wwPDB EM Validation Summary Report

Oct 28, 2026 – 02:40 PM JST

PDB ID	:	9WDP / pdb_00009wdp
EMDB ID	:	EMD-65894
Title	:	Cyro-EM structure of prefusion RSV fusion glycoprotein in complex with Ziresovir and motavizumab Fab
Authors	:	Zhang, W.; Yan, M.R.; Zou, J.J.; Peng, W.
Deposited on	:	2025-08-19
Resolution	:	3.27 Å(reported)

This is a wwPDB EM Validation Summary 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:

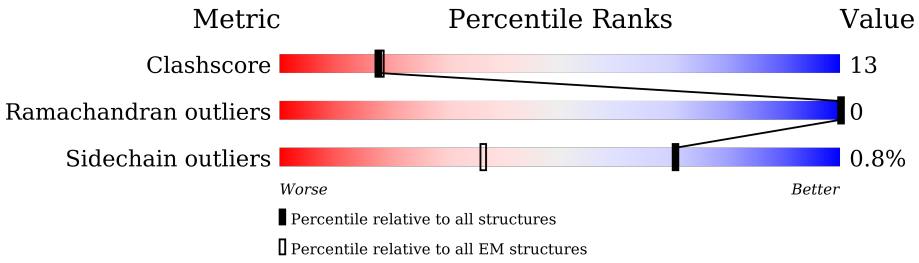
EMDB validation analysis	:	0.0.1.dev133
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4-5-2 with Phenix2.2
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using 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)	:	2.51

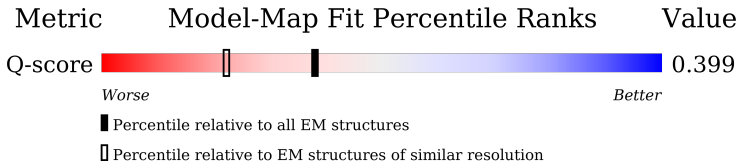
1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.27 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.





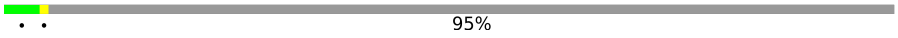
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14508 (2.77 - 3.77)

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 <=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	593	48% 22% 30%
1	B	593	48% 22% 30%
1	C	593	46% 24% 30%
2	D	225	20% 78%

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Mol	Chain	Length	Quality of chain
2	F	225	
2	H	225	
3	E	213	
3	G	213	
3	L	213	

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 11338 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 Fusion glycoprotein F0,Fibritin.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	415	Total	C	N	O	S	0	0
			3210	2037	524	628	21		
1	B	415	Total	C	N	O	S	0	0
			3218	2042	525	630	21		
1	C	415	Total	C	N	O	S	0	0
			3214	2040	524	629	21		

There are 177 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	ALA	PRO	variant	UNP P03420
A	155	CYS	SER	conflict	UNP P03420
A	190	PHE	SER	conflict	UNP P03420
A	207	LEU	VAL	conflict	UNP P03420
A	290	CYS	SER	conflict	UNP P03420
A	379	VAL	ILE	variant	UNP P03420
A	447	VAL	MET	variant	UNP P03420
A	514	GLY	-	linker	UNP P03420
A	515	SER	-	linker	UNP P03420
A	544	GLY	-	expression tag	UNP P10104
A	545	LEU	-	expression tag	UNP P10104
A	546	GLU	-	expression tag	UNP P10104
A	547	VAL	-	expression tag	UNP P10104
A	548	LEU	-	expression tag	UNP P10104
A	549	PHE	-	expression tag	UNP P10104
A	550	GLN	-	expression tag	UNP P10104
A	551	GLY	-	expression tag	UNP P10104
A	552	PRO	-	expression tag	UNP P10104
A	553	GLY	-	expression tag	UNP P10104
A	554	GLY	-	expression tag	UNP P10104
A	555	TRP	-	expression tag	UNP P10104
A	556	SER	-	expression tag	UNP P10104
A	557	HIS	-	expression tag	UNP P10104
A	558	PRO	-	expression tag	UNP P10104

