



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 04:30 AM UTC

PDB ID : 8T6K / pdb_00008t6k
EMDB ID : EMD-41070
Title : Cryo-EM structure of tetradecameric CaMKII beta holoenzyme T287A T306A T307A
Authors : Chien, C.-T.; Chiu, W.; Khan, S.
Deposited on : 2023-06-16
Resolution : 3.00 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:

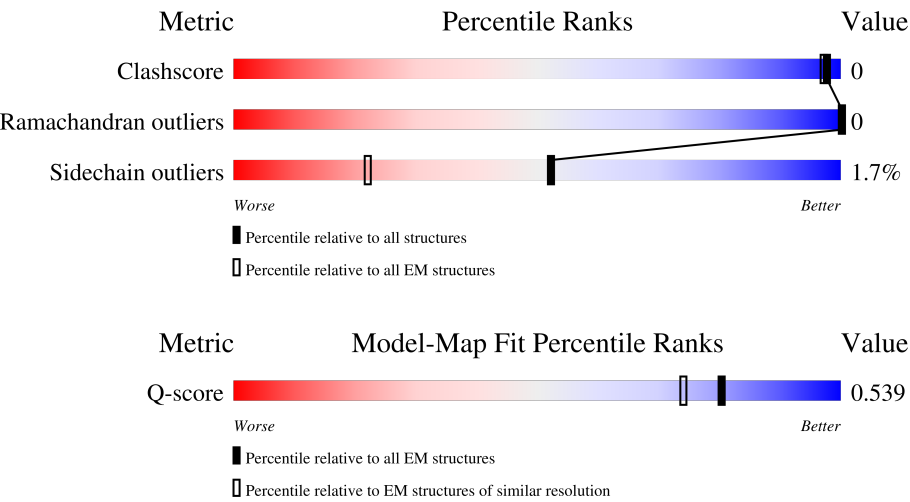
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance i

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

The reported resolution of this entry is 3.00 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.



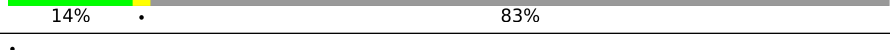
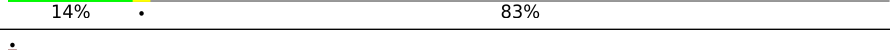
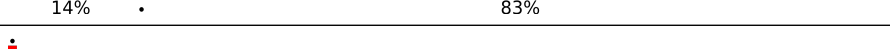
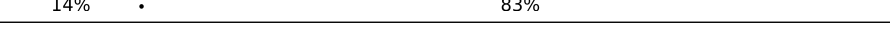
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14081 (2.50 - 3.50)

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	14% 83%
1	B	815	14% 83%
1	C	815	14% 83%
1	D	815	14% 83%

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Mol	Chain	Length	Quality of chain
1	E	815	
1	F	815	
1	G	815	
1	H	815	
1	I	815	
1	J	815	
1	K	815	
1	L	815	
1	M	815	
1	N	815	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 29722 atoms, of which 14616 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	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	B	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	C	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	D	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	E	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	F	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	G	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	H	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	I	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	J	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	K	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	L	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	M	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		
1	N	135	Total	C	H	N	O	S	0	0
			2123	681	1044	193	201	4		

There are 700 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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	SER	-	linker	UNP P42212
B	-9	ARG	-	linker	UNP P42212
B	-8	ALA	-	linker	UNP P42212
B	-7	GLN	-	linker	UNP P42212
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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
E	-272	MET	-	initiating methionine	UNP P42212
E	-271	GLY	-	expression tag	UNP P42212
E	-270	SER	-	expression tag	UNP P42212
E	-269	SER	-	expression tag	UNP P42212
E	-268	HIS	-	expression tag	UNP P42212
E	-267	HIS	-	expression tag	UNP P42212
E	-266	HIS	-	expression tag	UNP P42212
E	-265	HIS	-	expression tag	UNP P42212
E	-264	HIS	-	expression tag	UNP P42212
E	-263	HIS	-	expression tag	UNP P42212
E	-262	SER	-	expression tag	UNP P42212
E	-261	SER	-	expression tag	UNP P42212

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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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

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

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Chain	Residue	Modelled	Actual	Comment	Reference
M	-14	SER	-	linker	UNP P42212
M	-13	GLY	-	linker	UNP P42212
M	-12	LEU	-	linker	UNP P42212
M	-11	ARG	-	linker	UNP P42212
M	-10	SER	-	linker	UNP P42212
M	-9	ARG	-	linker	UNP P42212
M	-8	ALA	-	linker	UNP P42212
M	-7	GLN	-	linker	UNP P42212
M	-6	ALA	-	linker	UNP P42212
M	-5	SER	-	linker	UNP P42212
M	-4	ASN	-	linker	UNP P42212
M	-3	SER	-	linker	UNP P42212
M	-2	ALA	-	linker	UNP P42212
M	-1	VAL	-	linker	UNP P42212
M	0	ASP	-	linker	UNP P42212
M	287	ALA	THR	engineered mutation	UNP P08413
M	306	ALA	THR	engineered mutation	UNP P08413
M	307	ALA	THR	engineered mutation	UNP P08413
N	-272	MET	-	initiating methionine	UNP P42212
N	-271	GLY	-	expression tag	UNP P42212
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

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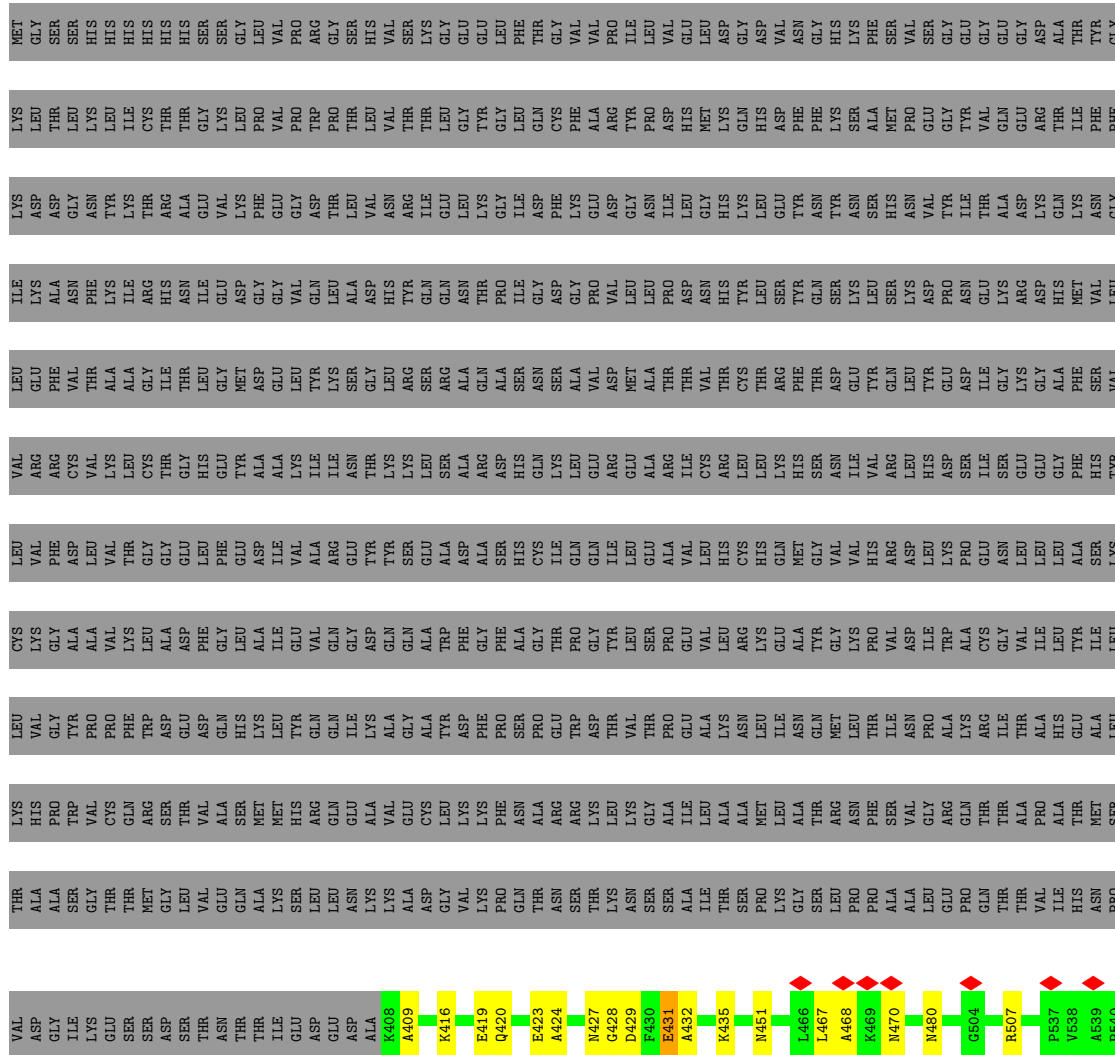
Chain	Residue	Modelled	Actual	Comment	Reference
N	-185	LEU	VAL	conflict	UNP P42212
N	-181	ALA	SER	conflict	UNP P42212
N	-100	THR	MET	conflict	UNP P42212
N	-90	ALA	VAL	conflict	UNP P42212
N	-78	GLY	SER	conflict	UNP P42212
N	-50	TYR	THR	conflict	UNP P42212
N	-47	LYS	ALA	conflict	UNP P42212
N	-22	LEU	HIS	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
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

