



Full wwPDB EM Validation Report ⓘ

Sep 24, 2026 – 12:33 PM EDT

PDB ID : 37BQ / pdb_000037bq
EMDB ID : EMD-78051
Title : Bacillus subtilis GerAA(A318C):GerAB:GerAC(S56C), in glyco-diosgenin
Authors : Cofsky, J.C.; Kruse, A.C.
Deposited on : 2026-07-11
Resolution : 2.02 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.dev133
Mogul : 2022.3.0, CSD as543be (2022)
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) : 2.50

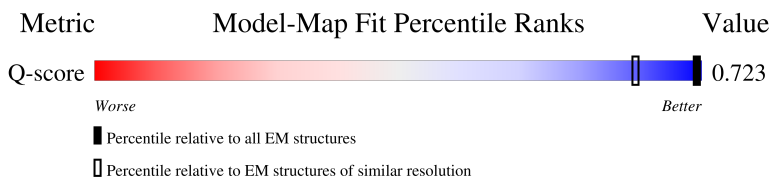
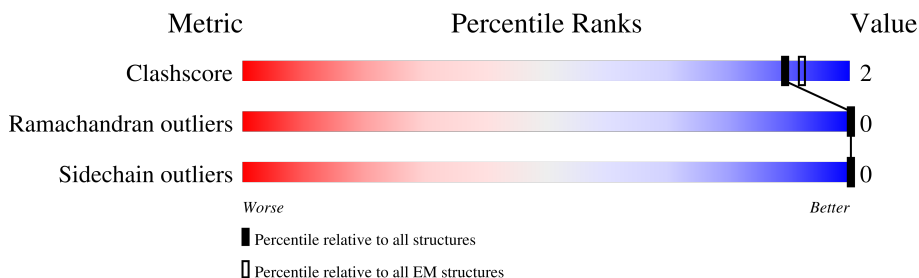
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.02 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	1728 (1.52 - 2.52)

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	505	
1	E	505	
1	I	505	
1	M	505	

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Mol	Chain	Length	Quality of chain
1	Q	505	
2	B	365	
2	F	365	
2	J	365	
2	N	365	
2	R	365	
3	C	356	
3	G	356	
3	K	356	
3	O	356	
3	S	356	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 46415 atoms, of which 0 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 Spore germination protein A1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	460	Total	C	N	O	S	11	0
			3703	2405	605	682	11		
1	E	460	Total	C	N	O	S	11	0
			3703	2405	605	682	11		
1	I	460	Total	C	N	O	S	11	0
			3703	2405	605	682	11		
1	M	460	Total	C	N	O	S	11	0
			3703	2405	605	682	11		
1	Q	460	Total	C	N	O	S	11	0
			3703	2405	605	682	11		

There are 125 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	ASP	-	expression tag	UNP P07868
A	-21	TYR	-	expression tag	UNP P07868
A	-20	LYS	-	expression tag	UNP P07868
A	-19	ASP	-	expression tag	UNP P07868
A	-18	ASP	-	expression tag	UNP P07868
A	-17	ASP	-	expression tag	UNP P07868
A	-16	ASP	-	expression tag	UNP P07868
A	-15	LYS	-	expression tag	UNP P07868
A	-14	GLY	-	expression tag	UNP P07868
A	-13	SER	-	expression tag	UNP P07868
A	-12	LEU	-	expression tag	UNP P07868
A	-11	GLU	-	expression tag	UNP P07868
A	-10	VAL	-	expression tag	UNP P07868
A	-9	LEU	-	expression tag	UNP P07868
A	-8	PHE	-	expression tag	UNP P07868
A	-7	GLN	-	expression tag	UNP P07868
A	-6	GLY	-	expression tag	UNP P07868
A	-5	PRO	-	expression tag	UNP P07868
A	-4	GLY	-	expression tag	UNP P07868
A	-3	GLY	-	expression tag	UNP P07868

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP P07868
A	-1	GLY	-	expression tag	UNP P07868
A	0	SER	-	expression tag	UNP P07868
A	299	ALA	THR	conflict	UNP P07868
A	318	CYS	ALA	engineered mutation	UNP P07868
E	-22	ASP	-	expression tag	UNP P07868
E	-21	TYR	-	expression tag	UNP P07868
E	-20	LYS	-	expression tag	UNP P07868
E	-19	ASP	-	expression tag	UNP P07868
E	-18	ASP	-	expression tag	UNP P07868
E	-17	ASP	-	expression tag	UNP P07868
E	-16	ASP	-	expression tag	UNP P07868
E	-15	LYS	-	expression tag	UNP P07868
E	-14	GLY	-	expression tag	UNP P07868
E	-13	SER	-	expression tag	UNP P07868
E	-12	LEU	-	expression tag	UNP P07868
E	-11	GLU	-	expression tag	UNP P07868
E	-10	VAL	-	expression tag	UNP P07868
E	-9	LEU	-	expression tag	UNP P07868
E	-8	PHE	-	expression tag	UNP P07868
E	-7	GLN	-	expression tag	UNP P07868
E	-6	GLY	-	expression tag	UNP P07868
E	-5	PRO	-	expression tag	UNP P07868
E	-4	GLY	-	expression tag	UNP P07868
E	-3	GLY	-	expression tag	UNP P07868
E	-2	SER	-	expression tag	UNP P07868
E	-1	GLY	-	expression tag	UNP P07868
E	0	SER	-	expression tag	UNP P07868
E	299	ALA	THR	conflict	UNP P07868
E	318	CYS	ALA	engineered mutation	UNP P07868
I	-22	ASP	-	expression tag	UNP P07868
I	-21	TYR	-	expression tag	UNP P07868
I	-20	LYS	-	expression tag	UNP P07868
I	-19	ASP	-	expression tag	UNP P07868
I	-18	ASP	-	expression tag	UNP P07868
I	-17	ASP	-	expression tag	UNP P07868
I	-16	ASP	-	expression tag	UNP P07868
I	-15	LYS	-	expression tag	UNP P07868
I	-14	GLY	-	expression tag	UNP P07868
I	-13	SER	-	expression tag	UNP P07868
I	-12	LEU	-	expression tag	UNP P07868
I	-11	GLU	-	expression tag	UNP P07868

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Chain	Residue	Modelled	Actual	Comment	Reference
I	-10	VAL	-	expression tag	UNP P07868
I	-9	LEU	-	expression tag	UNP P07868
I	-8	PHE	-	expression tag	UNP P07868
I	-7	GLN	-	expression tag	UNP P07868
I	-6	GLY	-	expression tag	UNP P07868
I	-5	PRO	-	expression tag	UNP P07868
I	-4	GLY	-	expression tag	UNP P07868
I	-3	GLY	-	expression tag	UNP P07868
I	-2	SER	-	expression tag	UNP P07868
I	-1	GLY	-	expression tag	UNP P07868
I	0	SER	-	expression tag	UNP P07868
I	299	ALA	THR	conflict	UNP P07868
I	318	CYS	ALA	engineered mutation	UNP P07868
M	-22	ASP	-	expression tag	UNP P07868
M	-21	TYR	-	expression tag	UNP P07868
M	-20	LYS	-	expression tag	UNP P07868
M	-19	ASP	-	expression tag	UNP P07868
M	-18	ASP	-	expression tag	UNP P07868
M	-17	ASP	-	expression tag	UNP P07868
M	-16	ASP	-	expression tag	UNP P07868
M	-15	LYS	-	expression tag	UNP P07868
M	-14	GLY	-	expression tag	UNP P07868
M	-13	SER	-	expression tag	UNP P07868
M	-12	LEU	-	expression tag	UNP P07868
M	-11	GLU	-	expression tag	UNP P07868
M	-10	VAL	-	expression tag	UNP P07868
M	-9	LEU	-	expression tag	UNP P07868
M	-8	PHE	-	expression tag	UNP P07868
M	-7	GLN	-	expression tag	UNP P07868
M	-6	GLY	-	expression tag	UNP P07868
M	-5	PRO	-	expression tag	UNP P07868
M	-4	GLY	-	expression tag	UNP P07868
M	-3	GLY	-	expression tag	UNP P07868
M	-2	SER	-	expression tag	UNP P07868
M	-1	GLY	-	expression tag	UNP P07868
M	0	SER	-	expression tag	UNP P07868
M	299	ALA	THR	conflict	UNP P07868
M	318	CYS	ALA	engineered mutation	UNP P07868
Q	-22	ASP	-	expression tag	UNP P07868
Q	-21	TYR	-	expression tag	UNP P07868
Q	-20	LYS	-	expression tag	UNP P07868
Q	-19	ASP	-	expression tag	UNP P07868

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Chain	Residue	Modelled	Actual	Comment	Reference
Q	-18	ASP	-	expression tag	UNP P07868
Q	-17	ASP	-	expression tag	UNP P07868
Q	-16	ASP	-	expression tag	UNP P07868
Q	-15	LYS	-	expression tag	UNP P07868
Q	-14	GLY	-	expression tag	UNP P07868
Q	-13	SER	-	expression tag	UNP P07868
Q	-12	LEU	-	expression tag	UNP P07868
Q	-11	GLU	-	expression tag	UNP P07868
Q	-10	VAL	-	expression tag	UNP P07868
Q	-9	LEU	-	expression tag	UNP P07868
Q	-8	PHE	-	expression tag	UNP P07868
Q	-7	GLN	-	expression tag	UNP P07868
Q	-6	GLY	-	expression tag	UNP P07868
Q	-5	PRO	-	expression tag	UNP P07868
Q	-4	GLY	-	expression tag	UNP P07868
Q	-3	GLY	-	expression tag	UNP P07868
Q	-2	SER	-	expression tag	UNP P07868
Q	-1	GLY	-	expression tag	UNP P07868
Q	0	SER	-	expression tag	UNP P07868
Q	299	ALA	THR	conflict	UNP P07868
Q	318	CYS	ALA	engineered mutation	UNP P07868

- Molecule 2 is a protein called Spore germination protein A2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	353	Total	C	N	O	S	5	0
			2874	1962	425	472	15		
2	F	353	Total	C	N	O	S	5	0
			2874	1962	425	472	15		
2	J	353	Total	C	N	O	S	5	0
			2874	1962	425	472	15		
2	N	353	Total	C	N	O	S	5	0
			2874	1962	425	472	15		
2	R	353	Total	C	N	O	S	5	0
			2874	1962	425	472	15		