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Chain	Residue	Modelled	Actual	Comment	Reference
A	559	GLN	-	expression tag	UNP P10104
A	560	PHE	-	expression tag	UNP P10104
A	561	GLU	-	expression tag	UNP P10104
A	562	LYS	-	expression tag	UNP P10104
A	563	GLY	-	expression tag	UNP P10104
A	564	GLY	-	expression tag	UNP P10104
A	565	GLY	-	expression tag	UNP P10104
A	566	SER	-	expression tag	UNP P10104
A	567	GLY	-	expression tag	UNP P10104
A	568	GLY	-	expression tag	UNP P10104
A	569	GLY	-	expression tag	UNP P10104
A	570	SER	-	expression tag	UNP P10104
A	571	GLY	-	expression tag	UNP P10104
A	572	GLY	-	expression tag	UNP P10104
A	573	SER	-	expression tag	UNP P10104
A	574	ALA	-	expression tag	UNP P10104
A	575	TRP	-	expression tag	UNP P10104
A	576	SER	-	expression tag	UNP P10104
A	577	HIS	-	expression tag	UNP P10104
A	578	PRO	-	expression tag	UNP P10104
A	579	GLN	-	expression tag	UNP P10104
A	580	PHE	-	expression tag	UNP P10104
A	581	GLU	-	expression tag	UNP P10104
A	582	LYS	-	expression tag	UNP P10104
A	583	GLY	-	expression tag	UNP P10104
A	584	GLY	-	expression tag	UNP P10104
A	585	SER	-	expression tag	UNP P10104
A	586	GLY	-	expression tag	UNP P10104
A	587	PRO	-	expression tag	UNP P10104
A	588	HIS	-	expression tag	UNP P10104
A	589	HIS	-	expression tag	UNP P10104
A	590	HIS	-	expression tag	UNP P10104
A	591	HIS	-	expression tag	UNP P10104
A	592	HIS	-	expression tag	UNP P10104
A	593	HIS	-	expression tag	UNP P10104
B	102	ALA	PRO	variant	UNP P03420
B	155	CYS	SER	conflict	UNP P03420
B	190	PHE	SER	conflict	UNP P03420
B	207	LEU	VAL	conflict	UNP P03420
B	290	CYS	SER	conflict	UNP P03420
B	379	VAL	ILE	variant	UNP P03420
B	447	VAL	MET	variant	UNP P03420

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Chain	Residue	Modelled	Actual	Comment	Reference
B	514	GLY	-	linker	UNP P03420
B	515	SER	-	linker	UNP P03420
B	544	GLY	-	expression tag	UNP P10104
B	545	LEU	-	expression tag	UNP P10104
B	546	GLU	-	expression tag	UNP P10104
B	547	VAL	-	expression tag	UNP P10104
B	548	LEU	-	expression tag	UNP P10104
B	549	PHE	-	expression tag	UNP P10104
B	550	GLN	-	expression tag	UNP P10104
B	551	GLY	-	expression tag	UNP P10104
B	552	PRO	-	expression tag	UNP P10104
B	553	GLY	-	expression tag	UNP P10104
B	554	GLY	-	expression tag	UNP P10104
B	555	TRP	-	expression tag	UNP P10104
B	556	SER	-	expression tag	UNP P10104
B	557	HIS	-	expression tag	UNP P10104
B	558	PRO	-	expression tag	UNP P10104
B	559	GLN	-	expression tag	UNP P10104
B	560	PHE	-	expression tag	UNP P10104
B	561	GLU	-	expression tag	UNP P10104
B	562	LYS	-	expression tag	UNP P10104
B	563	GLY	-	expression tag	UNP P10104
B	564	GLY	-	expression tag	UNP P10104
B	565	GLY	-	expression tag	UNP P10104
B	566	SER	-	expression tag	UNP P10104
B	567	GLY	-	expression tag	UNP P10104
B	568	GLY	-	expression tag	UNP P10104
B	569	GLY	-	expression tag	UNP P10104
B	570	SER	-	expression tag	UNP P10104
B	571	GLY	-	expression tag	UNP P10104
B	572	GLY	-	expression tag	UNP P10104
B	573	SER	-	expression tag	UNP P10104
B	574	ALA	-	expression tag	UNP P10104
B	575	TRP	-	expression tag	UNP P10104
B	576	SER	-	expression tag	UNP P10104
B	577	HIS	-	expression tag	UNP P10104
B	578	PRO	-	expression tag	UNP P10104
B	579	GLN	-	expression tag	UNP P10104
B	580	PHE	-	expression tag	UNP P10104
B	581	GLU	-	expression tag	UNP P10104
B	582	LYS	-	expression tag	UNP P10104
B	583	GLY	-	expression tag	UNP P10104

