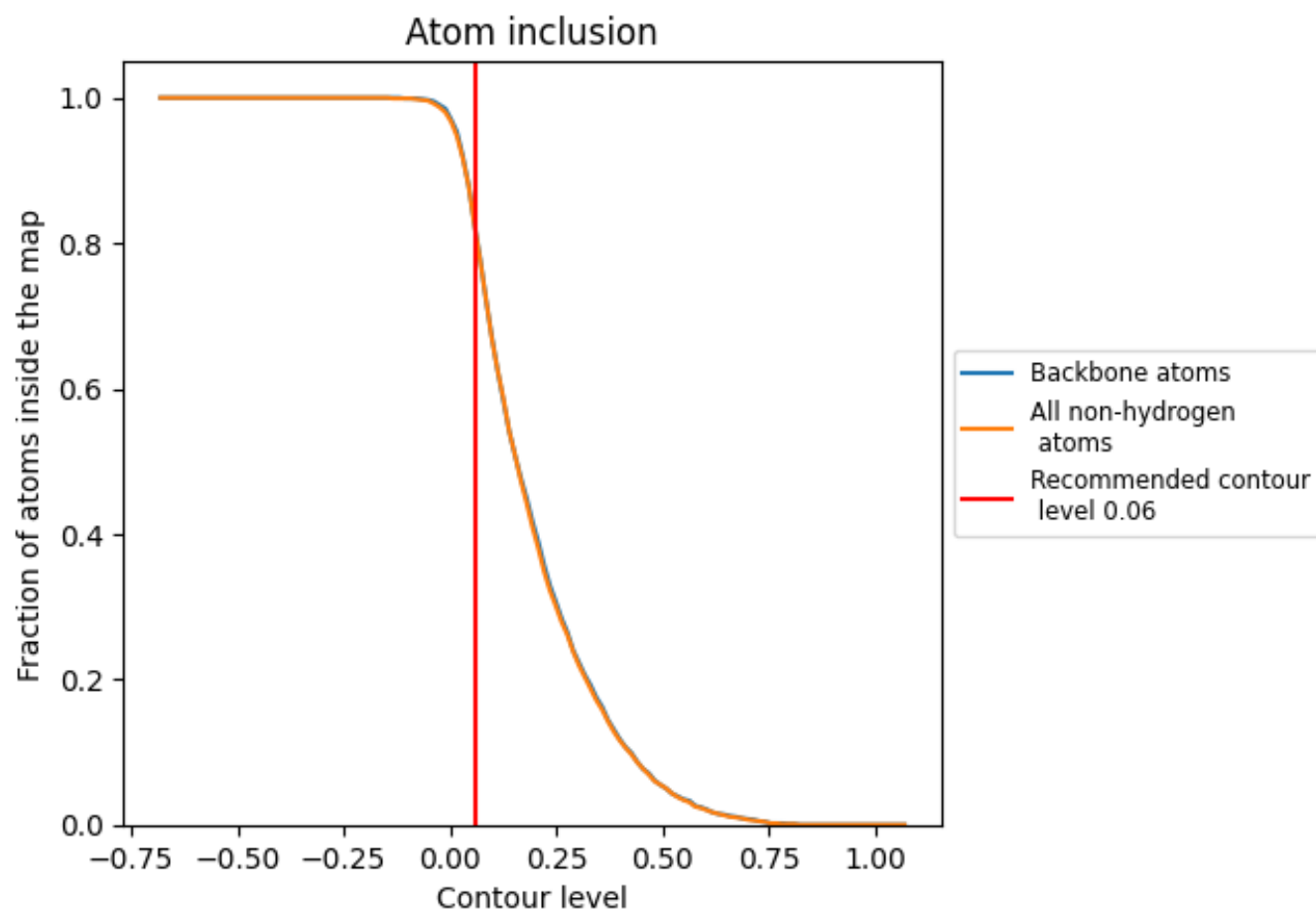


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 82% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8150	 0.3890
A	 0.5620	 0.2120
B	 0.9040	 0.4820
C	 0.8200	 0.3660
D	 0.9210	 0.5030
F	 0.8040	 0.3500
G	 0.8840	 0.4490
H	 0.8200	 0.3840
I	 0.9100	 0.4840
J	 0.6800	 0.2200
K	 0.9000	 0.4610
L	 0.7290	 0.2540
N	 0.9270	 0.5030