- Molecule 3 is a protein called Spore germination protein A3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	316	Total	C	N	O	S	2	0
			2550	1609	443	491	7		

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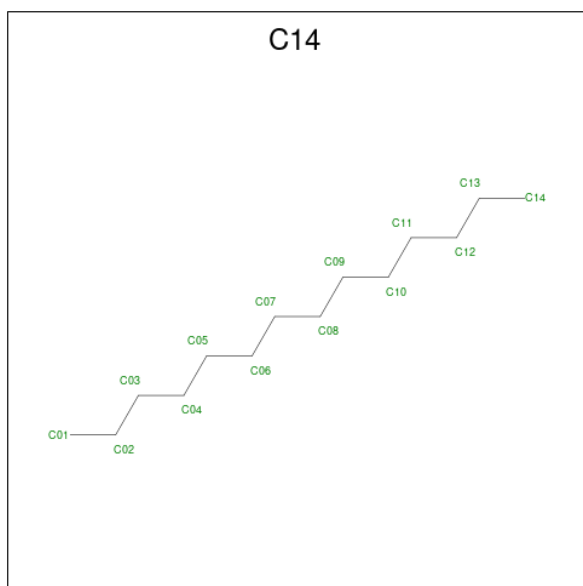
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Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	316	Total	C	N	O	S	2	0
			2550	1609	443	491	7		
3	K	316	Total	C	N	O	S	2	0
			2550	1609	443	491	7		
3	O	316	Total	C	N	O	S	2	0
			2550	1609	443	491	7		
3	S	316	Total	C	N	O	S	2	0
			2550	1609	443	491	7		

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	56	CYS	SER	engineered mutation	UNP P07870
G	56	CYS	SER	engineered mutation	UNP P07870
K	56	CYS	SER	engineered mutation	UNP P07870
O	56	CYS	SER	engineered mutation	UNP P07870
S	56	CYS	SER	engineered mutation	UNP P07870

- Molecule 4 is TETRADECANE (CCD ID: C14) (formula: C₁₄H₃₀).



Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	C	0
			14	14	
4	A	1	Total	C	0
			14	14	
4	I	1	Total	C	0
			14	14	

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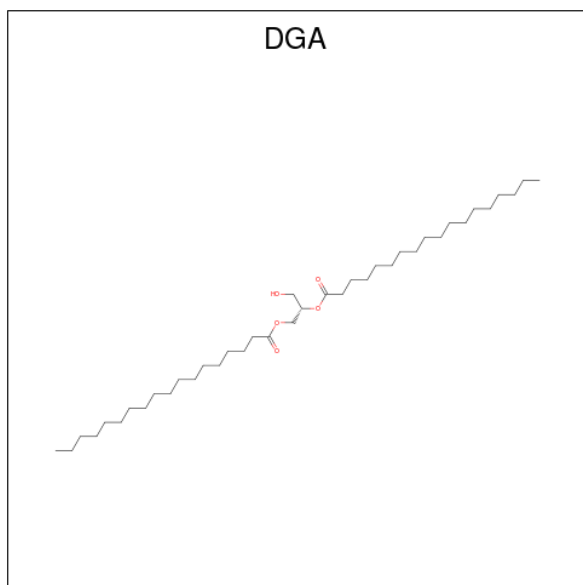
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Mol	Chain	Residues	Atoms	AltConf
4	M	1	Total C 14 14	0
4	Q	1	Total C 14 14	0

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	AltConf
5	B	1	Total Na 1 1	0
5	F	1	Total Na 1 1	0
5	J	1	Total Na 1 1	0
5	N	1	Total Na 1 1	0
5	R	1	Total Na 1 1	0

- Molecule 6 is DIACYL GLYCEROL (CCD ID: DGA) (formula: C₃₉H₇₆O₅).



Mol	Chain	Residues	Atoms	AltConf
6	C	1	Total C O 11 7 4	0
6	G	1	Total C O 11 7 4	0

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Mol	Chain	Residues	Atoms			AltConf
6	K	1	Total	C	O	0
			11	7	4	
6	O	1	Total	C	O	0
			11	7	4	
6	S	1	Total	C	O	0
			11	7	4	

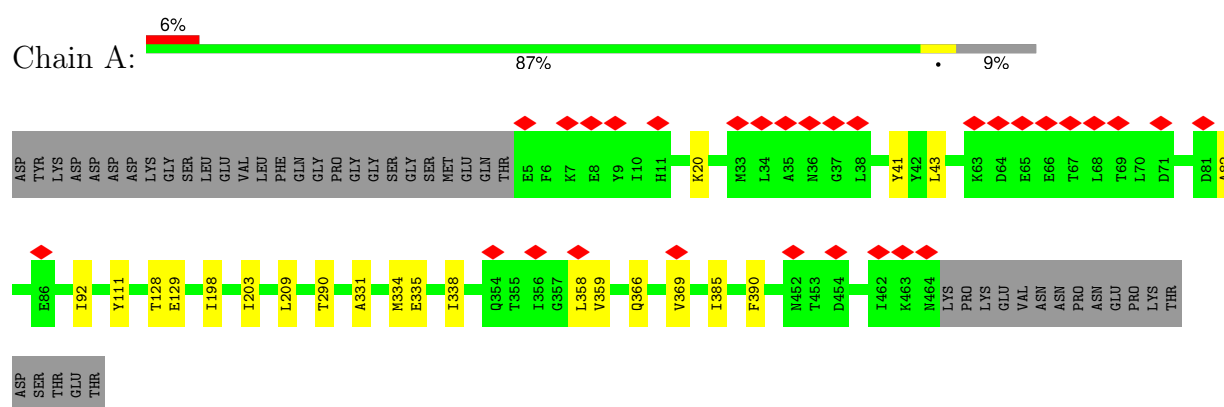
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		AltConf
7	A	62	Total	O	0
			62	62	
7	B	46	Total	O	0
			46	46	
7	C	22	Total	O	0
			22	22	
7	E	62	Total	O	0
			62	62	
7	F	46	Total	O	0
			46	46	
7	G	22	Total	O	0
			22	22	
7	I	62	Total	O	0
			62	62	
7	J	46	Total	O	0
			46	46	
7	K	22	Total	O	0
			22	22	
7	M	62	Total	O	0
			62	62	
7	N	46	Total	O	0
			46	46	
7	O	22	Total	O	0
			22	22	
7	Q	62	Total	O	0
			62	62	
7	R	46	Total	O	0
			46	46	
7	S	22	Total	O	0
			22	22	

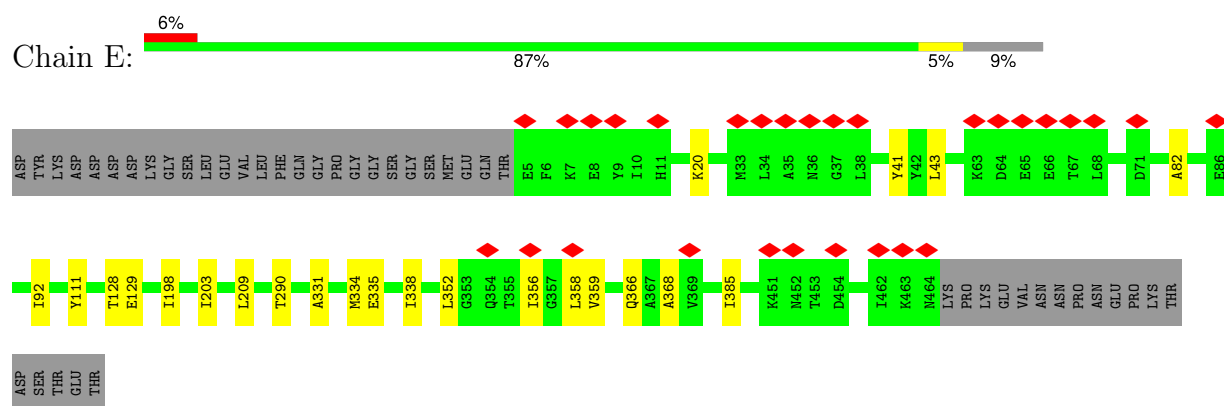
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

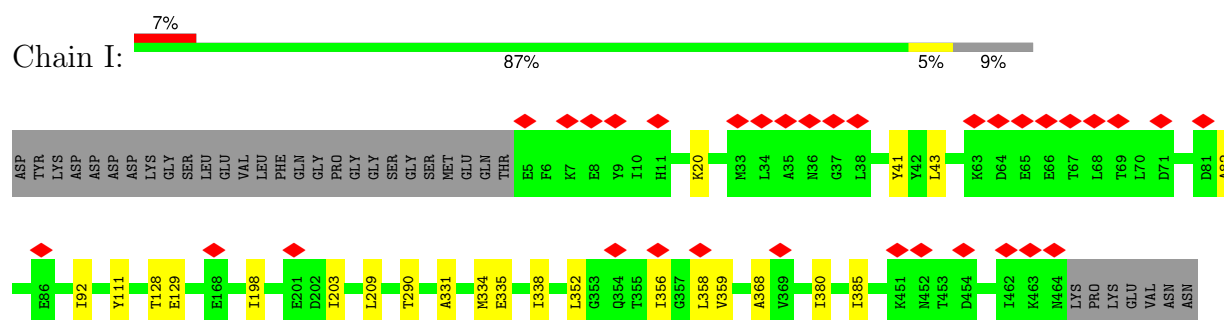
- Molecule 1: Spore germination protein A1



- Molecule 1: Spore germination protein A1



- Molecule 1: Spore germination protein A1



PRO
ASN
GLU
PRO
PRO
THR
THR
ASP
SER
THR
GLU
THR

- Molecule 1: Spore germination protein A1

Chain M: 6% 87% 9%

ASP
TYR
LYS
ASP
ASP
LYS
ASP
ASP
LYS
GLY
SER
LEU
GLU
VAL
LEU
PHE
GLN
GLY
PRO
GLY
GLY
GLY
SER
SER
MET
GLU
GLN
THR
E5
F6
K7
E8
Y9
I10
H11
K20
M33
L34
A35
N36
G37
L38
Y41
Y42
L43
D64
E65
E66
T67
L68
T69
D71
D81
A82