- Molecule 1: Venus-tagged CaMKII Beta Holoenzyme mutant

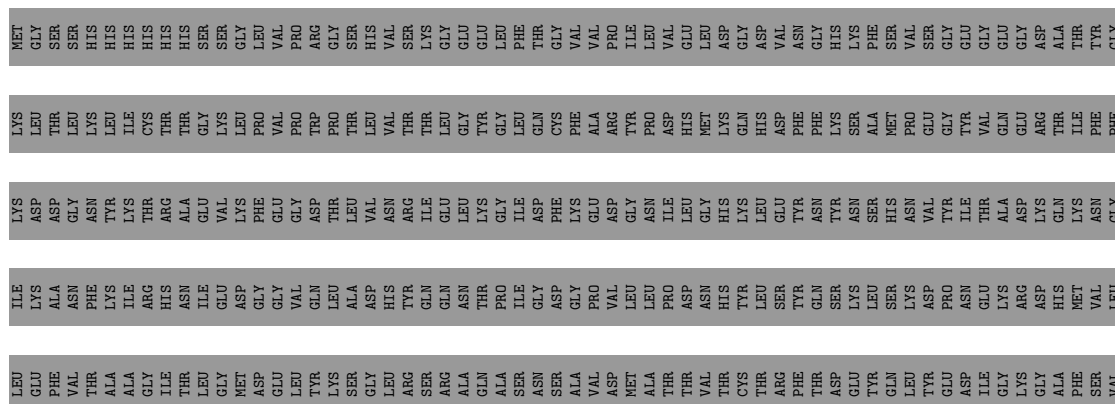


VAL	ASP	GLY	ALA	THR	LYS	LEU	VAL	CYS	LEU	VAL	LEU	VAL	ARG	VAL	LEU	THR	MET
ASP	GLY	THR	ALA	ALA	HIS	VAL	GLY	LYS	PHE	ARG	GLY	ARG	ARG	GLY	THR	GLY	GLY
ILE	ILE	SER	SER	SER	TRP	TRP	ASP	ALA	ASP	CYS	VAL	VAL	CYS	ASP	VAL	THR	SER
LYS	GLY	THR	GLY	THR	VAL	VAL	LEU	ALA	LEU	LYS	ALA	ALA	LYS	THR	ALA	LYS	HIS
THR	SER	THR	THR	THR	CYS	GLN	THR	LYS	THR	LYS	THR	THR	LYS	THR	ALA	HIS	HIS
SER	SER	MET	GLY	MET	ARG	THR	ASP	LEU	GLY	CYS	THR	THR	CYS	GLY	CYS	HIS	HIS
ASP	ASP	GLY	THR	GLY	SER	SER	ASP	ALA	GLY	THR	THR	THR	THR	THR	THR	HIS	HIS
SER	SER	LEU	LEU	LEU	THR	THR	GLU	PHE	GLY	GLY	GLY	GLY	HIS	GLY	THR	HIS	HIS
THR	THR	VAL	VAL	VAL	VAL	GLN	ASP	ALA	LEU	HIS	LEU	HIS	GLY	VAL	GLY	SER	SER
THR	THR	GLN	GLN	GLN	ALA	ALA	THR	LEU	GLY	THR	GLY	THR	THR	GLY	GLY	LEU	LEU
THR	THR	ALA	ALA	ALA	SER	THR	ASP	ALA	ASP	THR	ASP	ALA	THR	GLY	PRO	LEU	LEU
THR	THR	ALA	ALA	ALA	MET	LYS	LEU	LYS	LEU	LYS	LEU	LEU	LEU	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	MET	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	THR	THR	GLU	PHE	GLY	GLY	GLY	GLY	HIS	GLY	THR	HIS	HIS
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
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THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
THR	THR	ALA	ALA	ALA	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
THR	THR	ALA	ALA	ALA	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
THR	THR	ALA	ALA	ALA	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS	LYS
THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN
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THR	THR	ALA	ALA	ALA	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
THR	THR	ALA	ALA	ALA	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
THR	THR	ALA	ALA	ALA	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU	LEU
THR	THR	ALA	ALA	ALA	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GLN	GL				

Chain E: 14% 83%



Chain F: 14% 83%





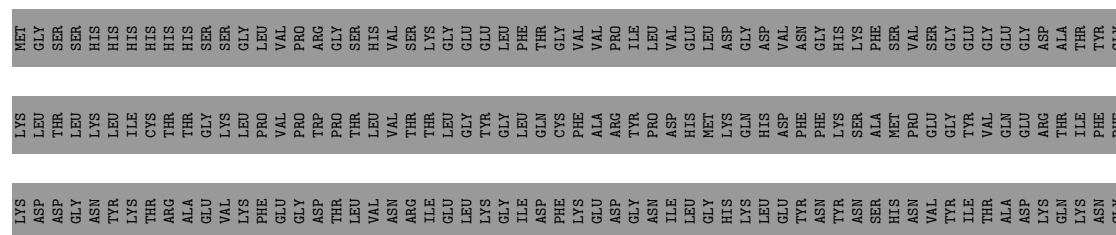
• Molecule 1: Venus-tagged CaMKII Beta Holoenzyme mutant

Chain H: 14% 83%



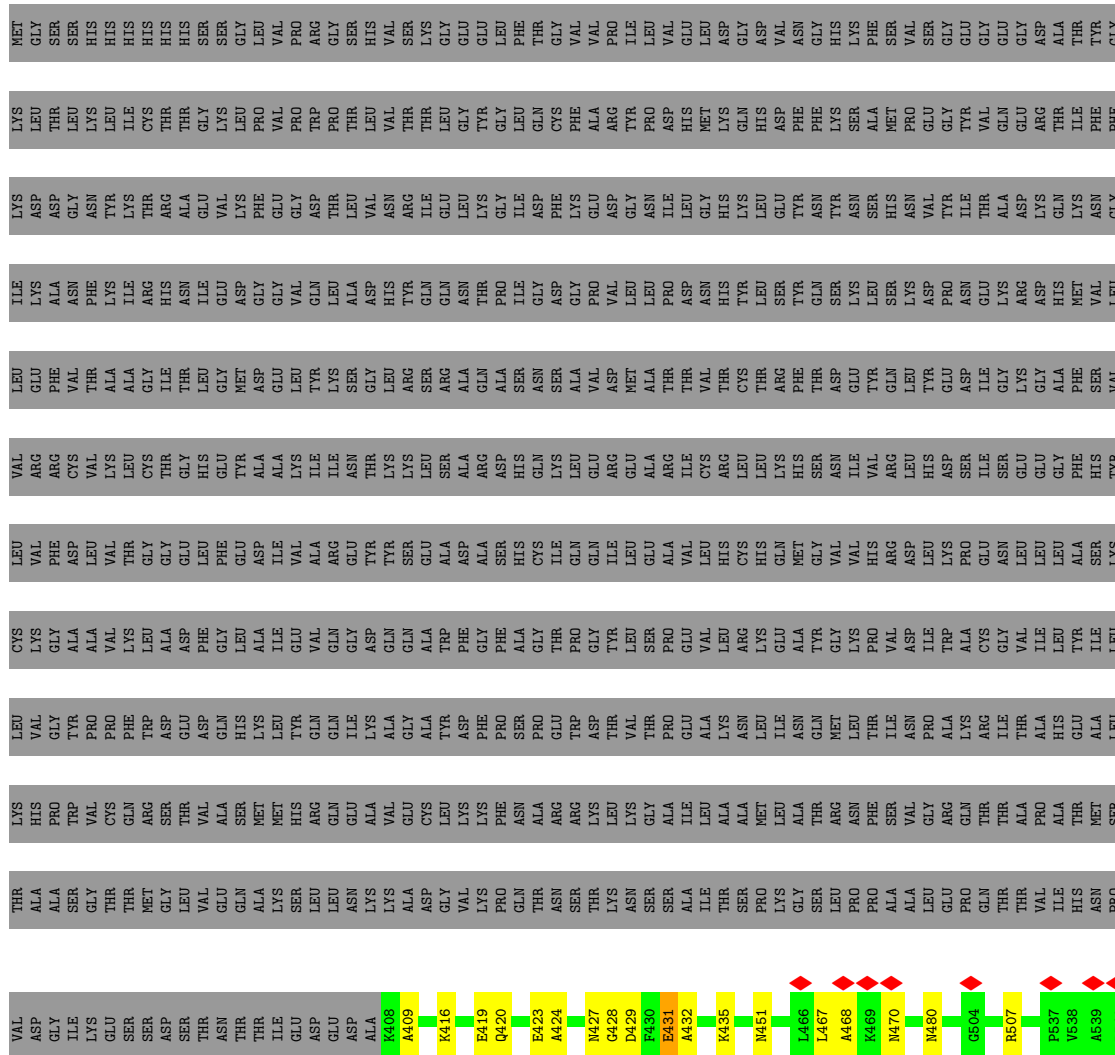
• Molecule 1: Venus-tagged CaMKII Beta Holoenzyme mutant

Chain I: 14% 83%



[illegible]

- Molecule 1: Venus-tagged CaMKII Beta Holoenzyme mutant



- Molecule 1: Venus-tagged CaMKII Beta Holoenzyme mutant



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, D7	Depositor
Number of particles used	138904	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	1.534	Depositor
Minimum map value	-1.191	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.030	Depositor
Recommended contour level	0.1	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	1.39	7/1107 (0.6%)	1.72	19/1501 (1.3%)
1	B	1.39	7/1107 (0.6%)	1.72	19/1501 (1.3%)
1	C	1.39	7/1107 (0.6%)	1.72	20/1501 (1.3%)
1	D	1.39	7/1107 (0.6%)	1.72	20/1501 (1.3%)
1	E	1.39	7/1107 (0.6%)	1.72	20/1501 (1.3%)
1	F	1.39	7/1107 (0.6%)	1.72	19/1501 (1.3%)
1	G	1.39	7/1107 (0.6%)	1.72	20/1501 (1.3%)
1	H	1.39	7/1107 (0.6%)	1.72	19/1501 (1.3%)
1	I	1.39	7/1107 (0.6%)	1.72	19/1501 (1.3%)
1	J	1.39	7/1107 (0.6%)	1.72	19/1501 (1.3%)
1	K	1.39	7/1107 (0.6%)	1.72	19/1501 (1.3%)
1	L	1.39	7/1107 (0.6%)	1.72	19/1501 (1.3%)
1	M	1.39	7/1107 (0.6%)	1.72	19/1501 (1.3%)
1	N	1.39	7/1107 (0.6%)	1.72	20/1501 (1.3%)
All	All	1.39	98/15498 (0.6%)	1.72	271/21014 (1.3%)