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Chain	Residue	Modelled	Actual	Comment	Reference
B	584	GLY	-	expression tag	UNP P10104
B	585	SER	-	expression tag	UNP P10104
B	586	GLY	-	expression tag	UNP P10104
B	587	PRO	-	expression tag	UNP P10104
B	588	HIS	-	expression tag	UNP P10104
B	589	HIS	-	expression tag	UNP P10104
B	590	HIS	-	expression tag	UNP P10104
B	591	HIS	-	expression tag	UNP P10104
B	592	HIS	-	expression tag	UNP P10104
B	593	HIS	-	expression tag	UNP P10104
C	102	ALA	PRO	variant	UNP P03420
C	155	CYS	SER	conflict	UNP P03420
C	190	PHE	SER	conflict	UNP P03420
C	207	LEU	VAL	conflict	UNP P03420
C	290	CYS	SER	conflict	UNP P03420
C	379	VAL	ILE	variant	UNP P03420
C	447	VAL	MET	variant	UNP P03420
C	514	GLY	-	linker	UNP P03420
C	515	SER	-	linker	UNP P03420
C	544	GLY	-	expression tag	UNP P10104
C	545	LEU	-	expression tag	UNP P10104
C	546	GLU	-	expression tag	UNP P10104
C	547	VAL	-	expression tag	UNP P10104
C	548	LEU	-	expression tag	UNP P10104
C	549	PHE	-	expression tag	UNP P10104
C	550	GLN	-	expression tag	UNP P10104
C	551	GLY	-	expression tag	UNP P10104
C	552	PRO	-	expression tag	UNP P10104
C	553	GLY	-	expression tag	UNP P10104
C	554	GLY	-	expression tag	UNP P10104
C	555	TRP	-	expression tag	UNP P10104
C	556	SER	-	expression tag	UNP P10104
C	557	HIS	-	expression tag	UNP P10104
C	558	PRO	-	expression tag	UNP P10104
C	559	GLN	-	expression tag	UNP P10104
C	560	PHE	-	expression tag	UNP P10104
C	561	GLU	-	expression tag	UNP P10104
C	562	LYS	-	expression tag	UNP P10104
C	563	GLY	-	expression tag	UNP P10104
C	564	GLY	-	expression tag	UNP P10104
C	565	GLY	-	expression tag	UNP P10104
C	566	SER	-	expression tag	UNP P10104

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Chain	Residue	Modelled	Actual	Comment	Reference
C	567	GLY	-	expression tag	UNP P10104
C	568	GLY	-	expression tag	UNP P10104
C	569	GLY	-	expression tag	UNP P10104
C	570	SER	-	expression tag	UNP P10104
C	571	GLY	-	expression tag	UNP P10104
C	572	GLY	-	expression tag	UNP P10104
C	573	SER	-	expression tag	UNP P10104
C	574	ALA	-	expression tag	UNP P10104
C	575	TRP	-	expression tag	UNP P10104
C	576	SER	-	expression tag	UNP P10104
C	577	HIS	-	expression tag	UNP P10104
C	578	PRO	-	expression tag	UNP P10104
C	579	GLN	-	expression tag	UNP P10104
C	580	PHE	-	expression tag	UNP P10104
C	581	GLU	-	expression tag	UNP P10104
C	582	LYS	-	expression tag	UNP P10104
C	583	GLY	-	expression tag	UNP P10104
C	584	GLY	-	expression tag	UNP P10104
C	585	SER	-	expression tag	UNP P10104
C	586	GLY	-	expression tag	UNP P10104
C	587	PRO	-	expression tag	UNP P10104
C	588	HIS	-	expression tag	UNP P10104
C	589	HIS	-	expression tag	UNP P10104
C	590	HIS	-	expression tag	UNP P10104
C	591	HIS	-	expression tag	UNP P10104
C	592	HIS	-	expression tag	UNP P10104
C	593	HIS	-	expression tag	UNP P10104

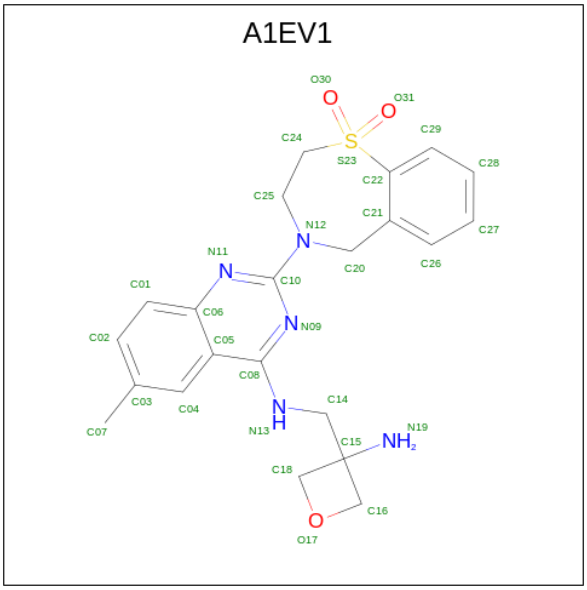
- Molecule 2 is a protein called Motavizumab Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	50	Total	C	N	O	S	0	0
			366	234	63	66	3		
2	F	50	Total	C	N	O	S	0	0
			432	288	70	71	3		
2	H	50	Total	C	N	O	S	0	0
			408	268	70	68	2		

- Molecule 3 is a protein called Motavizumab Fab light chain.