LYS
THR
ASP
SER
GLU
THR

- Molecule 1: Spore germination protein A1

Chain Q: 6% 87% 5% 9%

ASP
TYR
LYS
ASP
ASP
LYS
ASP
ASP
LYS
GLY
SER
LEU
GLU
VAL
LEU
PHE
GLN
GLY
PRO
GLY
GLY
SER
SER
MET
GLU
GLN
THR
E5
F6
K7
E8
Y9
I10
H11
K20
M33
L34
A35
N36
G37
L38
Y41
Y42
L43
K63
D64
E65
E66
T67
L68
T69
D71
A82

PRO
LYS
THR
ASP
SER
GLU
THR

- Molecule 2: Spore germination protein A2

Chain B: 92% 5%

MET
SER
GLN
LYS
GLN
THR
PRO
L8
L99
G134
M138
I196
S197
F198
M201
F225
L226
I229
L264
K265
G266
I267
F268
I269
E270
R271
F272
F275
I314
D327
V331
G339
R360
ARG
ILE
THR
THR
LYS

- Molecule 2: Spore germination protein A2

Chain F: 92% 6%

MET
SER
GLN
LYS
GLN
THR
PRO
L8
L99
I196
S197
F198
M201
F225
L226
I229
L264
K265
G266
I267
F268
I269
E270
R271
F272
F275
I314
D327
V331
G339
R360
ARG
ILE
THR
THR
LYS

- Molecule 2: Spore germination protein A2

Chain J: 91% 6%



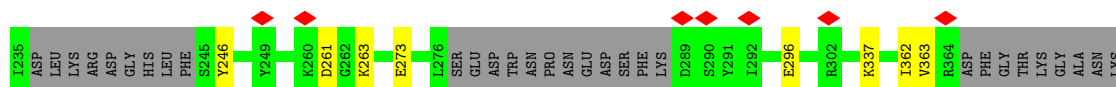
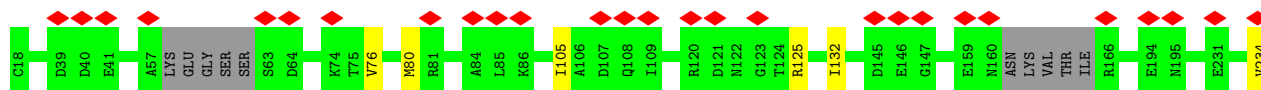
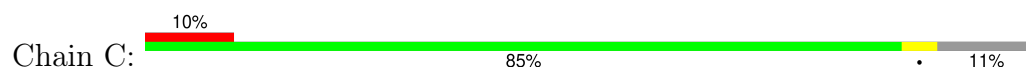
- Molecule 2: Spore germination protein A2



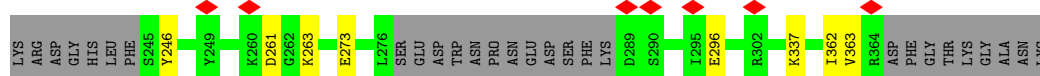
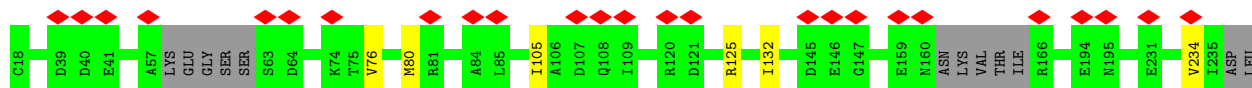
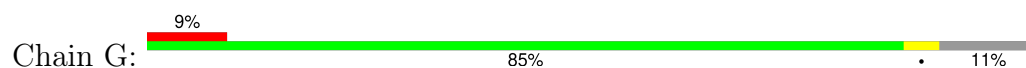
- Molecule 2: Spore germination protein A2



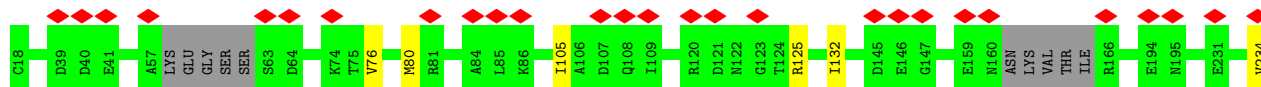
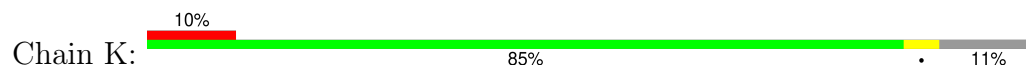
- Molecule 3: Spore germination protein A3



- Molecule 3: Spore germination protein A3

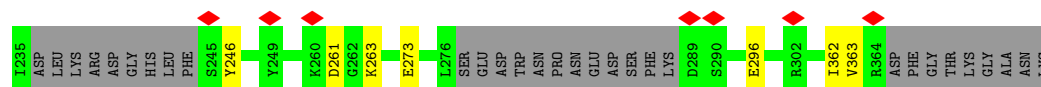
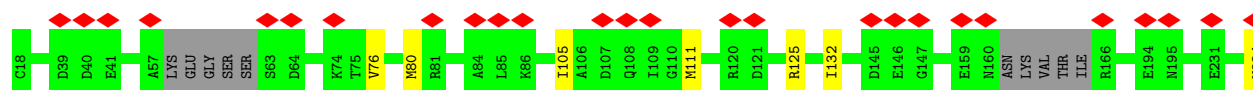
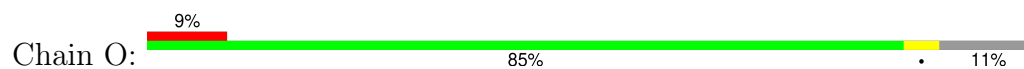


- Molecule 3: Spore germination protein A3

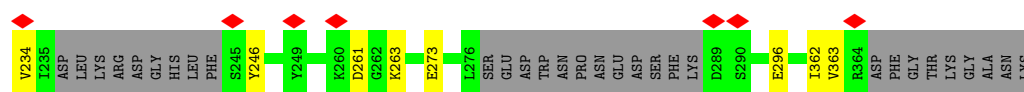
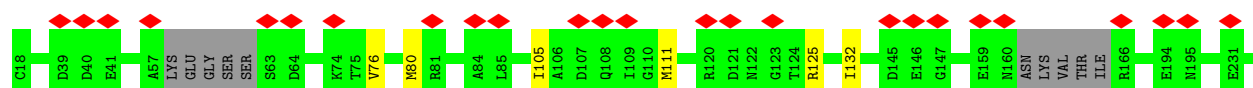
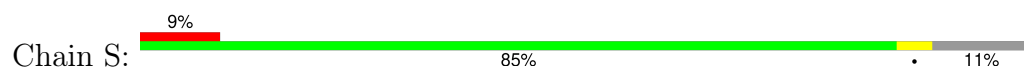




• Molecule 3: Spore germination protein A3



• Molecule 3: Spore germination protein A3



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C5	Depositor
Number of particles used	309046	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	52.4	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	165000	Depositor
Image detector	TFS FALCON 4i (4k x 4k)	Depositor
Maximum map value	52.649	Depositor
Minimum map value	-31.560	Depositor
Average map value	0.010	Depositor
Map value standard deviation	1.359	Depositor
Recommended contour level	6	Depositor
Map size (Å)	331.52002, 331.52002, 331.52002	wwPDB
Map dimensions	448, 448, 448	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.74000007, 0.74000007, 0.74000007	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: C14, NA, DGA

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	0.09	0/3770	0.24	0/5108
1	E	0.10	0/3770	0.24	0/5108
1	I	0.09	0/3770	0.24	0/5108
1	M	0.10	0/3770	0.24	0/5108
1	Q	0.10	0/3770	0.24	0/5108
2	B	0.12	0/2951	0.28	0/4006
2	F	0.13	0/2951	0.28	0/4006
2	J	0.13	0/2951	0.28	0/4006
2	N	0.13	0/2951	0.28	0/4006
2	R	0.13	0/2951	0.28	0/4006
3	C	0.10	0/2589	0.23	0/3488
3	G	0.10	0/2589	0.24	0/3488
3	K	0.10	0/2589	0.24	0/3488
3	O	0.10	0/2589	0.24	0/3488
3	S	0.10	0/2589	0.24	0/3488
All	All	0.11	0/46550	0.25	0/63010

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3703	0	3864	19	0
1	E	3703	0	3864	20	0
1	I	3703	0	3864	18	0
1	M	3703	0	3864	17	0
1	Q	3703	0	3864	18	0
2	B	2874	0	3024	11	0
2	F	2874	0	3024	10	0
2	J	2874	0	3024	12	0
2	N	2874	0	3024	12	0
2	R	2874	0	3024	11	0
3	C	2550	0	2576	8	0
3	G	2550	0	2576	8	0
3	K	2550	0	2576	8	0
3	O	2550	0	2576	8	0
3	S	2550	0	2576	8	0
4	A	28	0	60	6	0
4	I	14	0	30	3	0
4	M	14	0	30	3	0
4	Q	14	0	30	3	0
5	B	1	0	0	0	0
5	F	1	0	0	0	0
5	J	1	0	0	0	0
5	N	1	0	0	0	0
5	R	1	0	0	0	0
6	C	11	0	3	0	0
6	G	11	0	3	0	0
6	K	11	0	3	0	0
6	O	11	0	3	0	0
6	S	11	0	3	0	0
7	A	62	0	0	0	0
7	B	46	0	0	0	0
7	C	22	0	0	0	0
7	E	62	0	0	0	0
7	F	46	0	0	0	0
7	G	22	0	0	0	0
7	I	62	0	0	0	0
7	J	46	0	0	0	0
7	K	22	0	0	0	0
7	M	62	0	0	0	0
7	N	46	0	0	0	0
7	O	22	0	0	0	0
7	Q	62	0	0	0	0
7	R	46	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	S	22	0	0	0	0
All	All	46415	0	47485	169	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