All (98) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	480	ASN	CA-CB	-8.99	1.46	1.52
1	I	480	ASN	CA-CB	-8.96	1.46	1.52
1	N	480	ASN	CA-CB	-8.95	1.46	1.52
1	C	480	ASN	CA-CB	-8.94	1.46	1.52
1	K	480	ASN	CA-CB	-8.94	1.46	1.52
1	A	480	ASN	CA-CB	-8.94	1.46	1.52
1	G	480	ASN	CA-CB	-8.94	1.46	1.52
1	H	480	ASN	CA-CB	-8.94	1.46	1.52
1	J	480	ASN	CA-CB	-8.92	1.46	1.52
1	E	480	ASN	CA-CB	-8.90	1.46	1.52
1	L	480	ASN	CA-CB	-8.90	1.46	1.52
1	F	480	ASN	CA-CB	-8.88	1.46	1.52
1	M	480	ASN	CA-CB	-8.88	1.46	1.52
1	B	480	ASN	CA-CB	-8.87	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	424	ALA	CA-CB	-7.11	1.42	1.53
1	K	424	ALA	CA-CB	-7.10	1.42	1.53
1	J	424	ALA	CA-CB	-7.09	1.42	1.53
1	C	424	ALA	CA-CB	-7.09	1.42	1.53
1	A	424	ALA	CA-CB	-7.08	1.42	1.53
1	H	424	ALA	CA-CB	-7.08	1.42	1.53
1	E	424	ALA	CA-CB	-7.07	1.42	1.53
1	L	424	ALA	CA-CB	-7.07	1.42	1.53
1	B	424	ALA	CA-CB	-7.07	1.42	1.53
1	I	424	ALA	CA-CB	-7.07	1.42	1.53
1	E	432	ALA	CA-CB	-7.06	1.42	1.53
1	F	424	ALA	CA-CB	-7.06	1.42	1.53
1	M	424	ALA	CA-CB	-7.06	1.42	1.53
1	D	432	ALA	CA-CB	-7.06	1.42	1.53
1	G	424	ALA	CA-CB	-7.05	1.42	1.53
1	N	424	ALA	CA-CB	-7.05	1.42	1.53
1	K	432	ALA	CA-CB	-7.05	1.42	1.53
1	A	432	ALA	CA-CB	-7.05	1.42	1.53
1	H	432	ALA	CA-CB	-7.05	1.42	1.53
1	G	432	ALA	CA-CB	-7.04	1.42	1.53
1	N	432	ALA	CA-CB	-7.04	1.42	1.53
1	F	432	ALA	CA-CB	-7.03	1.42	1.53
1	M	432	ALA	CA-CB	-7.03	1.42	1.53
1	C	432	ALA	CA-CB	-7.03	1.42	1.53
1	L	432	ALA	CA-CB	-7.02	1.42	1.53
1	B	432	ALA	CA-CB	-7.00	1.42	1.53
1	I	432	ALA	CA-CB	-7.00	1.42	1.53
1	J	432	ALA	CA-CB	-6.99	1.42	1.53
1	F	468	ALA	CA-CB	-6.89	1.42	1.53
1	L	468	ALA	CA-CB	-6.88	1.42	1.53
1	D	468	ALA	CA-CB	-6.87	1.42	1.53
1	N	468	ALA	CA-CB	-6.87	1.42	1.53
1	M	468	ALA	CA-CB	-6.87	1.42	1.53
1	J	468	ALA	CA-CB	-6.86	1.42	1.53
1	A	468	ALA	CA-CB	-6.84	1.42	1.53
1	H	468	ALA	CA-CB	-6.84	1.42	1.53
1	B	468	ALA	CA-CB	-6.83	1.42	1.53
1	I	468	ALA	CA-CB	-6.83	1.42	1.53
1	E	468	ALA	CA-CB	-6.82	1.42	1.53
1	C	468	ALA	CA-CB	-6.81	1.42	1.53
1	K	468	ALA	CA-CB	-6.81	1.42	1.53
1	G	468	ALA	CA-CB	-6.81	1.42	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	K	480	ASN	N-CA	-6.52	1.41	1.46
1	E	409	ALA	CA-CB	-6.51	1.42	1.53
1	I	409	ALA	CA-CB	-6.51	1.42	1.53
1	C	480	ASN	N-CA	-6.50	1.41	1.46
1	F	480	ASN	N-CA	-6.49	1.41	1.46
1	G	409	ALA	CA-CB	-6.49	1.42	1.53
1	E	480	ASN	N-CA	-6.48	1.41	1.46
1	L	480	ASN	N-CA	-6.48	1.41	1.46
1	A	480	ASN	N-CA	-6.48	1.41	1.46
1	H	480	ASN	N-CA	-6.48	1.41	1.46
1	J	409	ALA	CA-CB	-6.48	1.42	1.53
1	L	409	ALA	CA-CB	-6.47	1.42	1.53
1	B	480	ASN	N-CA	-6.47	1.41	1.46
1	G	480	ASN	N-CA	-6.47	1.41	1.46
1	I	480	ASN	N-CA	-6.47	1.41	1.46
1	N	480	ASN	N-CA	-6.47	1.41	1.46
1	A	409	ALA	CA-CB	-6.47	1.42	1.53
1	C	409	ALA	CA-CB	-6.47	1.42	1.53
1	F	409	ALA	CA-CB	-6.47	1.42	1.53
1	H	409	ALA	CA-CB	-6.47	1.42	1.53
1	M	409	ALA	CA-CB	-6.47	1.42	1.53
1	N	409	ALA	CA-CB	-6.46	1.42	1.53
1	J	480	ASN	N-CA	-6.46	1.41	1.46
1	B	409	ALA	CA-CB	-6.46	1.42	1.53
1	D	409	ALA	CA-CB	-6.46	1.42	1.53
1	K	409	ALA	CA-CB	-6.46	1.42	1.53
1	D	480	ASN	N-CA	-6.45	1.41	1.46
1	M	480	ASN	N-CA	-6.45	1.41	1.46
1	J	428	GLY	N-CA	-5.59	1.37	1.45
1	E	428	GLY	N-CA	-5.57	1.37	1.45
1	L	428	GLY	N-CA	-5.57	1.37	1.45
1	D	428	GLY	N-CA	-5.56	1.37	1.45
1	K	428	GLY	N-CA	-5.56	1.37	1.45
1	A	428	GLY	N-CA	-5.55	1.37	1.45
1	H	428	GLY	N-CA	-5.55	1.37	1.45
1	G	428	GLY	N-CA	-5.54	1.37	1.45
1	N	428	GLY	N-CA	-5.54	1.37	1.45
1	F	428	GLY	N-CA	-5.54	1.37	1.45
1	M	428	GLY	N-CA	-5.54	1.37	1.45
1	C	428	GLY	N-CA	-5.53	1.37	1.45
1	I	428	GLY	N-CA	-5.50	1.38	1.45
1	B	428	GLY	N-CA	-5.50	1.38	1.45