Mol	Chain	Residues	Atoms	AltConf	Trace
3	E	30	Total C N O S 218 143 33 40 2	0	0
3	G	10	Total C N O S 91 64 14 12 1	0	0
3	L	19	Total C N O 150 101 23 26	0	0

- Molecule 4 is Ziresovir (CCD ID: A1EV1) (formula: $C_{22}H_{25}N_5O_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	AltConf
4	B	1	Total C N O S 31 22 5 3 1	0

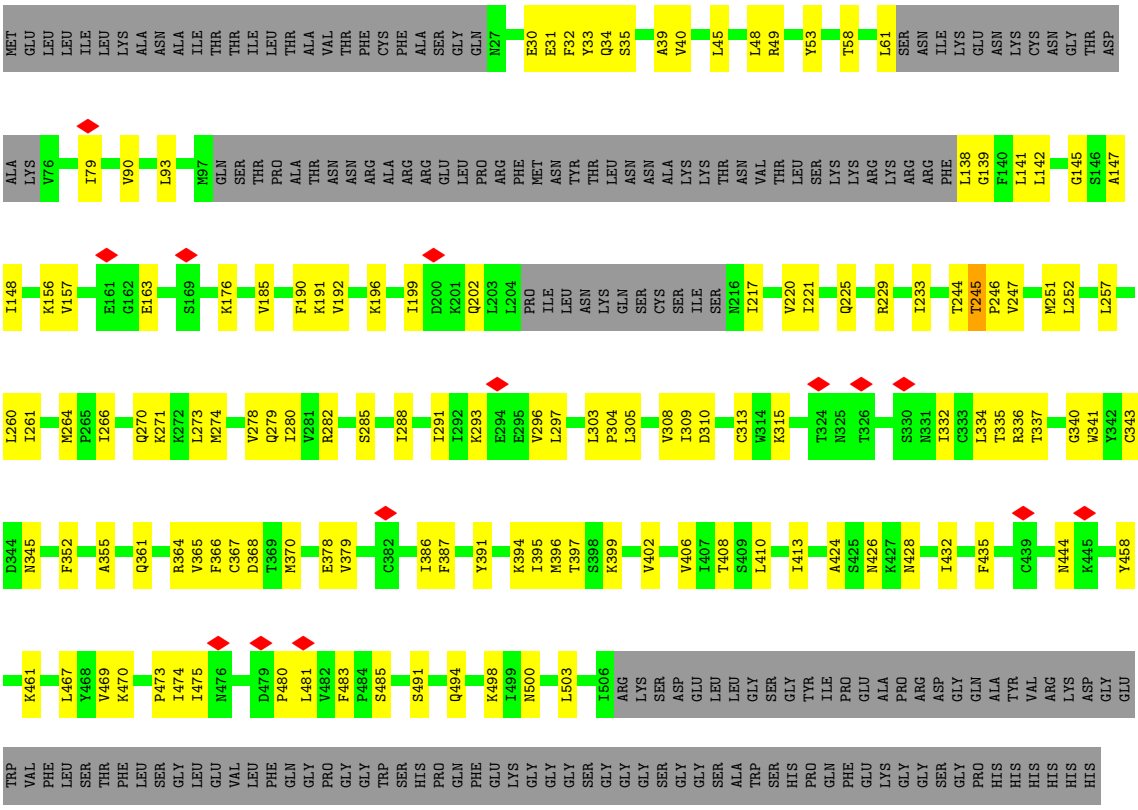
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and 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: Fusion glycoprotein F0,Fibritin

Chain A:

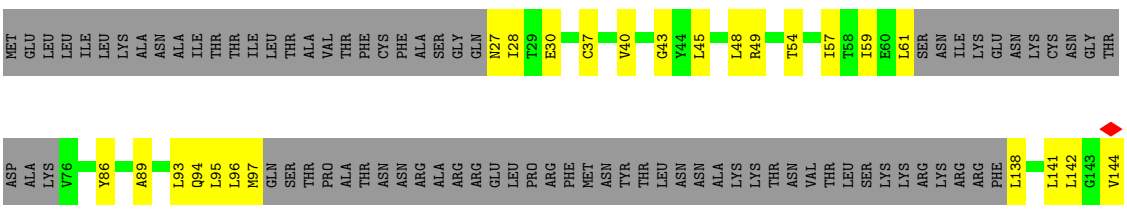


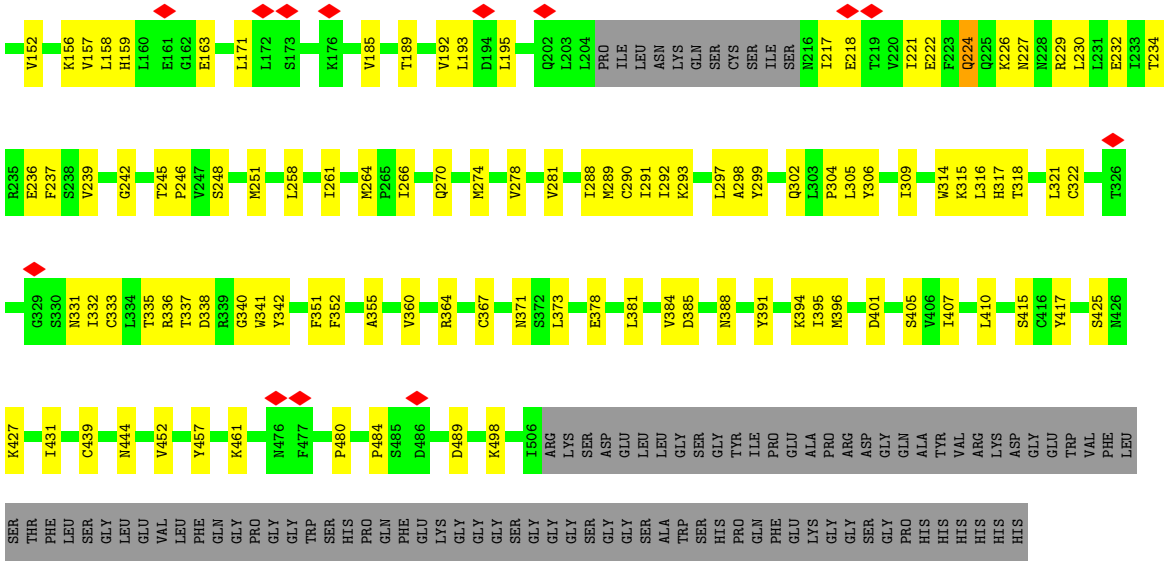


• Molecule 1: Fusion glycoprotein F0,Fibritin

Chain B:



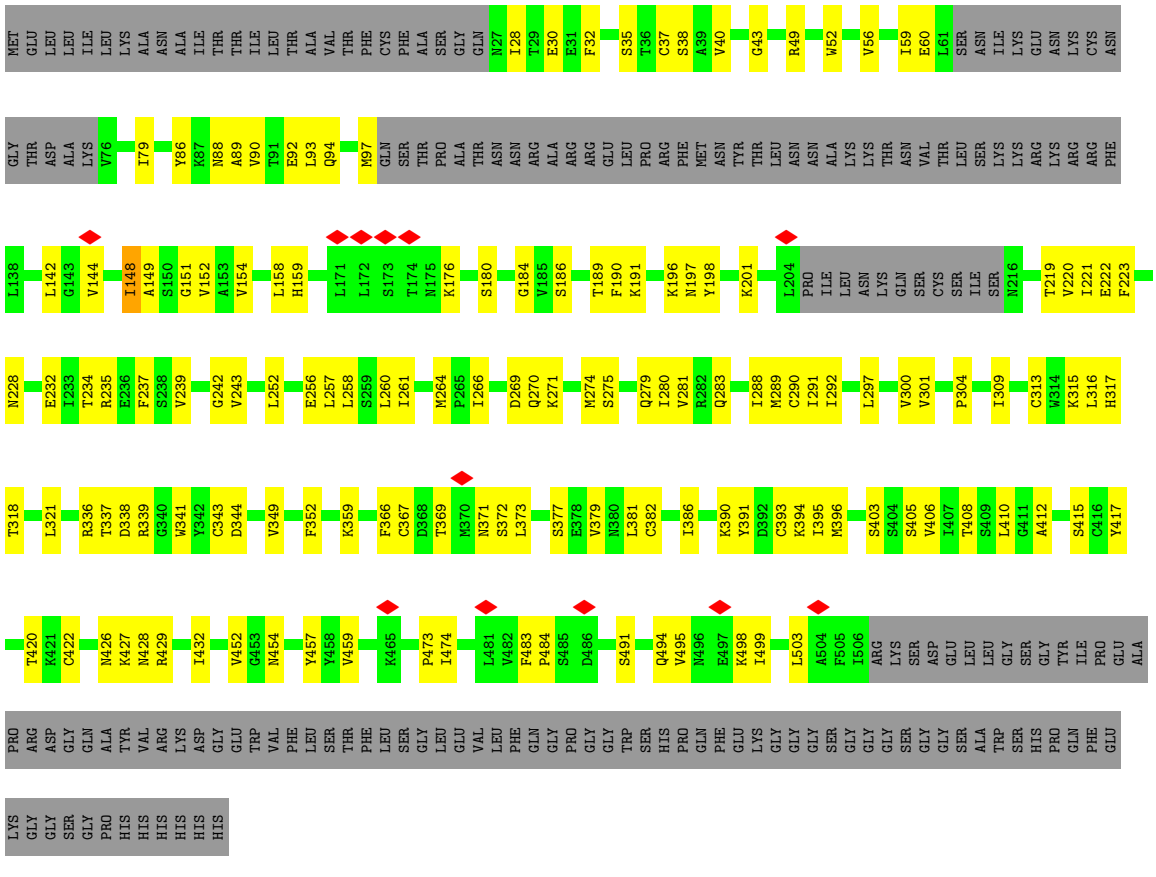




• Molecule 1: Fusion glycoprotein F0,Fibritin

Chain C:

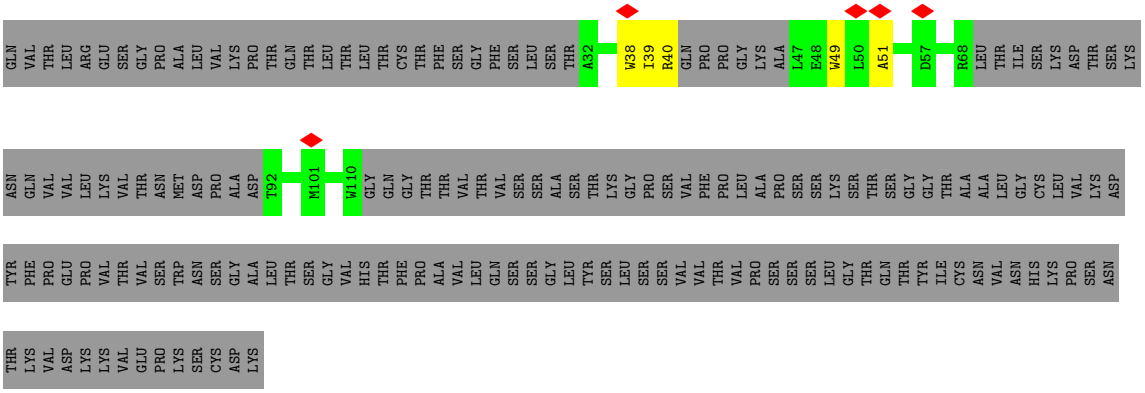




• Molecule 2: Motavizumab Fab heavy chain

Chain D:

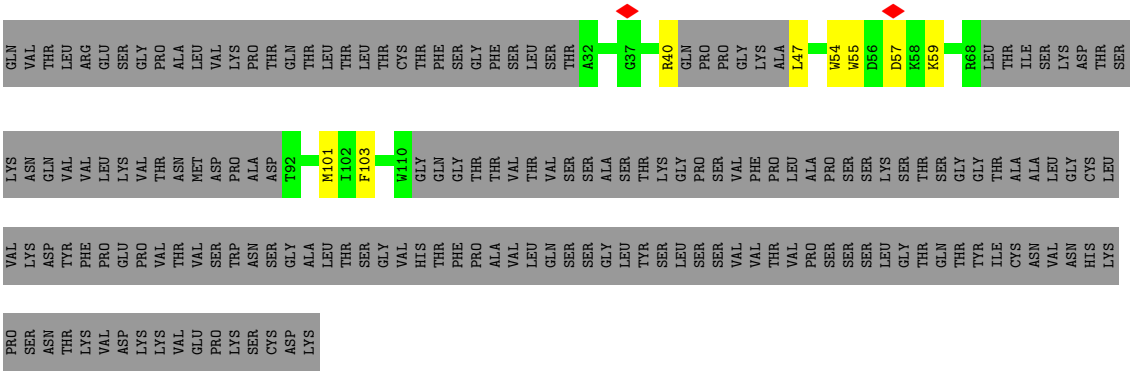




• Molecule 2: Motavizumab Fab heavy chain

Chain F:

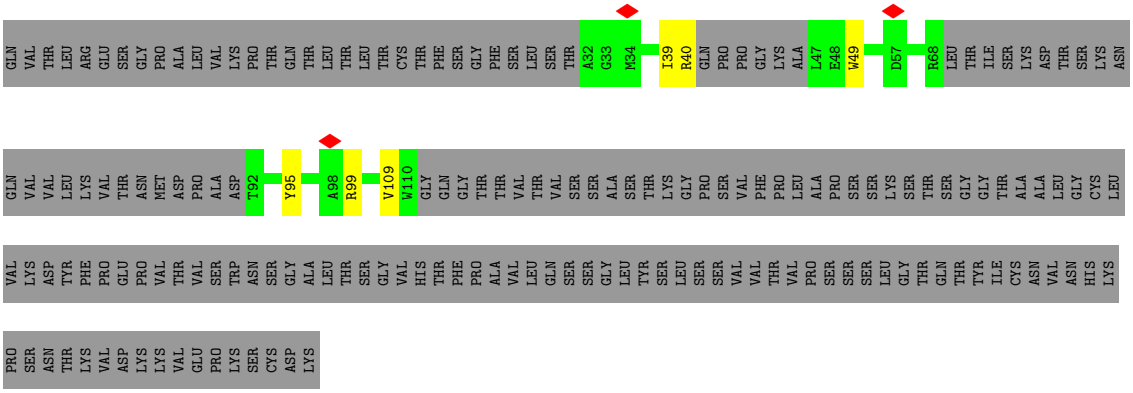




• Molecule 2: Motavizumab Fab heavy chain

Chain H:

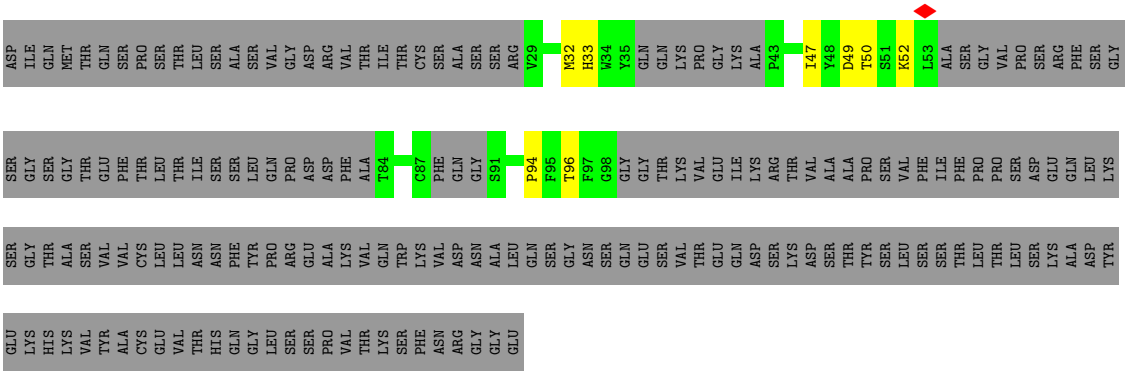




• Molecule 3: Motavizumab Fab light chain

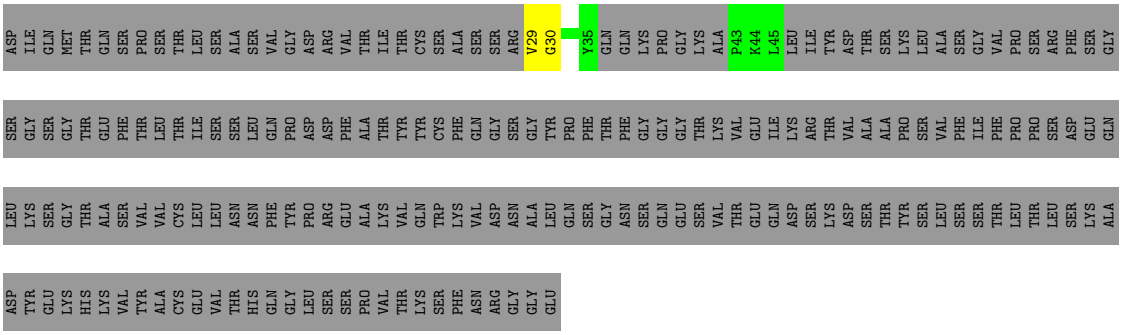
Chain E:





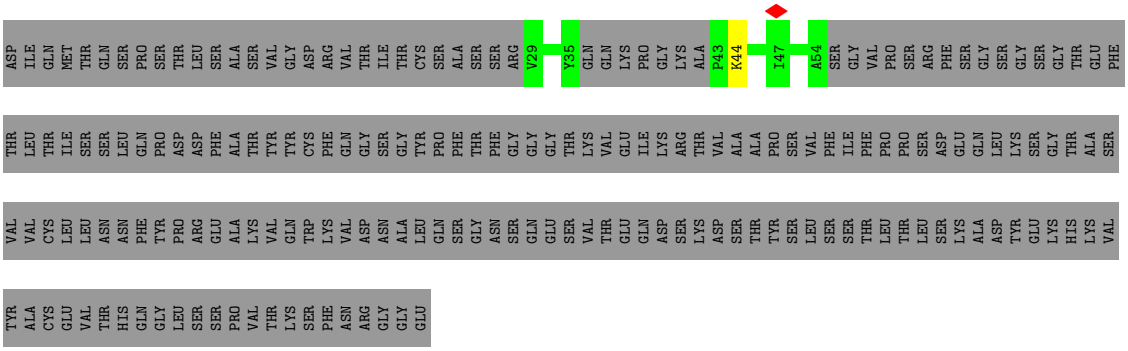
• Molecule 3: Motavizumab Fab light chain

Chain G: 95%



• Molecule 3: Motavizumab Fab light chain

Chain L: 8% 91%



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	81339	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 ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	0.346	Depositor
Minimum map value	-0.191	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.012	Depositor
Recommended contour level	0.054	Depositor
Map size (Å)	262.80002, 262.80002, 262.80002	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.7300001, 0.7300001, 0.7300001	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: A1EV1

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	0.15	0/3256	0.34	0/4414
1	B	0.12	0/3264	0.34	0/4423
1	C	0.14	0/3260	0.37	0/4418
2	D	0.24	0/377	0.34	0/516
2	F	0.11	0/448	0.34	0/607
2	H	0.09	0/422	0.30	0/574
3	E	0.60	0/224	0.76	0/302
3	G	0.19	0/95	0.35	0/126
3	L	0.12	0/155	0.38	0/210
All	All	0.16	0/11501	0.36	0/15590

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

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	3210	0	3241	92	0
1	B	3218	0	3256	101	0
1	C	3214	0	3250	109	0
2	D	366	0	274	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	432	0	393	6	0
2	H	408	0	354	3	0
3	E	218	0	167	3	0
3	G	91	0	86	1	0
3	L	150	0	128	0	0
4	B	31	0	0	0	0
All	All	11338	0	11149	303	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 13.