All (169) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:198:PHE:HB3	2:R:201[B]:MET:HE3	1.78	0.66
2:J:198:PHE:HB3	2:J:201[B]:MET:HE3	1.78	0.65
2:F:198:PHE:HB3	2:F:201[B]:MET:HE3	1.78	0.64
2:N:198:PHE:HB3	2:N:201[B]:MET:HE3	1.78	0.64
2:B:198:PHE:HB3	2:B:201[B]:MET:HE3	1.78	0.64
1:A:366:GLN:HE22	1:E:366:GLN:HB3	1.68	0.59
4:A:502:C14:H122	1:E:338[A]:ILE:HD12	1.85	0.58
1:Q:338[A]:ILE:HD12	4:Q:501:C14:H122	1.85	0.58
1:A:359:VAL:HG12	4:A:502:C14:H111	1.84	0.58
1:A:338[A]:ILE:HD12	4:A:501:C14:H122	1.85	0.58
1:I:338[A]:ILE:HD12	4:I:501:C14:H122	1.85	0.58
1:I:358:LEU:HD23	1:M:358:LEU:HD21	1.85	0.58
1:M:338[A]:ILE:HD12	4:M:501:C14:H122	1.85	0.58
1:I:359:VAL:HG12	4:M:501:C14:H111	1.88	0.55
1:M:331:ALA:O	1:M:335:GLU:HG2	2.09	0.53
1:M:359:VAL:HG12	4:Q:501:C14:H111	1.90	0.53
1:E:331:ALA:O	1:E:335:GLU:HG2	2.09	0.53
1:I:331:ALA:O	1:I:335:GLU:HG2	2.09	0.53
2:R:264:LEU:HB2	2:R:272:PHE:HE1	1.74	0.53
1:A:331:ALA:O	1:A:335:GLU:HG2	2.09	0.53
1:Q:331:ALA:O	1:Q:335:GLU:HG2	2.09	0.53
3:G:273:GLU:HG2	3:G:362:ILE:HB	1.92	0.52
1:I:41[B]:TYR:HE1	1:I:43:LEU:HD21	1.75	0.52
2:N:264:LEU:HB2	2:N:272:PHE:HE1	1.74	0.52
2:B:264:LEU:HB2	2:B:272:PHE:HE1	1.74	0.52
3:O:273:GLU:HG2	3:O:362:ILE:HB	1.92	0.52
4:A:501:C14:H111	1:Q:359:VAL:HG12	1.90	0.52
2:F:264:LEU:HB2	2:F:272:PHE:HE1	1.74	0.52
2:J:264:LEU:HB2	2:J:272:PHE:HE1	1.74	0.52
1:A:41[B]:TYR:HE1	1:A:43:LEU:HD21	1.75	0.51
1:E:359:VAL:HG12	4:I:501:C14:H111	1.91	0.51
1:Q:41[B]:TYR:HE1	1:Q:43:LEU:HD21	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:S:273:GLU:HG2	3:S:362:ILE:HB	1.92	0.51
2:J:269:ILE:HG12	2:J:275:PHE:HE2	1.76	0.51
3:K:273:GLU:HG2	3:K:362:ILE:HB	1.92	0.51
3:C:273:GLU:HG2	3:C:362:ILE:HB	1.92	0.51
2:N:269:ILE:HG12	2:N:275:PHE:HE2	1.76	0.51
1:M:41[B]:TYR:HE1	1:M:43:LEU:HD21	1.75	0.51
1:E:359:VAL:HG23	1:I:358:LEU:HD12	1.92	0.51
2:F:269:ILE:HG12	2:F:275:PHE:HE2	1.76	0.51
2:B:269:ILE:HG12	2:B:275:PHE:HE2	1.76	0.50
1:E:41[B]:TYR:HE1	1:E:43:LEU:HD21	1.75	0.50
2:R:269:ILE:HG12	2:R:275:PHE:HE2	1.76	0.50
2:N:196:ILE:HB	2:N:339:GLY:HA3	1.94	0.50
2:B:196:ILE:HB	2:B:339:GLY:HA3	1.94	0.49
2:F:196:ILE:HB	2:F:339:GLY:HA3	1.94	0.49
2:J:196:ILE:HB	2:J:339:GLY:HA3	1.94	0.49
2:F:99:LEU:HD21	2:F:314:ILE:HG12	1.95	0.49
2:R:196:ILE:HB	2:R:339:GLY:HA3	1.94	0.48
1:M:358:LEU:HB3	1:Q:358:LEU:HD21	1.95	0.48
2:N:99:LEU:HD21	2:N:314:ILE:HG12	1.95	0.48
2:J:99:LEU:HD21	2:J:314:ILE:HG12	1.95	0.48
2:R:99:LEU:HD21	2:R:314:ILE:HG12	1.95	0.47
3:K:105:ILE:HD13	3:K:132:ILE:HD12	1.96	0.47
3:O:105:ILE:HD13	3:O:132:ILE:HD12	1.96	0.47
2:B:99:LEU:HD21	2:B:314:ILE:HG12	1.95	0.47
3:G:234:VAL:HA	3:G:246:TYR:O	2.15	0.47
3:S:234:VAL:HA	3:S:246:TYR:O	2.15	0.47
3:C:105:ILE:HD13	3:C:132:ILE:HD12	1.96	0.47
3:O:234:VAL:HA	3:O:246:TYR:O	2.15	0.47
3:C:234:VAL:HA	3:C:246:TYR:O	2.15	0.47
3:G:105:ILE:HD13	3:G:132:ILE:HD12	1.96	0.47
2:R:226[B]:LEU:HA	2:R:229:ILE:HG22	1.97	0.47
3:S:105:ILE:HD13	3:S:132:ILE:HD12	1.96	0.47
3:K:234:VAL:HA	3:K:246:TYR:O	2.15	0.47
2:N:226[B]:LEU:HA	2:N:229:ILE:HG22	1.97	0.47
1:Q:203:ILE:HD13	1:Q:209:LEU:HD21	1.97	0.46
1:A:203:ILE:HD13	1:A:209:LEU:HD21	1.97	0.46
2:F:226[B]:LEU:HA	2:F:229:ILE:HG22	1.97	0.46
1:I:198:ILE:HD13	1:I:209:LEU:HD22	1.98	0.46
2:J:226[B]:LEU:HA	2:J:229:ILE:HG22	1.97	0.46
1:M:198:ILE:HD13	1:M:209:LEU:HD22	1.98	0.46
1:E:198:ILE:HD13	1:E:209:LEU:HD22	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:203:ILE:HD13	1:M:209:LEU:HD21	1.97	0.46
1:A:198:ILE:HD13	1:A:209:LEU:HD22	1.98	0.46
2:B:226[B]:LEU:HA	2:B:229:ILE:HG22	1.97	0.46
1:Q:198:ILE:HD13	1:Q:209:LEU:HD22	1.98	0.46
1:E:203:ILE:HD13	1:E:209:LEU:HD21	1.97	0.45
1:I:352:LEU:O	1:I:356:ILE:HG12	2.17	0.45
2:F:271:ARG:CZ	3:G:125:ARG:HD2	2.47	0.45
1:I:203:ILE:HD13	1:I:209:LEU:HD21	1.97	0.45
2:J:271:ARG:CZ	3:K:125:ARG:HD2	2.47	0.45
2:B:271:ARG:CZ	3:C:125:ARG:HD2	2.47	0.45
1:E:128:THR:HG22	1:E:129:GLU:HG3	1.99	0.45
1:Q:352:LEU:O	1:Q:356:ILE:HG12	2.17	0.45
1:M:359:VAL:HG23	1:Q:358:LEU:HD23	1.98	0.45
1:I:128:THR:HG22	1:I:129:GLU:HG3	1.99	0.44
2:N:271:ARG:CZ	3:O:125:ARG:HD2	2.47	0.44
1:M:128:THR:HG22	1:M:129:GLU:HG3	1.99	0.44
1:A:290:THR:HA	1:A:385:ILE:HG23	2.00	0.44
1:E:290:THR:HA	1:E:385:ILE:HG23	2.00	0.44
1:I:334:MET:O	1:I:338[A]:ILE:HG12	2.18	0.44
1:M:334:MET:O	1:M:338[A]:ILE:HG12	2.18	0.44
1:A:128:THR:HG22	1:A:129:GLU:HG3	1.99	0.44
1:A:334:MET:O	1:A:338[A]:ILE:HG12	2.18	0.44
1:I:290:THR:HA	1:I:385:ILE:HG23	2.00	0.44
2:R:271:ARG:CZ	3:S:125:ARG:HD2	2.47	0.44
1:M:290:THR:HA	1:M:385:ILE:HG23	2.00	0.44
1:Q:290:THR:HA	1:Q:385:ILE:HG23	2.00	0.44
2:B:201[A]:MET:SD	2:B:225:PHE:HB2	2.58	0.43
1:Q:128:THR:HG22	1:Q:129:GLU:HG3	1.99	0.43
2:F:201[A]:MET:SD	2:F:225:PHE:HB2	2.58	0.43
3:K:337[B]:LYS:HE3	3:K:337[B]:LYS:HB3	1.87	0.43
1:Q:334:MET:O	1:Q:338[A]:ILE:HG12	2.18	0.43
3:C:337[B]:LYS:HE3	3:C:337[B]:LYS:HB3	1.87	0.43
1:E:334:MET:O	1:E:338[A]:ILE:HG12	2.18	0.43
2:R:201[A]:MET:SD	2:R:225:PHE:HB2	2.58	0.43
3:C:296:GLU:HG2	3:C:363:VAL:HG23	2.00	0.43
3:O:296:GLU:HG2	3:O:363:VAL:HG23	2.00	0.43
1:M:82:ALA:HB1	1:M:111:TYR:HB3	2.01	0.43
1:E:20:LYS:HD3	1:E:92:ILE:HD11	2.01	0.43
1:Q:334:MET:HG2	4:Q:501:C14:H051	2.01	0.43
3:S:296:GLU:HG2	3:S:363:VAL:HG23	2.00	0.43
3:G:337[B]:LYS:HE3	3:G:337[B]:LYS:HB3	1.87	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:20:LYS:HD3	1:Q:92:ILE:HD11	2.01	0.43
1:A:20:LYS:HD3	1:A:92:ILE:HD11	2.01	0.43
1:A:358:LEU:HD22	1:E:358:LEU:HD11	2.01	0.43
3:G:261:ASP:OD1	3:G:263:LYS:HB3	2.19	0.43
3:O:261:ASP:OD1	3:O:263:LYS:HB3	2.19	0.43
2:J:201[A]:MET:SD	2:J:225:PHE:HB2	2.58	0.42
3:G:296:GLU:HG2	3:G:363:VAL:HG23	2.00	0.42
3:O:111:MET:HE2	3:O:111:MET:HB3	1.95	0.42
1:I:368:ALA:HA	1:M:380:ILE:HD11	2.01	0.42
3:S:261:ASP:OD1	3:S:263:LYS:HB3	2.19	0.42
1:A:369:VAL:HG11	1:Q:370:GLU:HG2	2.02	0.42
4:A:502:C14:H051	1:E:334:MET:HG2	2.01	0.42
3:K:296:GLU:HG2	3:K:363:VAL:HG23	2.00	0.42
1:M:334:MET:HG2	4:M:501:C14:H051	2.01	0.42
1:I:82:ALA:HB1	1:I:111:TYR:HB3	2.01	0.42
3:K:76:VAL:O	3:K:80:MET:HG3	2.20	0.42
3:C:76:VAL:O	3:C:80:MET:HG3	2.20	0.42
1:I:334:MET:HG2	4:I:501:C14:H051	2.01	0.42
3:S:111:MET:HE2	3:S:111:MET:HB3	1.95	0.42
1:E:82:ALA:HB1	1:E:111:TYR:HB3	2.01	0.42
1:A:334:MET:HG2	4:A:501:C14:H051	2.01	0.42
1:A:390:PHE:CZ	1:Q:356:ILE:HD12	2.55	0.42
3:K:261:ASP:OD1	3:K:263:LYS:HB3	2.19	0.42
2:N:201[A]:MET:SD	2:N:225:PHE:HB2	2.58	0.42
3:C:261:ASP:OD1	3:C:263:LYS:HB3	2.19	0.41
1:I:20:LYS:HD3	1:I:92:ILE:HD11	2.01	0.41
1:A:82:ALA:HB1	1:A:111:TYR:HB3	2.01	0.41
1:M:20:LYS:HD3	1:M:92:ILE:HD11	2.01	0.41
2:N:269:ILE:HG12	2:N:275:PHE:CE2	2.55	0.41
1:Q:82:ALA:HB1	1:Q:111:TYR:HB3	2.01	0.41
1:A:358:LEU:HB3	1:E:358:LEU:HD11	2.02	0.41
3:O:76:VAL:O	3:O:80:MET:HG3	2.20	0.41
2:R:134:GLY:O	2:R:138:MET:HG3	2.21	0.41
2:B:269:ILE:HG12	2:B:275:PHE:CE2	2.55	0.41
2:F:269:ILE:HG12	2:F:275:PHE:CE2	2.55	0.41
1:E:352:LEU:O	1:E:356:ILE:HG12	2.20	0.41
3:S:76:VAL:O	3:S:80:MET:HG3	2.20	0.41
2:J:327:ASP:O	2:J:331:VAL:HG23	2.20	0.41
2:B:134:GLY:O	2:B:138:MET:HG3	2.21	0.41
1:E:368:ALA:HA	1:I:380:ILE:HD11	2.03	0.41
2:F:327:ASP:O	2:F:331:VAL:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:76:VAL:O	3:G:80:MET:HG3	2.20	0.41
2:J:134:GLY:O	2:J:138:MET:HG3	2.21	0.41
2:R:327:ASP:O	2:R:331:VAL:HG23	2.20	0.41
1:A:358:LEU:HD13	1:E:358:LEU:HD21	2.01	0.40
2:B:327:ASP:O	2:B:331:VAL:HG23	2.20	0.40
2:N:327:ASP:O	2:N:331:VAL:HG23	2.20	0.40
1:A:366:GLN:HG2	1:Q:366:GLN:OE1	2.21	0.40
2:J:207:LEU:HD23	2:J:207:LEU:HA	1.94	0.40
2:J:8:LEU:HD13	2:J:143:SER:HA	2.04	0.40
2:N:134:GLY:O	2:N:138:MET:HG3	2.21	0.40
1:E:358:LEU:HB3	1:I:358:LEU:HD11	2.04	0.40
1:M:352:LEU:O	1:M:356:ILE:HG12	2.21	0.40
2:N:8:LEU:HD13	2:N:143:SER:HA	2.04	0.40
2:R:169:LYS:HE3	2:R:169:LYS:HB2	1.97	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	469/505 (93%)	467 (100%)	2 (0%)	0	100	100
1	E	469/505 (93%)	467 (100%)	2 (0%)	0	100	100
1	I	469/505 (93%)	467 (100%)	2 (0%)	0	100	100
1	M	469/505 (93%)	467 (100%)	2 (0%)	0	100	100
1	Q	469/505 (93%)	467 (100%)	2 (0%)	0	100	100
2	B	356/365 (98%)	352 (99%)	4 (1%)	0	100	100
2	F	356/365 (98%)	352 (99%)	4 (1%)	0	100	100
2	J	356/365 (98%)	352 (99%)	4 (1%)	0	100	100
2	N	356/365 (98%)	352 (99%)	4 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	R	356/365 (98%)	352 (99%)	4 (1%)	0	100	100
3	C	308/356 (86%)	298 (97%)	10 (3%)	0	100	100
3	G	308/356 (86%)	298 (97%)	10 (3%)	0	100	100
3	K	308/356 (86%)	298 (97%)	10 (3%)	0	100	100
3	O	308/356 (86%)	298 (97%)	10 (3%)	0	100	100
3	S	308/356 (86%)	298 (97%)	10 (3%)	0	100	100
All	All	5665/6130 (92%)	5585 (99%)	80 (1%)	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	419/448 (94%)	419 (100%)	0	100	100
1	E	419/448 (94%)	419 (100%)	0	100	100
1	I	419/448 (94%)	419 (100%)	0	100	100
1	M	419/448 (94%)	419 (100%)	0	100	100
1	Q	419/448 (94%)	419 (100%)	0	100	100
2	B	316/323 (98%)	316 (100%)	0	100	100
2	F	316/323 (98%)	316 (100%)	0	100	100
2	J	316/323 (98%)	316 (100%)	0	100	100
2	N	316/323 (98%)	316 (100%)	0	100	100
2	R	316/323 (98%)	316 (100%)	0	100	100
3	C	290/323 (90%)	290 (100%)	0	100	100
3	G	290/323 (90%)	290 (100%)	0	100	100
3	K	290/323 (90%)	290 (100%)	0	100	100
3	O	290/323 (90%)	290 (100%)	0	100	100
3	S	290/323 (90%)	290 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	5125/5470 (94%)	5125 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	216	ASN
1	A	226	ASN
1	A	354	GLN
1	A	366	GLN
1	A	434	GLN
2	B	174	ASN
2	B	260	GLN
3	C	47	GLN
3	C	87	HIS
1	E	216	ASN
1	E	226	ASN
1	E	366	GLN
1	E	434	GLN
2	F	174	ASN
2	F	260	GLN
3	G	47	GLN
3	G	87	HIS
1	I	105	ASN
1	I	216	ASN
1	I	434	GLN
2	J	174	ASN
2	J	260	GLN
3	K	47	GLN
3	K	87	HIS
1	M	216	ASN
1	M	434	GLN
2	N	66	GLN
2	N	174	ASN
2	N	260	GLN
3	O	47	GLN
3	O	87	HIS
1	Q	216	ASN
1	Q	226	ASN
1	Q	434	GLN
2	R	174	ASN