All (271) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	427	ASN	CA-CB-CG	7.70	120.30	112.60
1	C	427	ASN	CA-CB-CG	7.70	120.30	112.60
1	L	427	ASN	CA-CB-CG	7.69	120.29	112.60
1	B	427	ASN	CA-CB-CG	7.67	120.27	112.60
1	D	427	ASN	CA-CB-CG	7.67	120.27	112.60
1	K	427	ASN	CA-CB-CG	7.66	120.26	112.60
1	I	427	ASN	CA-CB-CG	7.65	120.25	112.60
1	N	427	ASN	CA-CB-CG	7.65	120.25	112.60
1	A	427	ASN	CA-CB-CG	7.65	120.25	112.60
1	H	427	ASN	CA-CB-CG	7.65	120.25	112.60
1	G	427	ASN	CA-CB-CG	7.65	120.25	112.60
1	M	427	ASN	CA-CB-CG	7.65	120.25	112.60
1	E	427	ASN	CA-CB-CG	7.64	120.24	112.60
1	J	427	ASN	CA-CB-CG	7.62	120.22	112.60
1	J	429	ASP	CA-CB-CG	6.89	119.50	112.60
1	B	429	ASP	CA-CB-CG	6.89	119.49	112.60
1	C	429	ASP	CA-CB-CG	6.89	119.49	112.60
1	G	429	ASP	CA-CB-CG	6.89	119.49	112.60
1	N	429	ASP	CA-CB-CG	6.89	119.49	112.60
1	I	429	ASP	CA-CB-CG	6.88	119.48	112.60
1	F	429	ASP	CA-CB-CG	6.87	119.47	112.60
1	M	429	ASP	CA-CB-CG	6.87	119.47	112.60
1	A	429	ASP	CA-CB-CG	6.86	119.46	112.60
1	H	429	ASP	CA-CB-CG	6.86	119.46	112.60
1	N	470	ASN	CA-CB-CG	6.86	119.46	112.60
1	D	429	ASP	CA-CB-CG	6.85	119.45	112.60
1	K	429	ASP	CA-CB-CG	6.85	119.45	112.60
1	F	470	ASN	CA-CB-CG	6.84	119.44	112.60
1	M	470	ASN	CA-CB-CG	6.84	119.44	112.60
1	E	429	ASP	CA-CB-CG	6.84	119.44	112.60
1	D	470	ASN	CA-CB-CG	6.83	119.44	112.60
1	K	470	ASN	CA-CB-CG	6.83	119.44	112.60
1	B	470	ASN	CA-CB-CG	6.83	119.43	112.60
1	G	470	ASN	CA-CB-CG	6.83	119.43	112.60
1	A	470	ASN	CA-CB-CG	6.82	119.42	112.60
1	C	470	ASN	CA-CB-CG	6.82	119.42	112.60
1	H	470	ASN	CA-CB-CG	6.82	119.42	112.60
1	J	470	ASN	CA-CB-CG	6.82	119.42	112.60
1	L	429	ASP	CA-CB-CG	6.82	119.42	112.60
1	E	470	ASN	CA-CB-CG	6.81	119.41	112.60
1	L	470	ASN	CA-CB-CG	6.81	119.41	112.60
1	I	470	ASN	CA-CB-CG	6.80	119.40	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	451	ASN	CA-CB-CG	6.55	119.15	112.60
1	N	451	ASN	CA-CB-CG	6.55	119.15	112.60
1	E	451	ASN	CA-CB-CG	6.54	119.14	112.60
1	F	451	ASN	CA-CB-CG	6.54	119.14	112.60
1	M	451	ASN	CA-CB-CG	6.54	119.14	112.60
1	C	451	ASN	CA-CB-CG	6.54	119.14	112.60
1	A	451	ASN	CA-CB-CG	6.53	119.13	112.60
1	H	451	ASN	CA-CB-CG	6.53	119.13	112.60
1	I	451	ASN	CA-CB-CG	6.53	119.13	112.60
1	D	451	ASN	CA-CB-CG	6.53	119.13	112.60
1	K	451	ASN	CA-CB-CG	6.53	119.13	112.60
1	L	451	ASN	CA-CB-CG	6.53	119.13	112.60
1	J	451	ASN	CA-CB-CG	6.51	119.11	112.60
1	B	451	ASN	CA-CB-CG	6.49	119.09	112.60
1	B	480	ASN	CA-CB-CG	6.24	118.83	112.60
1	F	480	ASN	CA-CB-CG	6.23	118.83	112.60
1	G	480	ASN	CA-CB-CG	6.23	118.83	112.60
1	I	480	ASN	CA-CB-CG	6.23	118.83	112.60
1	M	480	ASN	CA-CB-CG	6.23	118.83	112.60
1	C	480	ASN	CA-CB-CG	6.21	118.81	112.60
1	L	480	ASN	CA-CB-CG	6.21	118.81	112.60
1	K	480	ASN	CA-CB-CG	6.21	118.81	112.60
1	N	480	ASN	CA-CB-CG	6.21	118.81	112.60
1	A	480	ASN	CA-CB-CG	6.21	118.81	112.60
1	H	480	ASN	CA-CB-CG	6.21	118.81	112.60
1	E	480	ASN	CA-CB-CG	6.20	118.80	112.60
1	J	480	ASN	CA-CB-CG	6.20	118.80	112.60
1	D	480	ASN	CA-CB-CG	6.18	118.78	112.60
1	J	431	GLU	CA-C-N	5.35	127.40	120.44
1	J	431	GLU	C-N-CA	5.35	127.40	120.44
1	D	431	GLU	CA-C-N	5.35	127.39	120.44
1	D	431	GLU	C-N-CA	5.35	127.39	120.44
1	F	431	GLU	CA-C-N	5.33	127.38	120.44
1	F	431	GLU	C-N-CA	5.33	127.38	120.44
1	M	431	GLU	CA-C-N	5.33	127.38	120.44
1	M	431	GLU	C-N-CA	5.33	127.38	120.44
1	B	431	GLU	CA-C-N	5.33	127.36	120.44
1	B	431	GLU	C-N-CA	5.33	127.36	120.44
1	A	431	GLU	CA-C-N	5.32	127.36	120.44
1	A	431	GLU	C-N-CA	5.32	127.36	120.44
1	H	431	GLU	CA-C-N	5.32	127.36	120.44
1	H	431	GLU	C-N-CA	5.32	127.36	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	431	GLU	CA-C-N	5.32	127.36	120.44
1	G	431	GLU	C-N-CA	5.32	127.36	120.44
1	N	431	GLU	CA-C-N	5.32	127.35	120.44
1	N	431	GLU	C-N-CA	5.32	127.35	120.44
1	E	431	GLU	CA-C-N	5.31	127.34	120.44
1	E	431	GLU	C-N-CA	5.31	127.34	120.44
1	I	431	GLU	CA-C-N	5.31	127.34	120.44
1	I	431	GLU	C-N-CA	5.31	127.34	120.44
1	L	431	GLU	CA-C-N	5.30	127.33	120.44
1	L	431	GLU	C-N-CA	5.30	127.33	120.44
1	C	431	GLU	CA-C-N	5.29	127.32	120.44
1	C	431	GLU	C-N-CA	5.29	127.32	120.44
1	K	431	GLU	CA-C-N	5.29	127.31	120.44
1	K	431	GLU	C-N-CA	5.29	127.31	120.44
1	D	419	GLU	CA-C-N	5.25	127.26	120.44
1	D	419	GLU	C-N-CA	5.25	127.26	120.44
1	I	419	GLU	CA-C-N	5.24	127.26	120.44
1	I	419	GLU	C-N-CA	5.24	127.26	120.44
1	M	416	LYS	N-CA-CB	5.23	117.60	110.01
1	E	416	LYS	N-CA-CB	5.23	117.59	110.01
1	G	419	GLU	CA-C-N	5.23	127.24	120.44
1	G	419	GLU	C-N-CA	5.23	127.24	120.44
1	L	416	LYS	N-CA-CB	5.23	117.59	110.01
1	N	419	GLU	CA-C-N	5.23	127.24	120.44
1	N	419	GLU	C-N-CA	5.23	127.24	120.44
1	L	419	GLU	CA-C-N	5.23	127.24	120.44
1	L	419	GLU	C-N-CA	5.23	127.24	120.44
1	B	419	GLU	CA-C-N	5.23	127.24	120.44
1	B	419	GLU	C-N-CA	5.23	127.24	120.44
1	J	419	GLU	CA-C-N	5.23	127.24	120.44
1	J	419	GLU	C-N-CA	5.23	127.24	120.44
1	A	419	GLU	CA-C-N	5.22	127.23	120.44
1	A	419	GLU	C-N-CA	5.22	127.23	120.44
1	B	416	LYS	N-CA-CB	5.22	117.58	110.01
1	G	416	LYS	N-CA-CB	5.22	117.59	110.01
1	H	419	GLU	CA-C-N	5.22	127.23	120.44
1	H	419	GLU	C-N-CA	5.22	127.23	120.44
1	I	416	LYS	N-CA-CB	5.22	117.58	110.01
1	N	416	LYS	N-CA-CB	5.22	117.59	110.01
1	E	419	GLU	CA-C-N	5.22	127.23	120.44
1	E	419	GLU	C-N-CA	5.22	127.23	120.44
1	K	419	GLU	CA-C-N	5.22	127.22	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	419	GLU	C-N-CA	5.22	127.22	120.44
1	A	416	LYS	N-CA-CB	5.21	117.57	110.01
1	H	416	LYS	N-CA-CB	5.21	117.57	110.01
1	J	423	GLU	CA-C-N	5.21	127.22	120.44
1	J	423	GLU	C-N-CA	5.21	127.22	120.44
1	F	419	GLU	CA-C-N	5.21	127.21	120.44
1	F	419	GLU	C-N-CA	5.21	127.21	120.44
1	M	419	GLU	CA-C-N	5.21	127.21	120.44
1	M	419	GLU	C-N-CA	5.21	127.21	120.44
1	C	419	GLU	CA-C-N	5.21	127.21	120.44
1	C	419	GLU	C-N-CA	5.21	127.21	120.44
1	K	416	LYS	N-CA-CB	5.21	117.56	110.01
1	A	423	GLU	CA-C-N	5.21	127.21	120.44
1	A	423	GLU	C-N-CA	5.21	127.21	120.44
1	H	423	GLU	CA-C-N	5.21	127.21	120.44
1	H	423	GLU	C-N-CA	5.21	127.21	120.44
1	K	423	GLU	CA-C-N	5.21	127.21	120.44
1	K	423	GLU	C-N-CA	5.21	127.21	120.44
1	C	416	LYS	N-CA-CB	5.21	117.56	110.01
1	J	416	LYS	N-CA-CB	5.21	117.56	110.01
1	C	423	GLU	CA-C-N	5.20	127.20	120.44
1	C	423	GLU	C-N-CA	5.20	127.20	120.44
1	I	423	GLU	CA-C-N	5.20	127.20	120.44
1	I	423	GLU	C-N-CA	5.20	127.20	120.44
1	D	416	LYS	N-CA-CB	5.20	117.55	110.01
1	E	423	GLU	CA-C-N	5.20	127.20	120.44
1	E	423	GLU	C-N-CA	5.20	127.20	120.44
1	L	423	GLU	CA-C-N	5.20	127.20	120.44
1	L	423	GLU	C-N-CA	5.20	127.20	120.44
1	F	416	LYS	N-CA-CB	5.20	117.55	110.01
1	F	423	GLU	CA-C-N	5.20	127.20	120.44
1	F	423	GLU	C-N-CA	5.20	127.20	120.44
1	M	423	GLU	CA-C-N	5.20	127.20	120.44
1	M	423	GLU	C-N-CA	5.20	127.20	120.44
1	D	423	GLU	CA-C-N	5.20	127.19	120.44
1	D	423	GLU	C-N-CA	5.20	127.19	120.44
1	G	423	GLU	CA-C-N	5.19	127.19	120.44
1	G	423	GLU	C-N-CA	5.19	127.19	120.44
1	N	423	GLU	CA-C-N	5.19	127.19	120.44
1	N	423	GLU	C-N-CA	5.19	127.19	120.44
1	B	423	GLU	CA-C-N	5.18	127.17	120.44
1	B	423	GLU	C-N-CA	5.18	127.17	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	420	GLN	CA-CB-CG	5.17	124.44	114.10
1	L	420	GLN	CA-CB-CG	5.17	124.44	114.10
1	N	420	GLN	CA-CB-CG	5.17	124.44	114.10
1	I	420	GLN	CA-CB-CG	5.16	124.42	114.10
1	A	420	GLN	CA-CB-CG	5.16	124.41	114.10
1	D	420	GLN	CA-CB-CG	5.16	124.41	114.10
1	H	420	GLN	CA-CB-CG	5.16	124.41	114.10
1	J	420	GLN	CA-CB-CG	5.15	124.41	114.10
1	E	420	GLN	CA-CB-CG	5.15	124.41	114.10
1	K	420	GLN	CA-CB-CG	5.15	124.41	114.10
1	C	420	GLN	CA-CB-CG	5.15	124.39	114.10
1	F	420	GLN	CA-CB-CG	5.14	124.39	114.10
1	E	467	LEU	CA-C-N	5.13	127.16	120.28
1	E	467	LEU	C-N-CA	5.13	127.16	120.28
1	L	467	LEU	CA-C-N	5.13	127.16	120.28
1	L	467	LEU	C-N-CA	5.13	127.16	120.28
1	I	428	GLY	CA-C-N	5.13	130.37	122.77
1	I	428	GLY	C-N-CA	5.13	130.37	122.77
1	I	467	LEU	CA-C-N	5.13	127.16	120.28
1	I	467	LEU	C-N-CA	5.13	127.16	120.28
1	K	428	GLY	CA-C-N	5.13	130.37	122.77
1	K	428	GLY	C-N-CA	5.13	130.37	122.77
1	B	420	GLN	CA-CB-CG	5.13	124.36	114.10
1	F	428	GLY	CA-C-N	5.13	130.37	122.77
1	F	428	GLY	C-N-CA	5.13	130.37	122.77
1	C	467	LEU	CA-C-N	5.13	127.15	120.28
1	C	467	LEU	C-N-CA	5.13	127.15	120.28
1	D	467	LEU	CA-C-N	5.13	127.15	120.28
1	D	467	LEU	C-N-CA	5.13	127.15	120.28
1	M	420	GLN	CA-CB-CG	5.13	124.36	114.10
1	B	467	LEU	CA-C-N	5.13	127.15	120.28
1	B	467	LEU	C-N-CA	5.13	127.15	120.28
1	F	467	LEU	CA-C-N	5.12	127.15	120.28
1	F	467	LEU	C-N-CA	5.12	127.15	120.28
1	N	428	GLY	CA-C-N	5.12	130.35	122.77
1	N	428	GLY	C-N-CA	5.12	130.35	122.77
1	B	428	GLY	CA-C-N	5.12	130.35	122.77
1	B	428	GLY	C-N-CA	5.12	130.35	122.77
1	A	467	LEU	CA-C-N	5.12	127.14	120.28
1	A	467	LEU	C-N-CA	5.12	127.14	120.28
1	H	467	LEU	CA-C-N	5.12	127.14	120.28
1	H	467	LEU	C-N-CA	5.12	127.14	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	428	GLY	CA-C-N	5.12	130.34	122.77
1	J	428	GLY	C-N-CA	5.12	130.34	122.77
1	C	428	GLY	CA-C-N	5.11	130.34	122.77
1	C	428	GLY	C-N-CA	5.11	130.34	122.77
1	E	428	GLY	CA-C-N	5.11	130.33	122.77
1	E	428	GLY	C-N-CA	5.11	130.33	122.77
1	G	428	GLY	CA-C-N	5.11	130.33	122.77
1	G	428	GLY	C-N-CA	5.11	130.33	122.77
1	A	428	GLY	CA-C-N	5.11	130.33	122.77
1	A	428	GLY	C-N-CA	5.11	130.33	122.77
1	G	467	LEU	CA-C-N	5.11	127.12	120.28
1	G	467	LEU	C-N-CA	5.11	127.12	120.28
1	H	428	GLY	CA-C-N	5.11	130.33	122.77
1	H	428	GLY	C-N-CA	5.11	130.33	122.77
1	M	467	LEU	CA-C-N	5.11	127.12	120.28
1	M	467	LEU	C-N-CA	5.11	127.12	120.28
1	N	467	LEU	CA-C-N	5.11	127.12	120.28
1	N	467	LEU	C-N-CA	5.11	127.12	120.28
1	D	428	GLY	CA-C-N	5.11	130.33	122.77
1	D	428	GLY	C-N-CA	5.11	130.33	122.77
1	M	428	GLY	CA-C-N	5.10	130.32	122.77
1	M	428	GLY	C-N-CA	5.10	130.32	122.77
1	K	467	LEU	CA-C-N	5.10	127.11	120.28
1	K	467	LEU	C-N-CA	5.10	127.11	120.28
1	L	428	GLY	CA-C-N	5.10	130.31	122.77
1	L	428	GLY	C-N-CA	5.10	130.31	122.77
1	J	467	LEU	CA-C-N	5.09	127.10	120.28
1	J	467	LEU	C-N-CA	5.09	127.10	120.28
1	E	431	GLU	CA-CB-CG	5.06	124.22	114.10
1	K	431	GLU	CA-CB-CG	5.05	124.20	114.10
1	L	431	GLU	CA-CB-CG	5.05	124.19	114.10
1	J	431	GLU	CA-CB-CG	5.04	124.18	114.10
1	G	423	GLU	N-CA-CB	5.04	117.31	110.01
1	G	431	GLU	CA-CB-CG	5.04	124.17	114.10
1	N	431	GLU	CA-CB-CG	5.04	124.17	114.10
1	A	431	GLU	CA-CB-CG	5.04	124.17	114.10
1	H	431	GLU	CA-CB-CG	5.04	124.17	114.10
1	B	431	GLU	CA-CB-CG	5.03	124.17	114.10
1	F	431	GLU	CA-CB-CG	5.03	124.16	114.10
1	I	431	GLU	CA-CB-CG	5.03	124.16	114.10
1	M	431	GLU	CA-CB-CG	5.03	124.16	114.10
1	C	419	GLU	N-CA-CB	5.03	117.51	110.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	431	GLU	CA-CB-CG	5.03	124.16	114.10
1	L	423	GLU	N-CA-CB	5.03	117.30	110.01
1	D	431	GLU	CA-CB-CG	5.02	124.15	114.10
1	J	423	GLU	N-CA-CB	5.02	117.29	110.01
1	E	423	GLU	N-CA-CB	5.02	117.29	110.01
1	N	423	GLU	N-CA-CB	5.02	117.28	110.01
1	A	423	GLU	N-CA-CB	5.01	117.28	110.01
1	C	423	GLU	N-CA-CB	5.01	117.28	110.01
1	E	419	GLU	N-CA-CB	5.01	117.49	110.12
1	H	423	GLU	N-CA-CB	5.01	117.28	110.01
1	G	419	GLU	N-CA-CB	5.01	117.49	110.12
1	N	419	GLU	N-CA-CB	5.01	117.49	110.12
1	B	423	GLU	N-CA-CB	5.01	117.27	110.01
1	F	423	GLU	N-CA-CB	5.01	117.27	110.01
1	I	423	GLU	N-CA-CB	5.01	117.27	110.01
1	M	423	GLU	N-CA-CB	5.01	117.27	110.01
1	D	423	GLU	N-CA-CB	5.00	117.27	110.01
1	K	423	GLU	N-CA-CB	5.00	117.27	110.01
1	D	419	GLU	N-CA-CB	5.00	117.47	110.12

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	1044	1043	2	0
1	B	1079	1044	1043	2	0
1	C	1079	1044	1043	2	0
1	D	1079	1044	1043	2	0
1	E	1079	1044	1043	2	0
1	F	1079	1044	1043	2	0
1	G	1079	1044	1043	2	0
1	H	1079	1044	1043	2	0
1	I	1079	1044	1043	2	0
1	J	1079	1044	1043	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	K	1079	1044	1043	2	0
1	L	1079	1044	1043	2	0
1	M	1079	1044	1043	2	0
1	N	1079	1044	1043	2	0
All	All	15106	14616	14602	14	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 0.