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Mol	Chain	Res	Type
2	R	260	GLN
3	S	47	GLN
3	S	87	HIS
3	S	212	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](#)

Of 15 ligands modelled in this entry, 5 are monoatomic - leaving 10 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	C14	A	501	-	13,13,13	0.22	0	12,12,12	0.20	0
6	DGA	G	401	3	10,10,43	0.23	0	12,12,45	0.28	0
4	C14	A	502	-	13,13,13	0.22	0	12,12,12	0.20	0
4	C14	M	501	-	13,13,13	0.22	0	12,12,12	0.20	0
6	DGA	C	401	3	10,10,43	0.24	0	12,12,45	0.28	0
6	DGA	S	401	3	10,10,43	0.24	0	12,12,45	0.28	0
4	C14	I	501	-	13,13,13	0.22	0	12,12,12	0.20	0
6	DGA	K	401	3	10,10,43	0.24	0	12,12,45	0.28	0
6	DGA	O	401	3	10,10,43	0.24	0	12,12,45	0.28	0
4	C14	Q	501	-	13,13,13	0.22	0	12,12,12	0.20	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	C14	A	501	-	-	6/11/11/11	-
6	DGA	G	401	3	-	0/9/9/45	-
4	C14	A	502	-	-	6/11/11/11	-
4	C14	M	501	-	-	6/11/11/11	-
6	DGA	C	401	3	-	0/9/9/45	-
6	DGA	S	401	3	-	0/9/9/45	-
4	C14	I	501	-	-	6/11/11/11	-
6	DGA	K	401	3	-	0/9/9/45	-
6	DGA	O	401	3	-	0/9/9/45	-
4	C14	Q	501	-	-	6/11/11/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	501	C14	C07-C08-C09-C10
4	I	501	C14	C07-C08-C09-C10
4	M	501	C14	C07-C08-C09-C10
4	Q	501	C14	C07-C08-C09-C10
4	A	502	C14	C07-C08-C09-C10
4	A	501	C14	C09-C10-C11-C12
4	A	502	C14	C09-C10-C11-C12
4	I	501	C14	C09-C10-C11-C12
4	M	501	C14	C09-C10-C11-C12
4	Q	501	C14	C09-C10-C11-C12
4	I	501	C14	C10-C11-C12-C13
4	A	502	C14	C10-C11-C12-C13
4	M	501	C14	C10-C11-C12-C13
4	Q	501	C14	C10-C11-C12-C13
4	A	501	C14	C10-C11-C12-C13
4	I	501	C14	C04-C05-C06-C07
4	M	501	C14	C04-C05-C06-C07
4	A	501	C14	C04-C05-C06-C07
4	A	502	C14	C04-C05-C06-C07

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Mol	Chain	Res	Type	Atoms
4	Q	501	C14	C04-C05-C06-C07
4	A	502	C14	C11-C12-C13-C14
4	M	501	C14	C11-C12-C13-C14
4	A	501	C14	C11-C12-C13-C14
4	I	501	C14	C11-C12-C13-C14
4	Q	501	C14	C11-C12-C13-C14
4	A	502	C14	C05-C06-C07-C08
4	M	501	C14	C05-C06-C07-C08
4	Q	501	C14	C05-C06-C07-C08
4	A	501	C14	C05-C06-C07-C08
4	I	501	C14	C05-C06-C07-C08

There are no ring outliers.