All (14) 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:E:542:GLN:OE1	1:F:507:ARG:NH1	2.51	0.44
1:J:542:GLN:OE1	1:K:507:ARG:NH1	2.51	0.44
1:I:542:GLN:OE1	1:J:507:ARG:NH1	2.51	0.43
1:C:542:GLN:OE1	1:D:507:ARG:NH1	2.51	0.43
1:K:542:GLN:OE1	1:L:507:ARG:NH1	2.52	0.43
1:F:542:GLN:OE1	1:G:507:ARG:NH1	2.52	0.43
1:H:542:GLN:OE1	1:I:507:ARG:NH1	2.52	0.43
1:A:507:ARG:NH1	1:G:542:GLN:OE1	2.52	0.43
1:A:542:GLN:OE1	1:B:507:ARG:NH1	2.52	0.43
1:D:542:GLN:OE1	1:E:507:ARG:NH1	2.52	0.43
1:H:507:ARG:NH1	1:N:542:GLN:OE1	2.52	0.43
1:L:542:GLN:OE1	1:M:507:ARG:NH1	2.52	0.43
1:M:542:GLN:OE1	1:N:507:ARG:NH1	2.52	0.43
1:B:542:GLN:OE1	1:C:507:ARG:NH1	2.52	0.42

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%)	130 (98%)	3 (2%)	0	100	100
1	B	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	C	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	D	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	E	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	F	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	G	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	H	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	I	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	J	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	K	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	L	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	M	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
1	N	133/815 (16%)	130 (98%)	3 (2%)	0	100	100
All	All	1862/11410 (16%)	1820 (98%)	42 (2%)	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%)	113 (98%)	2 (2%)	53	78
1	B	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	C	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	D	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	E	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	F	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	G	115/689 (17%)	113 (98%)	2 (2%)	53	78

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	I	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	J	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	K	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	L	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	M	115/689 (17%)	113 (98%)	2 (2%)	53	78
1	N	115/689 (17%)	113 (98%)	2 (2%)	53	78
All	All	1610/9646 (17%)	1582 (98%)	28 (2%)	52	78

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	431	GLU
1	A	435	LYS
1	B	431	GLU
1	B	435	LYS
1	C	431	GLU
1	C	435	LYS
1	D	431	GLU
1	D	435	LYS
1	E	431	GLU
1	E	435	LYS
1	F	431	GLU
1	F	435	LYS
1	G	431	GLU
1	G	435	LYS
1	H	431	GLU
1	H	435	LYS
1	I	431	GLU
1	I	435	LYS
1	J	431	GLU
1	J	435	LYS
1	K	431	GLU
1	K	435	LYS
1	L	431	GLU
1	L	435	LYS
1	M	431	GLU
1	M	435	LYS
1	N	431	GLU
1	N	435	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	484	HIS
1	G	484	HIS
1	H	484	HIS

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 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.

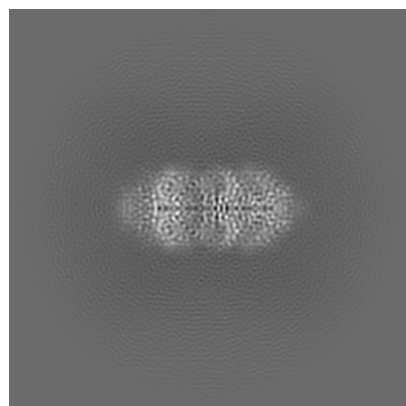
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-41070. These allow visual inspection of the internal detail of the map and identification of artifacts.

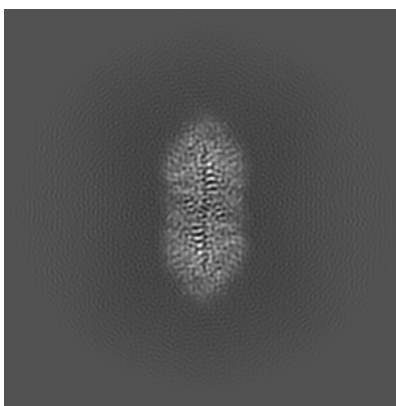
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

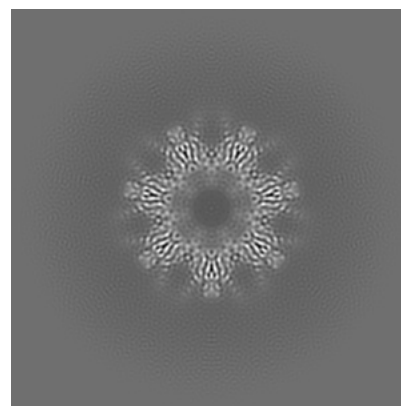
6.1.1 Primary map



X

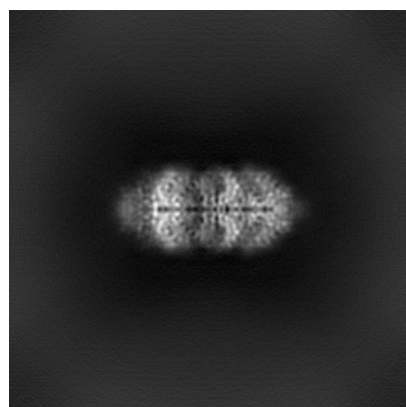


Y

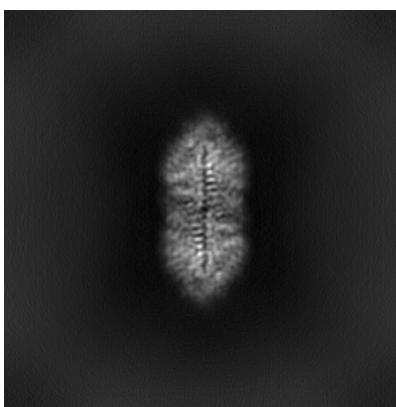


Z

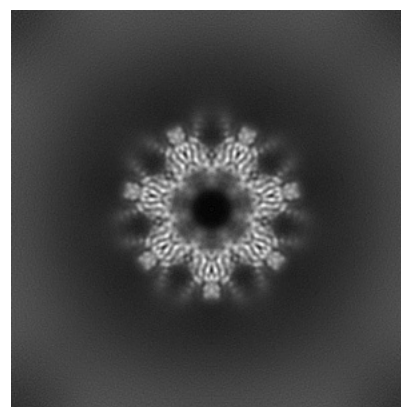
6.1.2 Raw map



X



Y

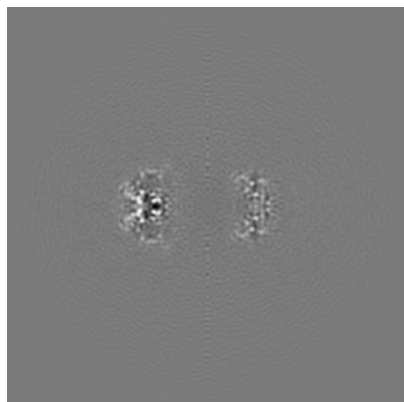


Z

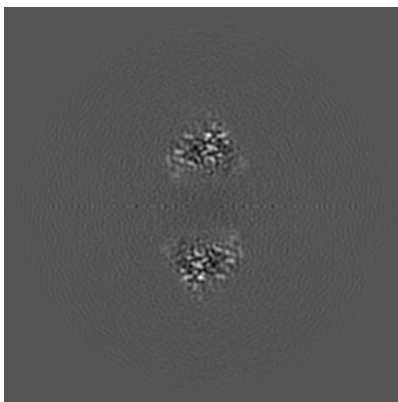
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

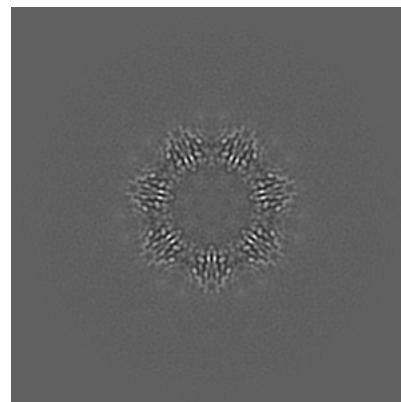
6.2.1 Primary map



X Index: 160

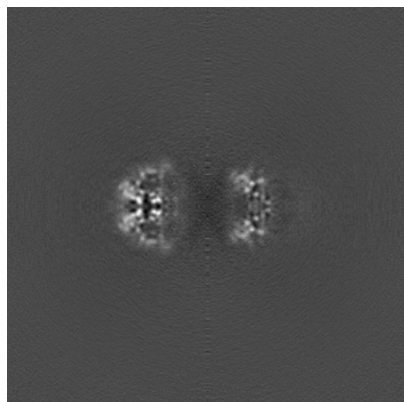


Y Index: 160

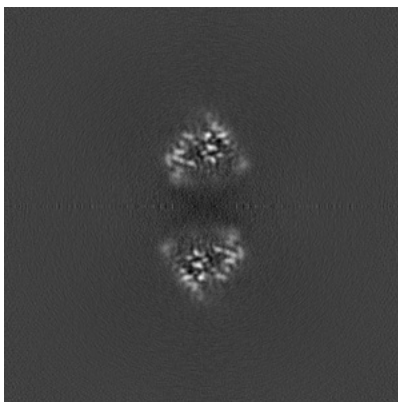


Z Index: 160

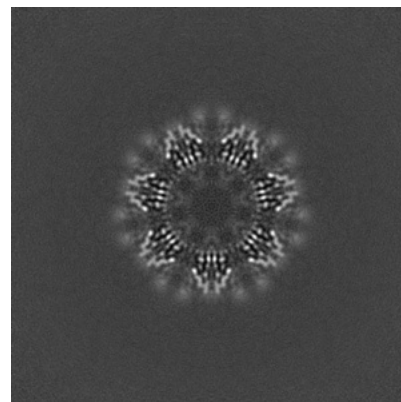
6.2.2 Raw map



X Index: 160



Y Index: 160

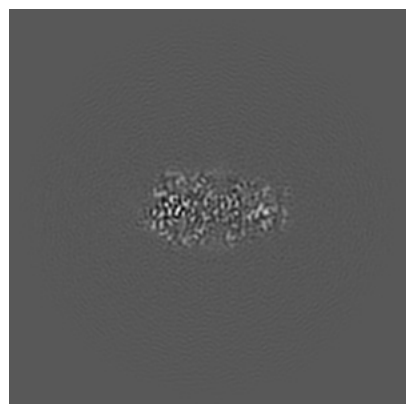


Z Index: 160

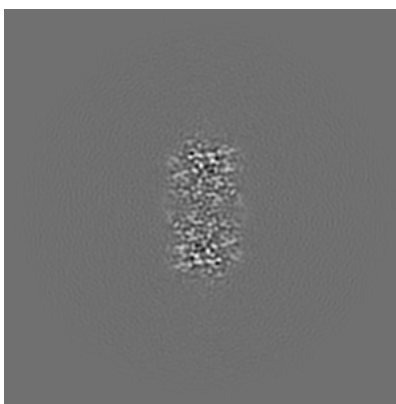
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