5 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	501	C14	3	0
4	A	502	C14	3	0
4	M	501	C14	3	0
4	I	501	C14	3	0
4	Q	501	C14	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

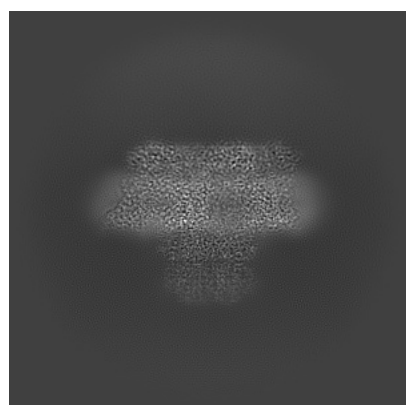
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-78051. These allow visual inspection of the internal detail of the map and identification of artifacts.

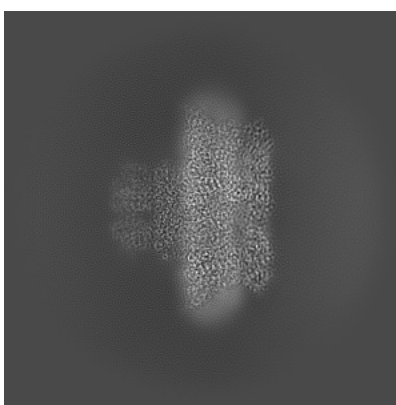
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

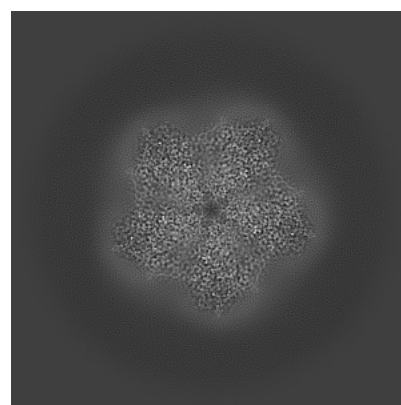
6.1.1 Primary map



X



Y

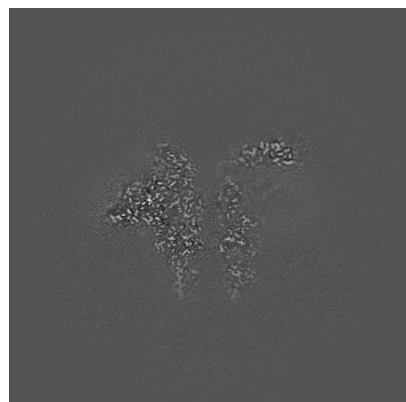


Z

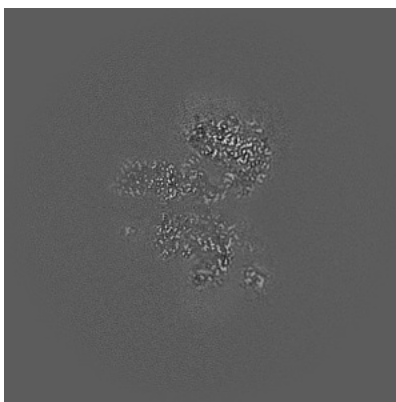
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

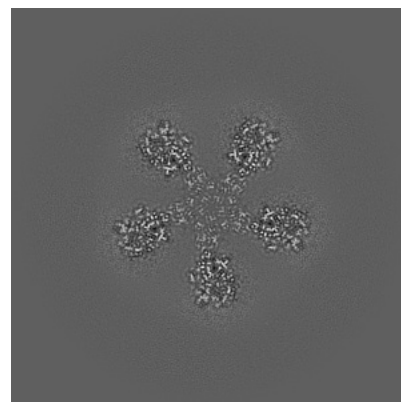
6.2.1 Primary map



X Index: 224



Y Index: 224

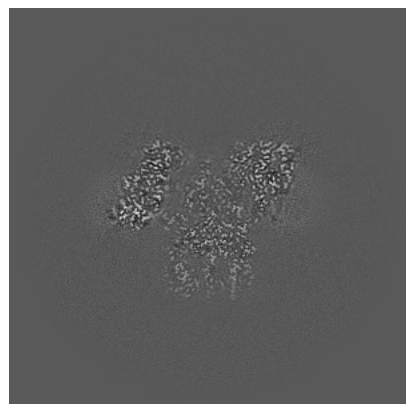


Z Index: 224

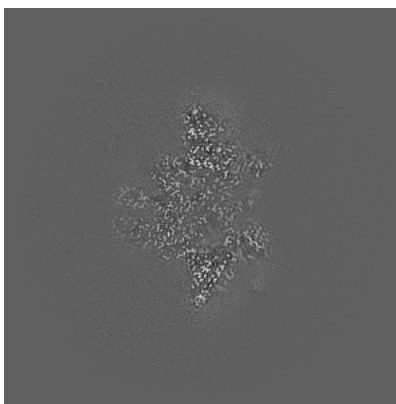
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

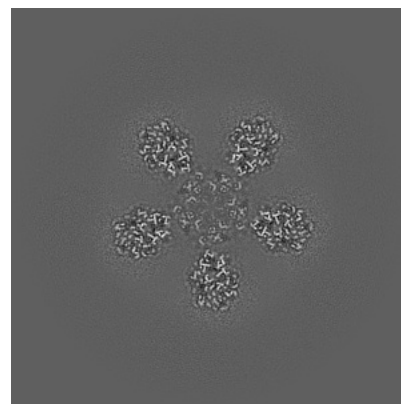
6.3.1 Primary map



X Index: 243



Y Index: 205

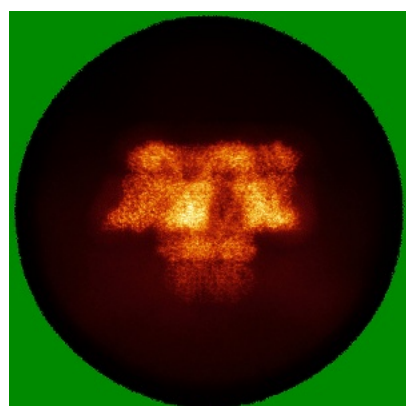


Z Index: 219

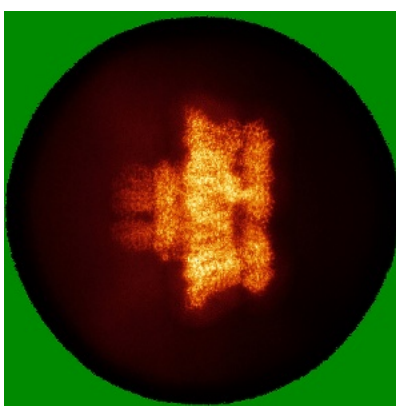
The images above show the largest variance slices of the map in three orthogonal directions.

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

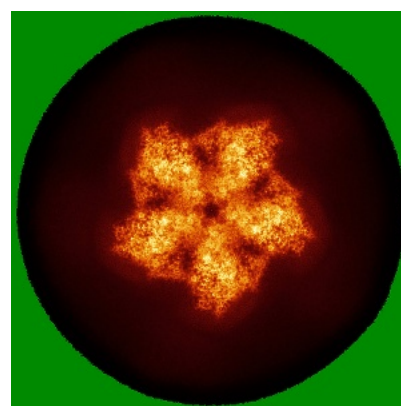
6.4.1 Primary map



X



Y

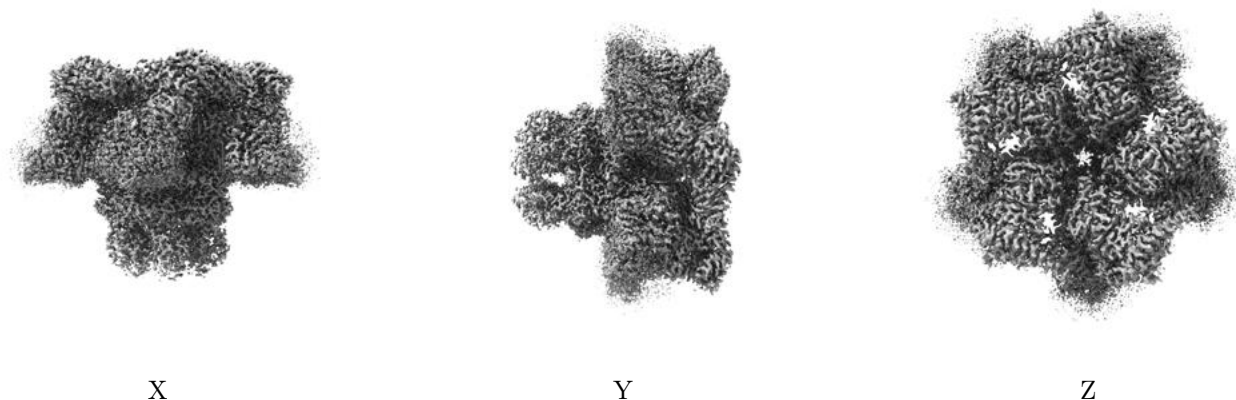


Z

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 6.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