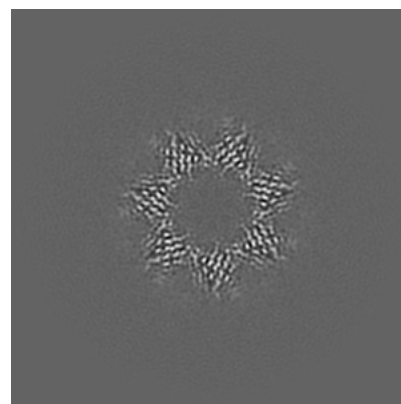
6.3.1 Primary map



X Index: 125

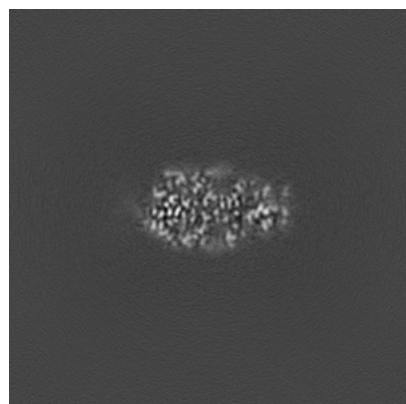


Y Index: 126

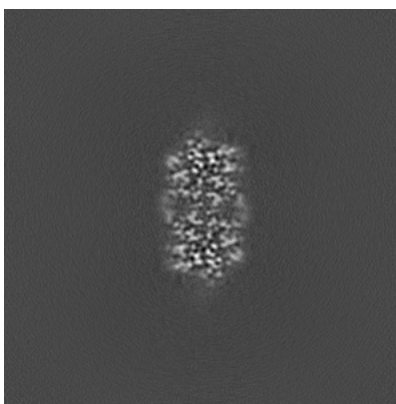


Z Index: 156

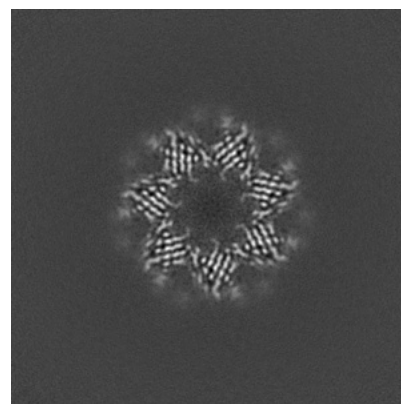
6.3.2 Raw map



X Index: 125



Y Index: 126

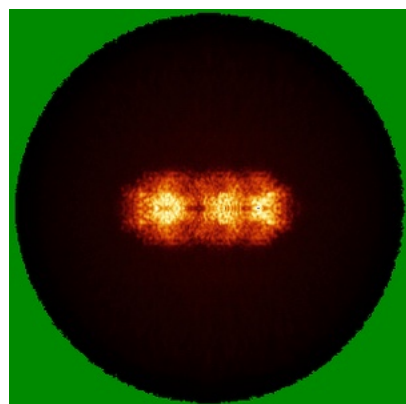


Z Index: 156

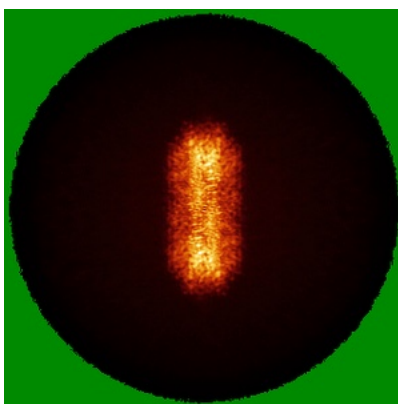
The images above show the largest variance slices of the map in three orthogonal directions.

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

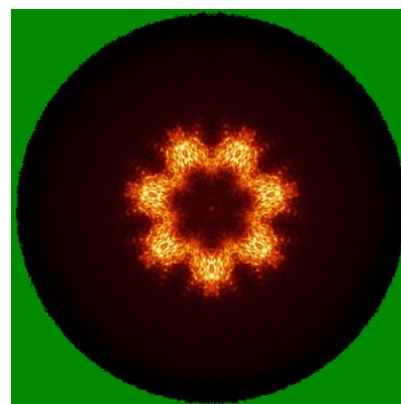
6.4.1 Primary map



X

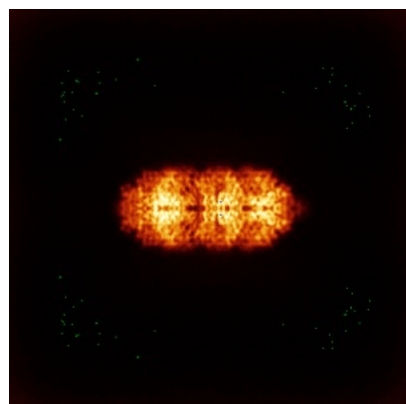


Y

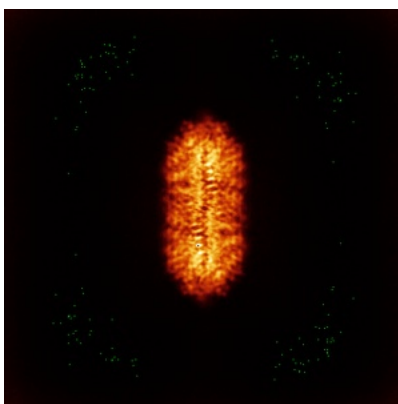


Z

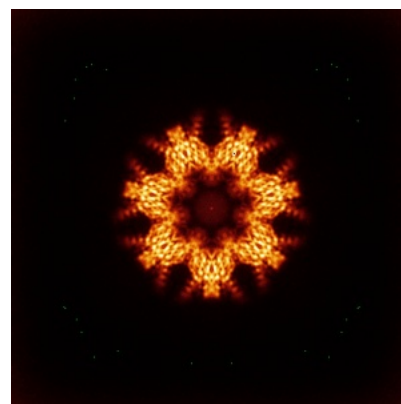
6.4.2 Raw map



X



Y

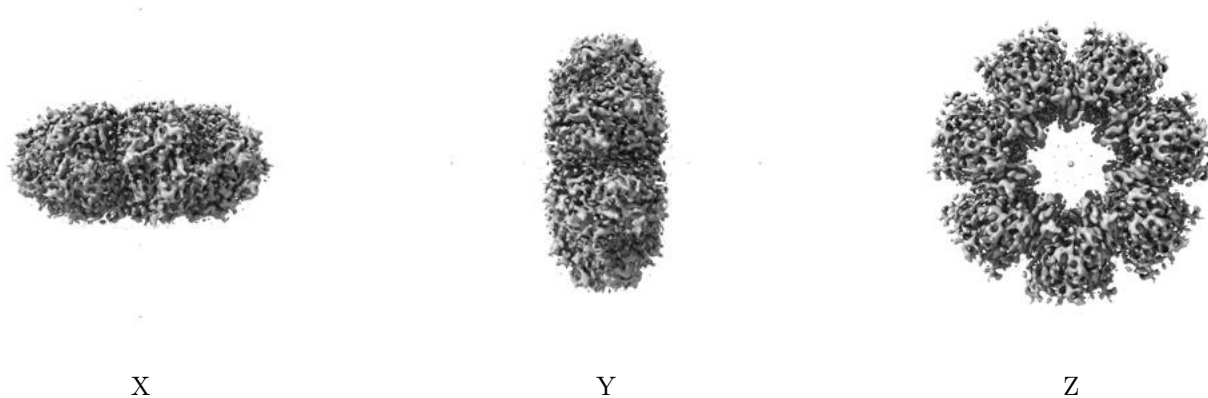


Z

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.

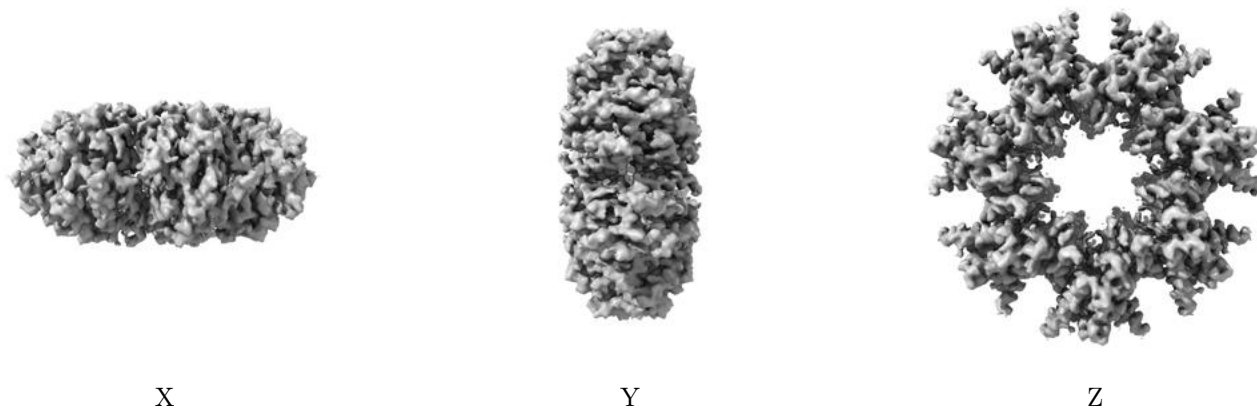
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

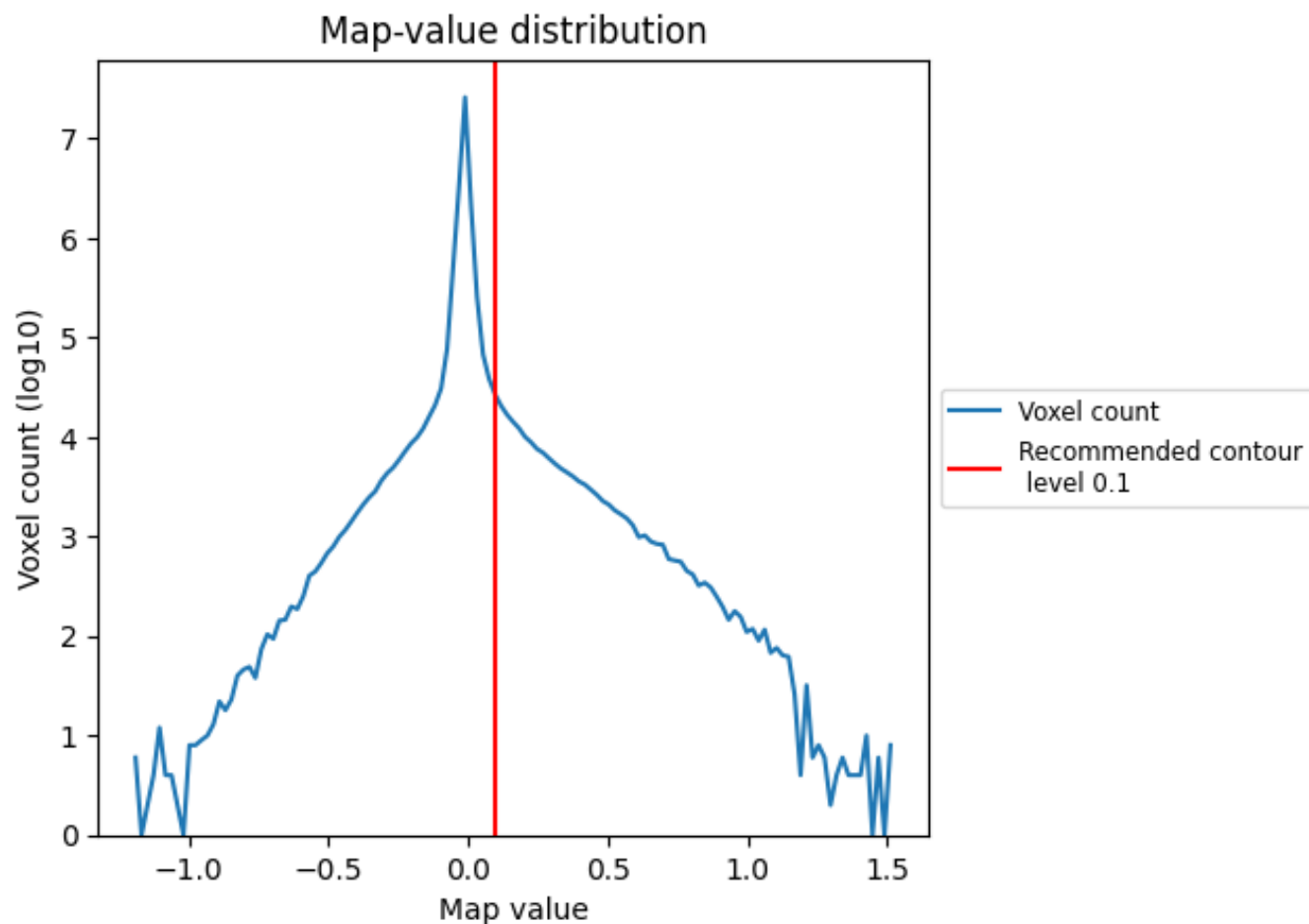
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

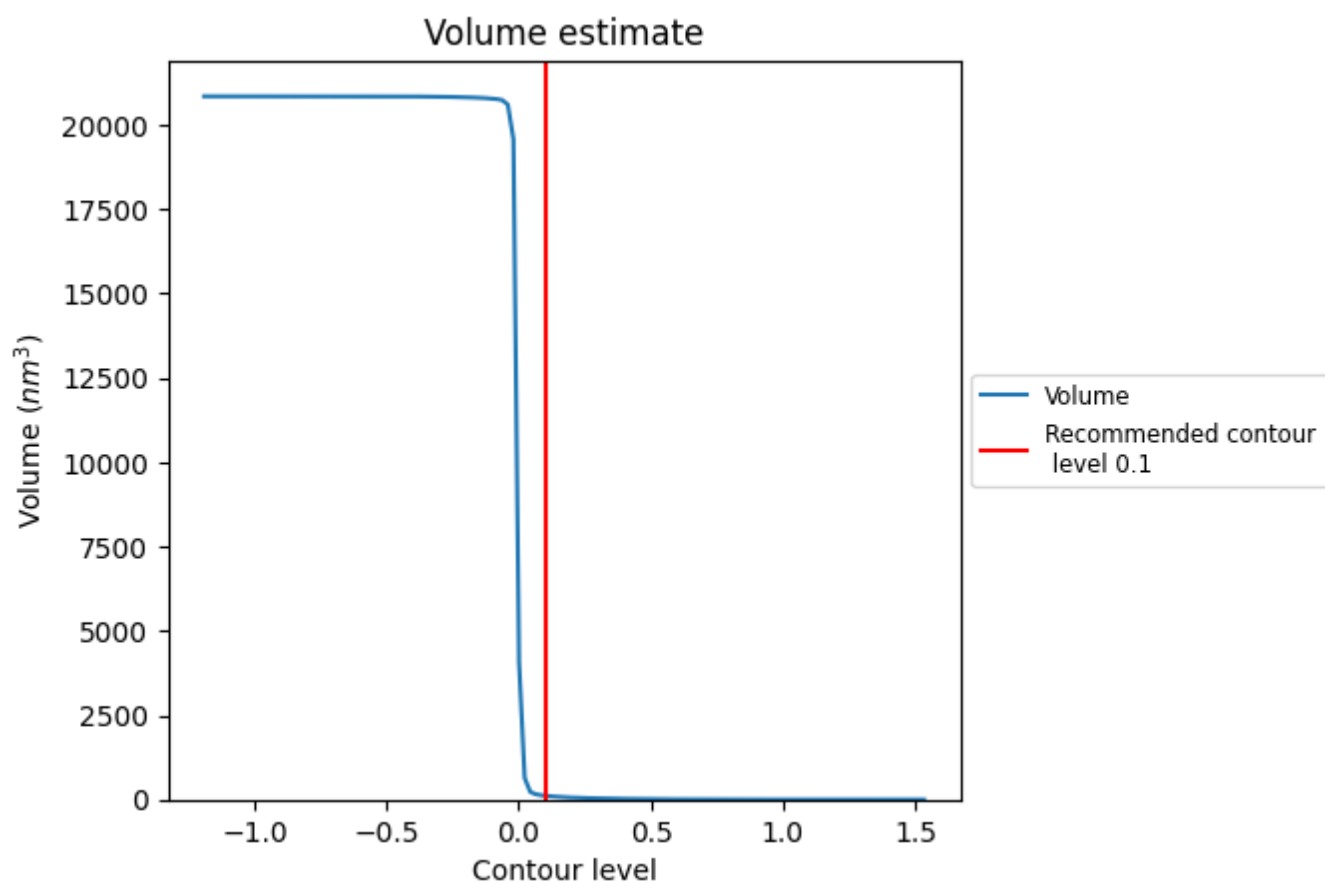
This section contains the results of statistical analysis of the map.

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.

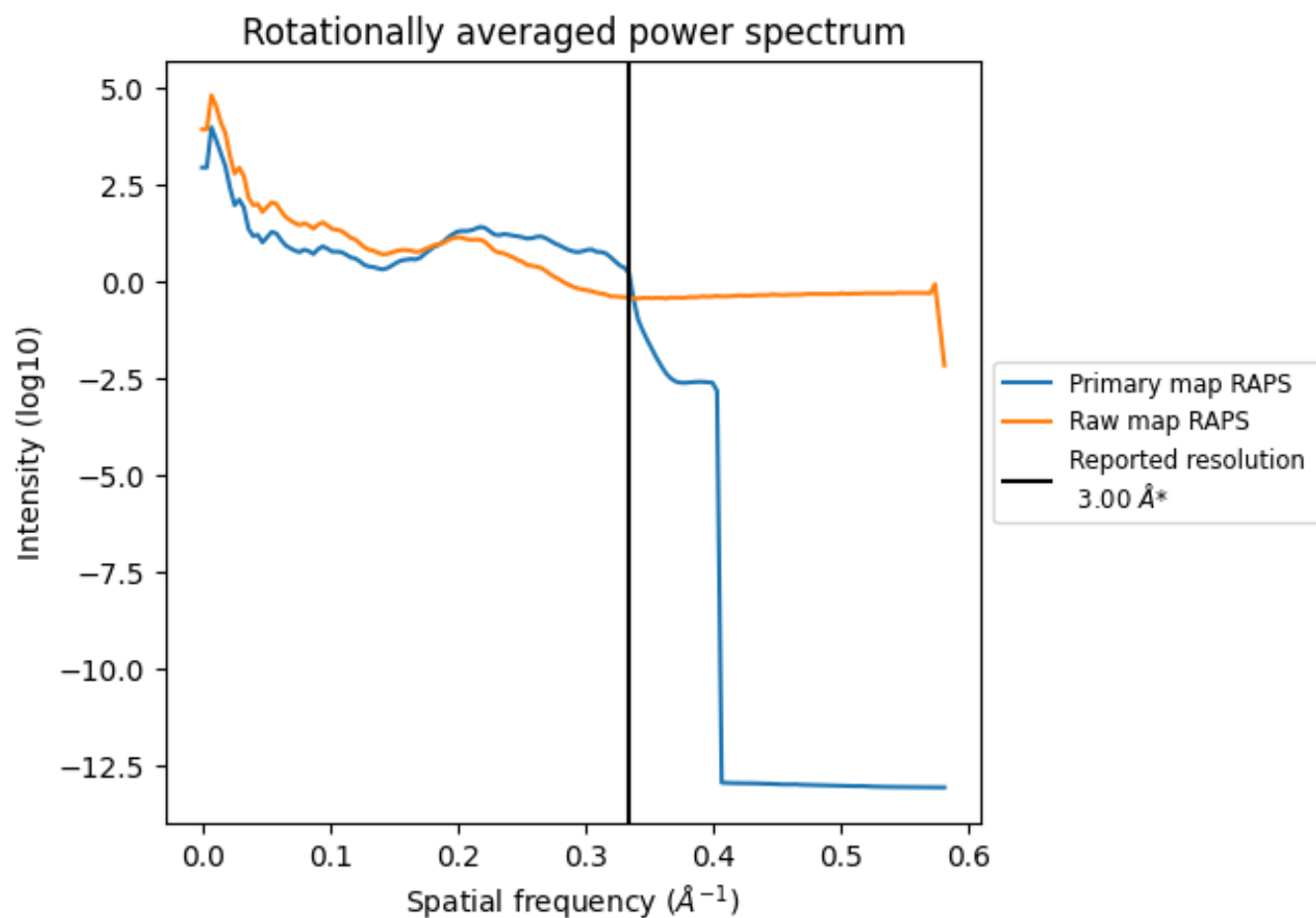
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 113 nm³; this corresponds to an approximate mass of 102 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 ⓘ

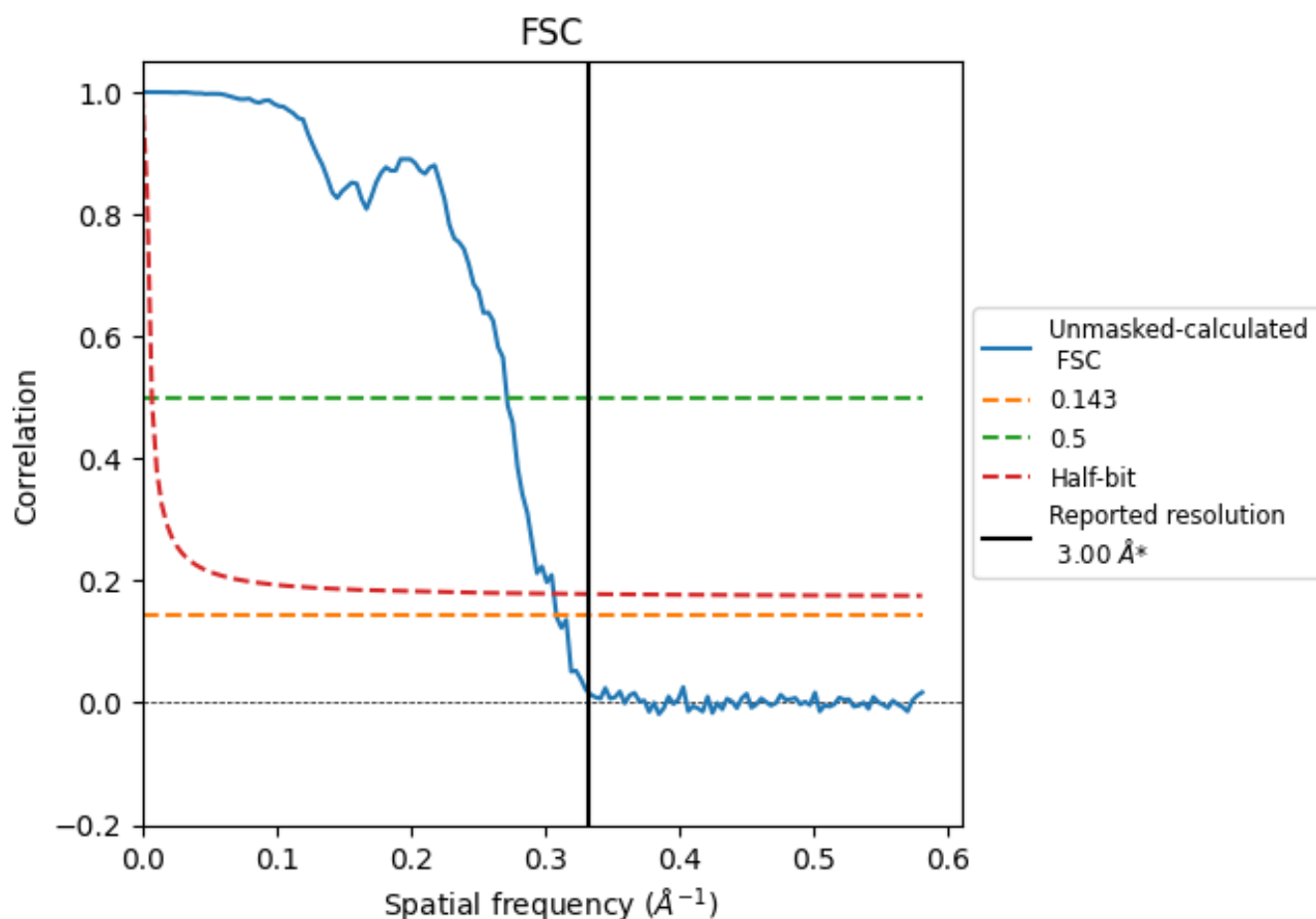


*Reported resolution corresponds to spatial frequency of 0.333 Å⁻¹

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 $0.333 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