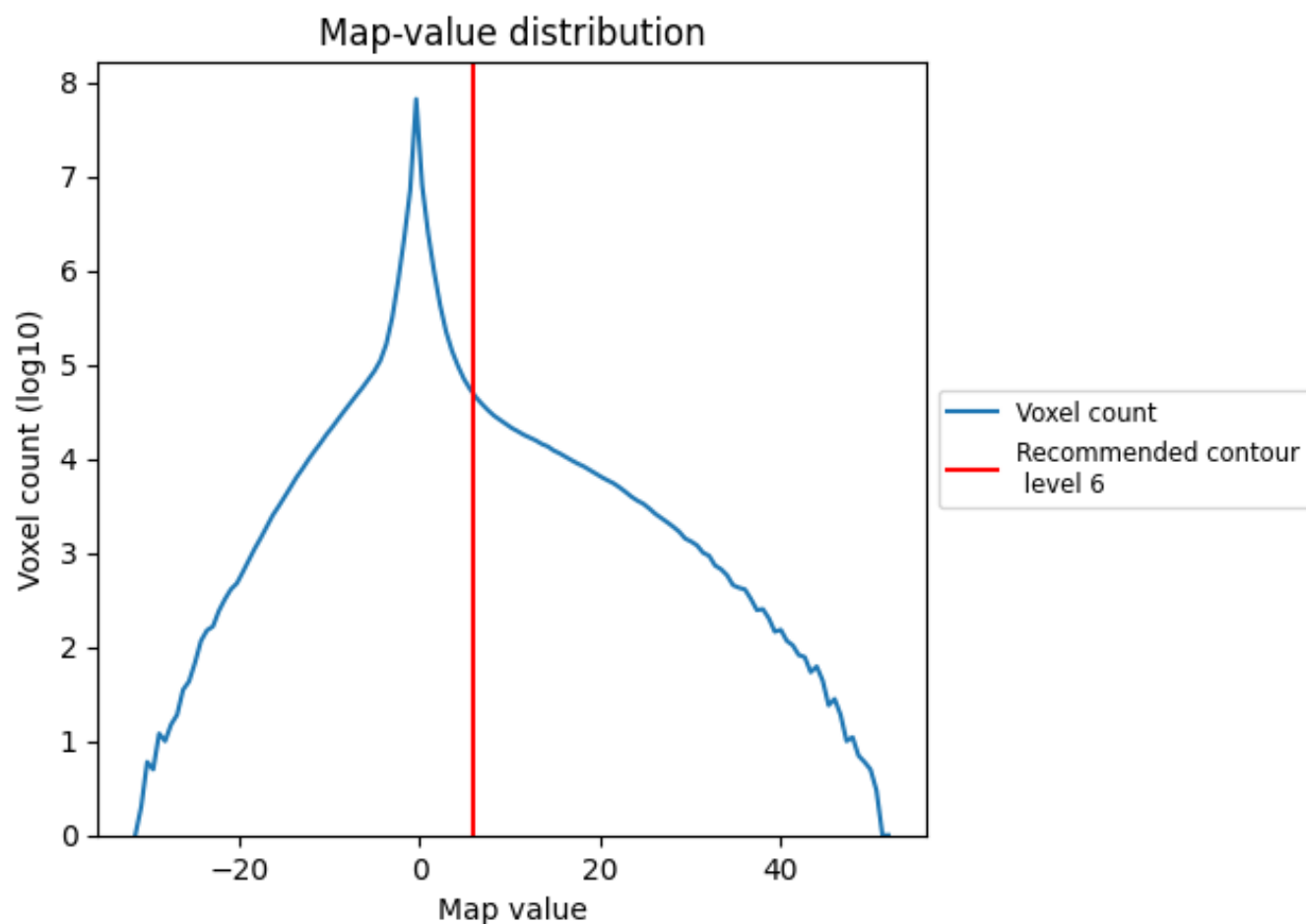
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

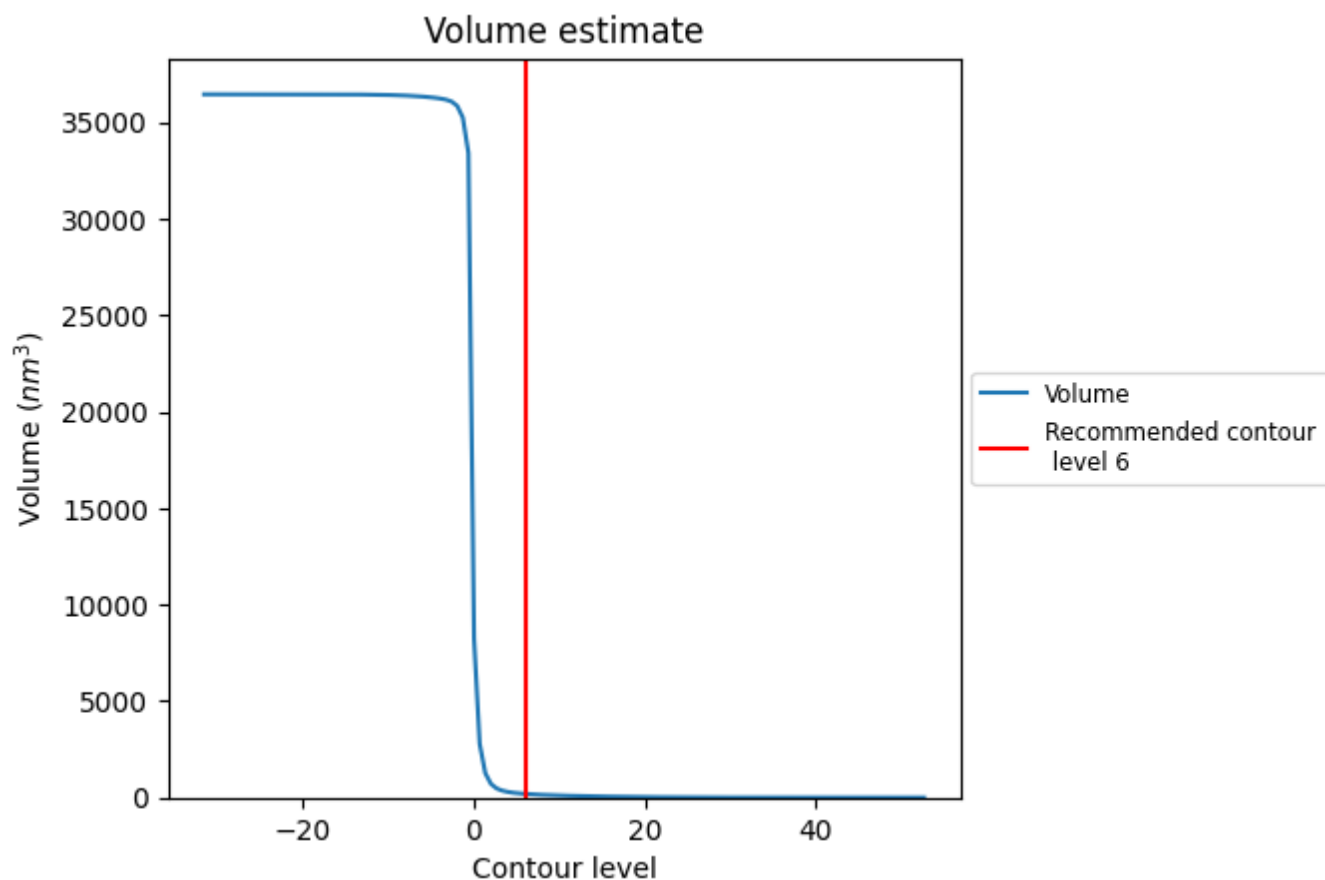
This section contains the results of statistical analysis of the map.

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.

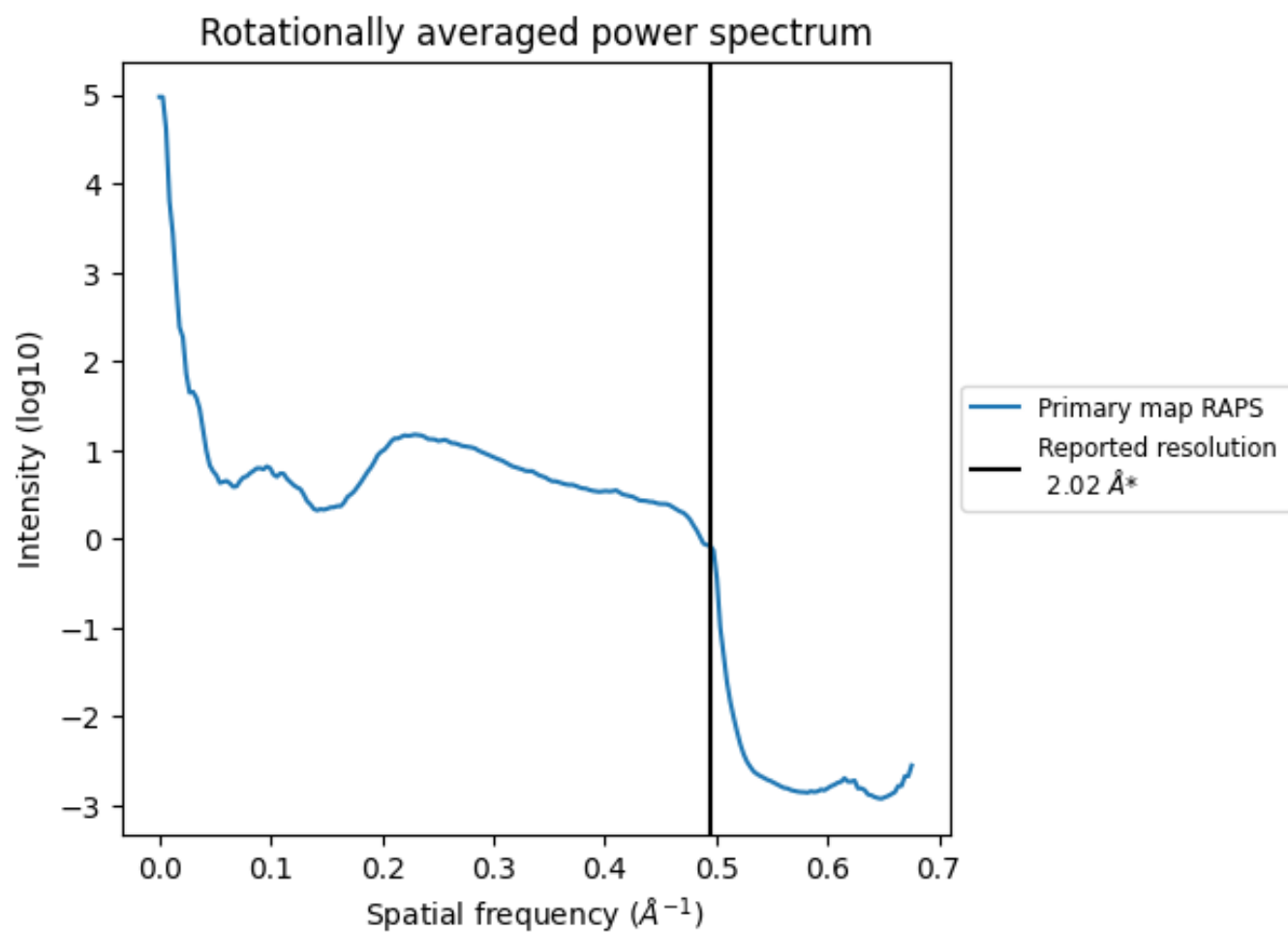
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 190 nm^3 ; this corresponds to an approximate mass of 171 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 0.495 Å⁻¹