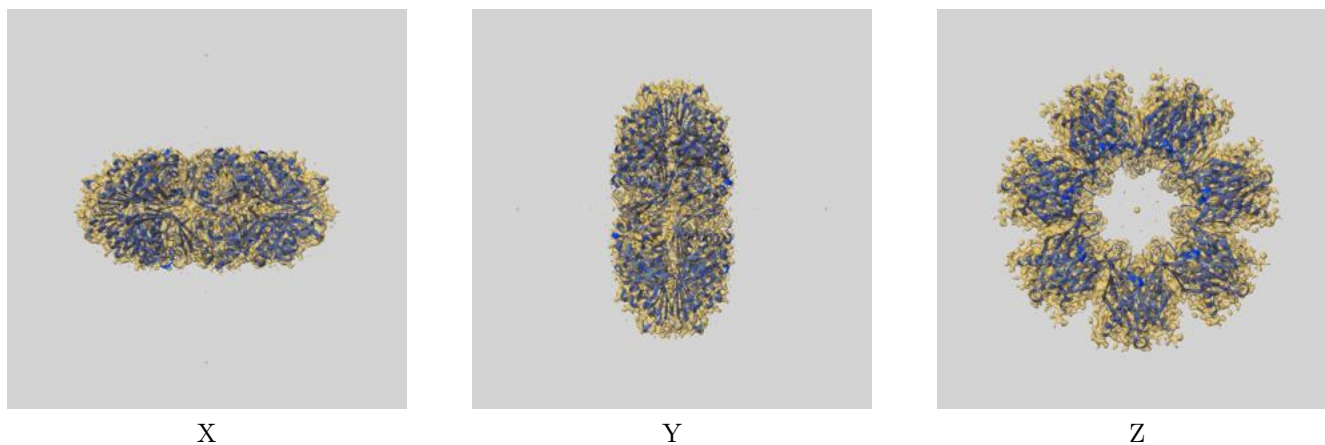
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.00	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.24	3.68	3.26

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

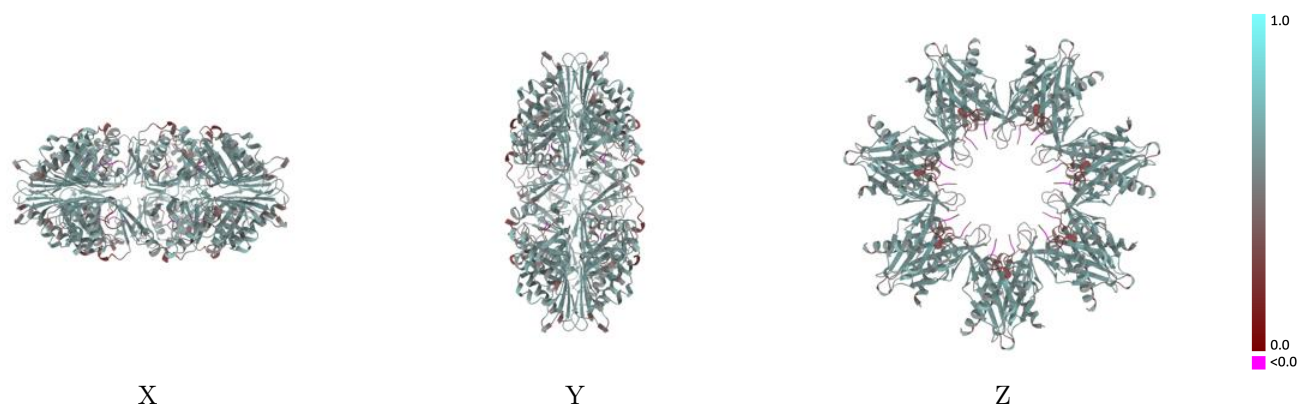
This section contains information regarding the fit between EMDB map EMD-41070 and PDB model 8T6K. Per-residue inclusion information can be found in [section 3](#) on [page 22](#).

9.1 Map-model overlay [i](#)



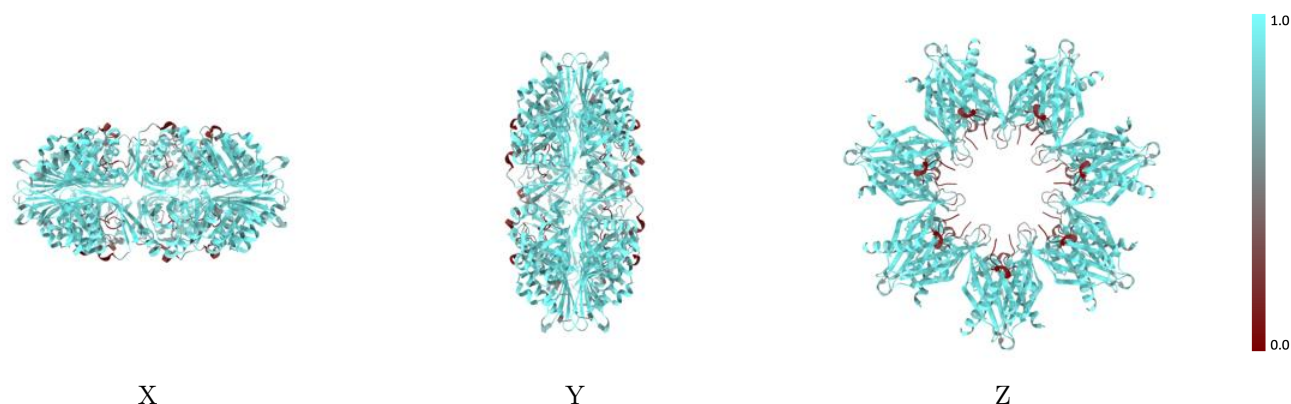
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



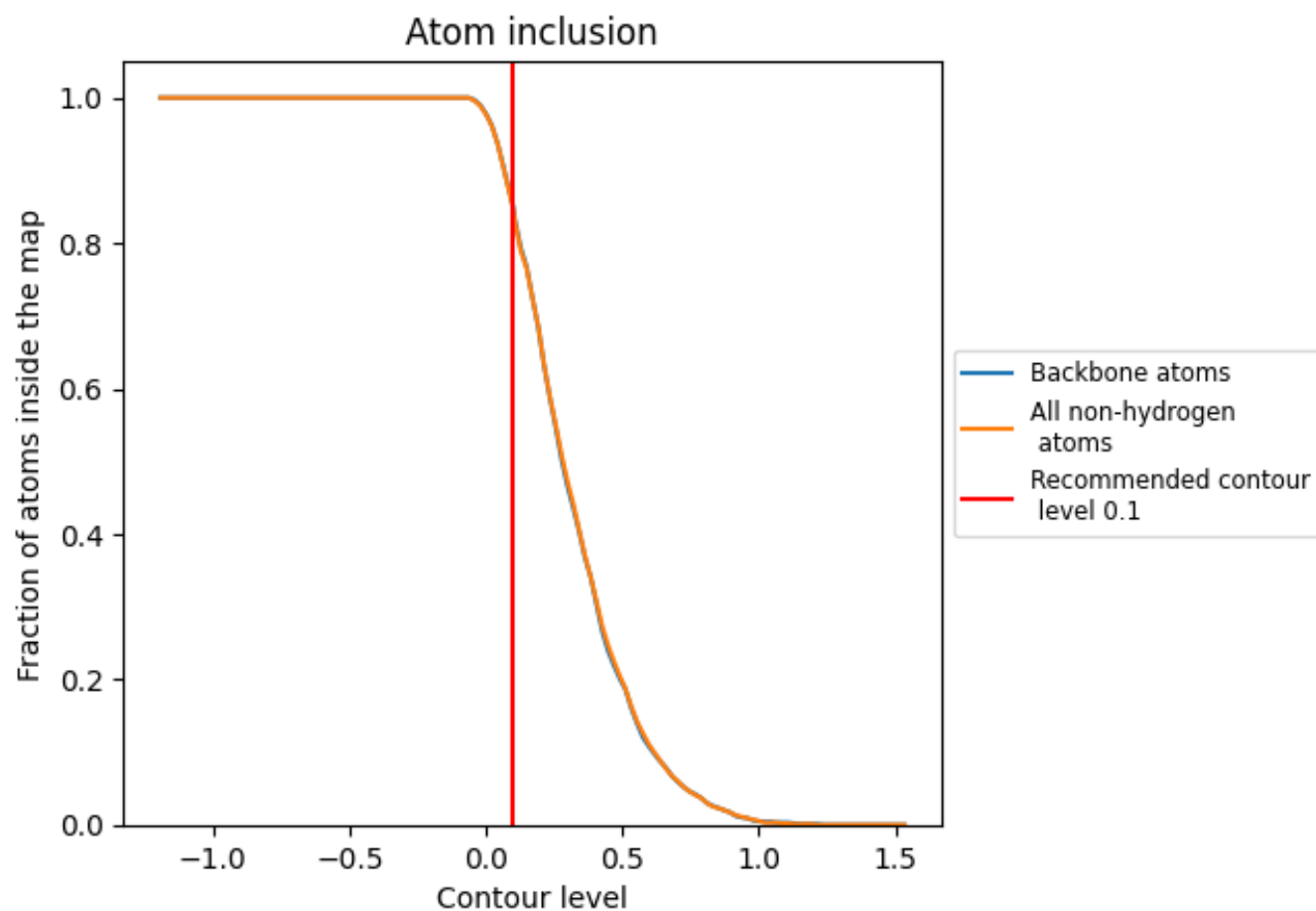
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 85% 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 (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8490	0.5390
A	0.8520	0.5400
B	0.8490	0.5380
C	0.8500	0.5400
D	0.8490	0.5400
E	0.8470	0.5400
F	0.8520	0.5390
G	0.8500	0.5390
H	0.8520	0.5400
I	0.8490	0.5400
J	0.8520	0.5400
K	0.8480	0.5390
L	0.8480	0.5380
M	0.8530	0.5380
N	0.8500	0.5380

1.0

0.0

<0.0