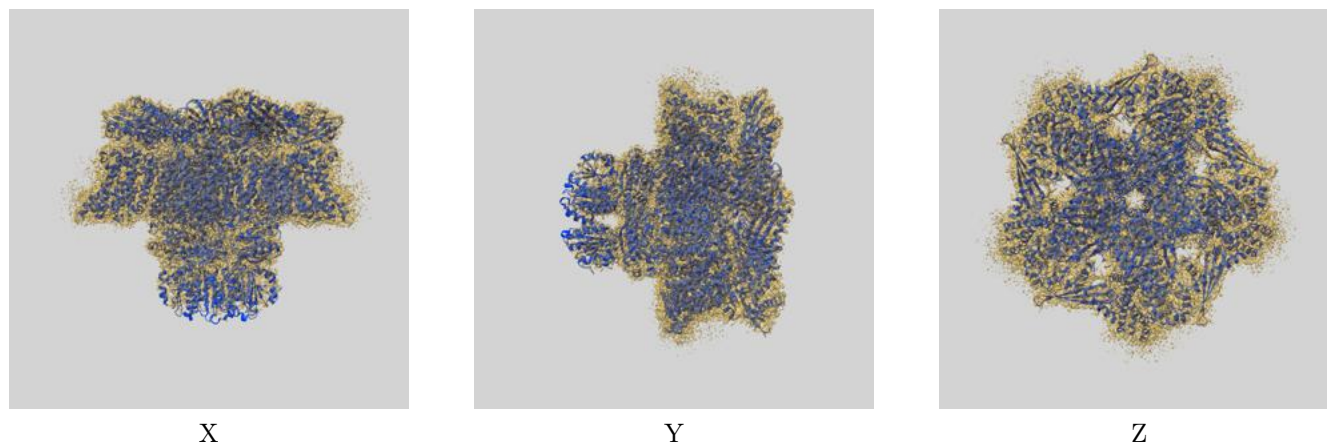
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

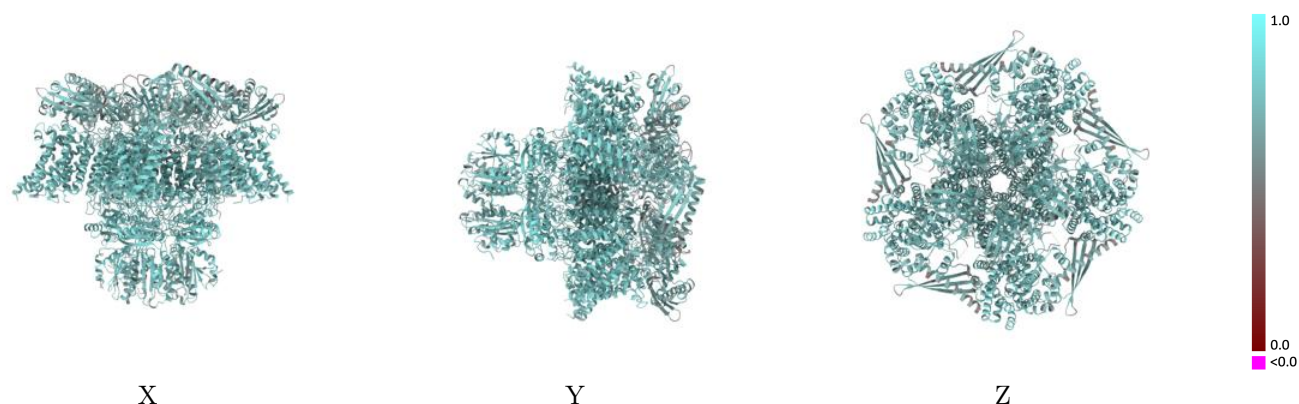
This section contains information regarding the fit between EMDB map EMD-78051 and PDB model 37BQ. Per-residue inclusion information can be found in section [3](#) on page [11](#).

9.1 Map-model overlay [i](#)



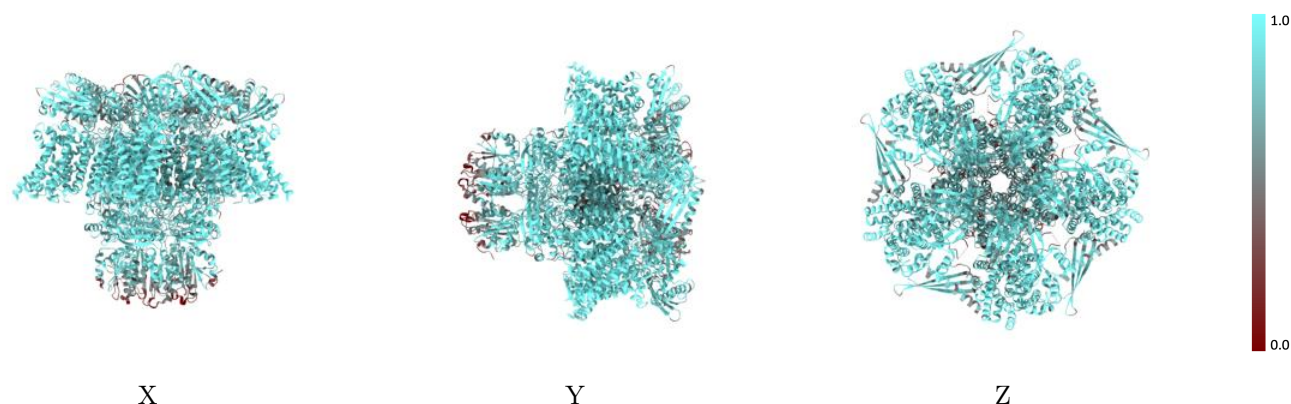
The images above show the 3D surface view of the map at the recommended contour level 6.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



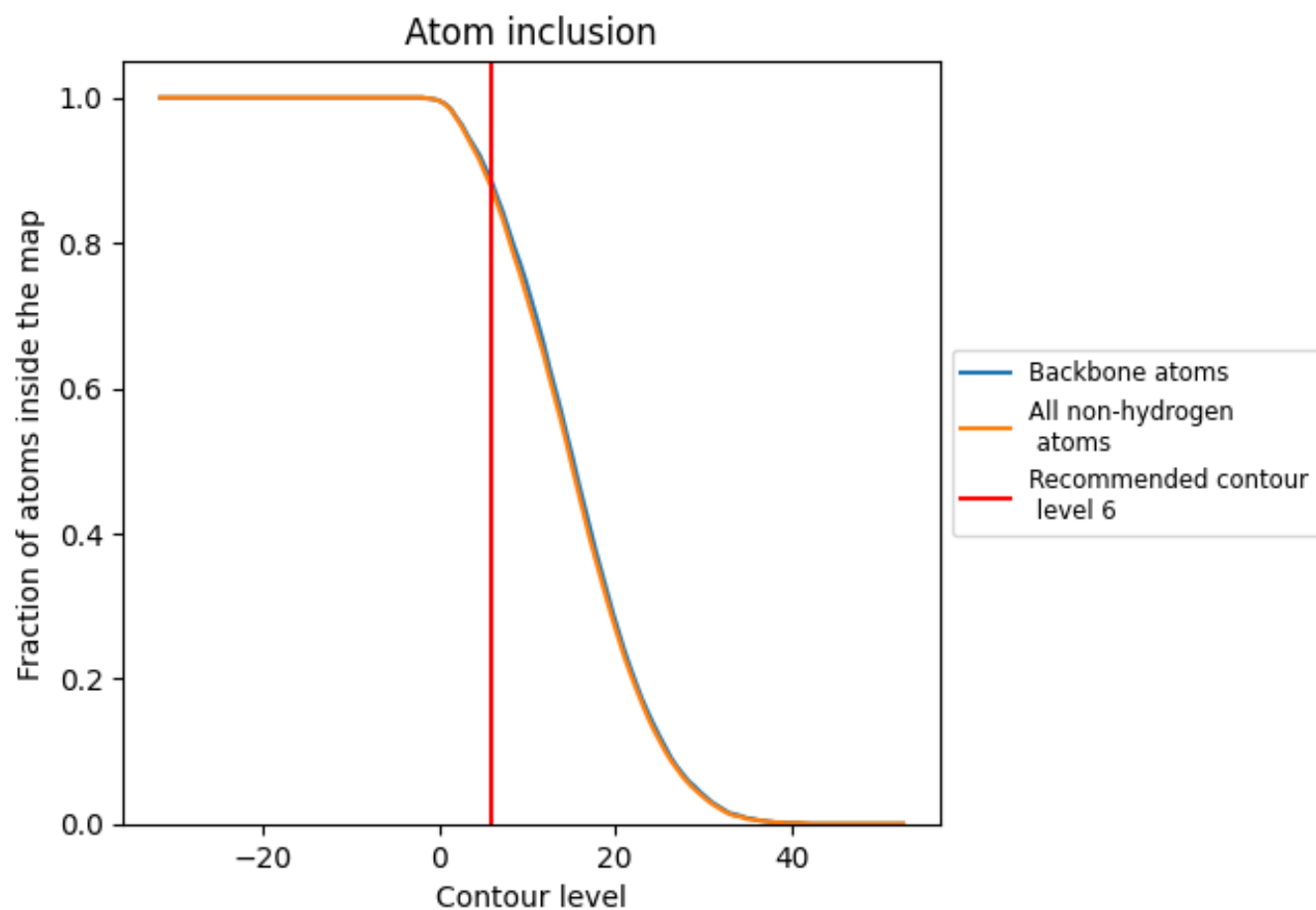
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (6).

9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 88% 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 (6) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8760	0.7230
A	0.8700	0.7600
B	0.9610	0.7490
C	0.8060	0.6460
E	0.8670	0.7590
F	0.9610	0.7460
G	0.8000	0.6460
I	0.8640	0.7580
J	0.9610	0.7460
K	0.8010	0.6460
M	0.8660	0.7590
N	0.9610	0.7470
O	0.8030	0.6460
Q	0.8660	0.7580
R	0.9600	0.7470
S	0.8030	0.6450