The worst 5 of 303 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:335:THR:HB	1:B:396:MET:HB2	1.60	0.82
1:A:313:CYS:HB2	1:A:343:CYS:HA	1.62	0.81
1:C:336:ARG:HA	1:C:395:ILE:HA	1.61	0.80
1:B:336:ARG:HA	1:B:395:ILE:HA	1.63	0.78
1:C:344:ASP:HB3	1:C:349:VAL:HG22	1.67	0.77

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	407/593 (69%)	382 (94%)	25 (6%)	0	100	100
1	B	407/593 (69%)	384 (94%)	23 (6%)	0	100	100
1	C	407/593 (69%)	396 (97%)	11 (3%)	0	100	100
2	D	44/225 (20%)	42 (96%)	2 (4%)	0	100	100
2	F	44/225 (20%)	39 (89%)	5 (11%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	44/225 (20%)	41 (93%)	3 (7%)	0	100	100
3	E	22/213 (10%)	20 (91%)	2 (9%)	0	100	100
3	G	6/213 (3%)	6 (100%)	0	0	100	100
3	L	15/213 (7%)	13 (87%)	2 (13%)	0	100	100
All	All	1396/3093 (45%)	1323 (95%)	73 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	376/524 (72%)	374 (100%)	2 (0%)	81	83
1	B	378/524 (72%)	377 (100%)	1 (0%)	86	86
1	C	377/524 (72%)	376 (100%)	1 (0%)	86	86
2	D	27/197 (14%)	25 (93%)	2 (7%)	13	39
2	F	42/197 (21%)	42 (100%)	0	100	100
2	H	37/197 (19%)	37 (100%)	0	100	100
3	E	19/184 (10%)	16 (84%)	3 (16%)	2	12
3	G	9/184 (5%)	9 (100%)	0	100	100
3	L	13/184 (7%)	12 (92%)	1 (8%)	12	37
All	All	1278/2715 (47%)	1268 (99%)	10 (1%)	70	79

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	E	50	THR
3	E	96	THR
3	L	44	LYS
1	C	148	ILE
2	D	39	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	462	GLN
1	C	165	ASN
1	C	225	GLN
1	B	361	GLN
1	B	388	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is modelled in this entry.

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	A1EV1	B	601	-	30,35,35	1.90	4 (13%)	29,53,53	2.19	11 (37%)

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	A1EV1	B	601	-	-	2/7/34/34	0/4/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	601	A1EV1	C08-N13	6.30	1.46	1.34
4	B	601	A1EV1	C10-N12	5.22	1.45	1.35
4	B	601	A1EV1	C05-C06	-3.24	1.37	1.42
4	B	601	A1EV1	C20-C21	2.26	1.54	1.51

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	601	A1EV1	C08-C05-C06	5.60	119.40	115.88
4	B	601	A1EV1	C04-C05-C08	-4.77	120.58	124.88
4	B	601	A1EV1	C05-C06-N11	-3.89	118.68	122.81
4	B	601	A1EV1	O31-S23-C22	-3.15	105.42	108.89
4	B	601	A1EV1	O30-S23-C22	-3.14	105.43	108.89

There are no chirality outliers.

All (2) torsion outliers are listed below:

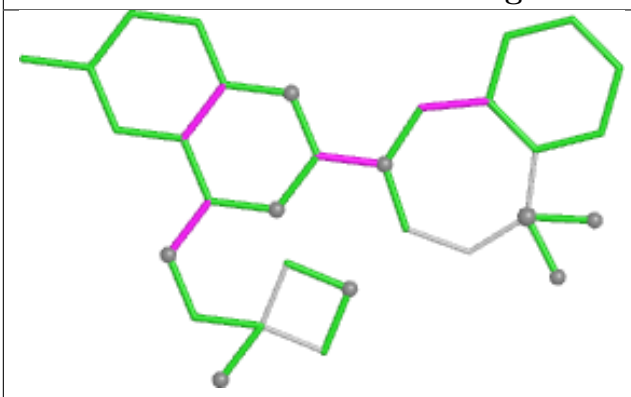
Mol	Chain	Res	Type	Atoms
4	B	601	A1EV1	N09-C08-N13-C14
4	B	601	A1EV1	C05-C08-N13-C14

There are no ring outliers.

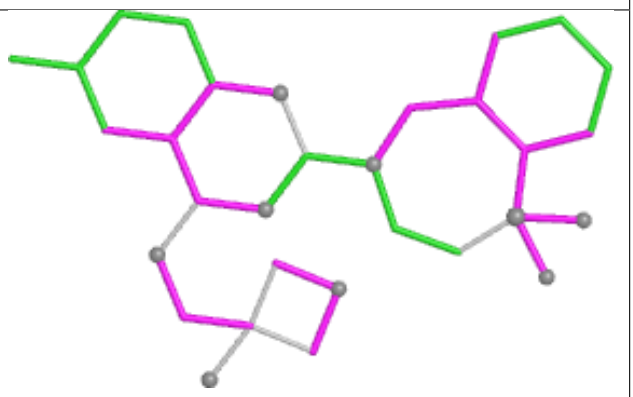
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

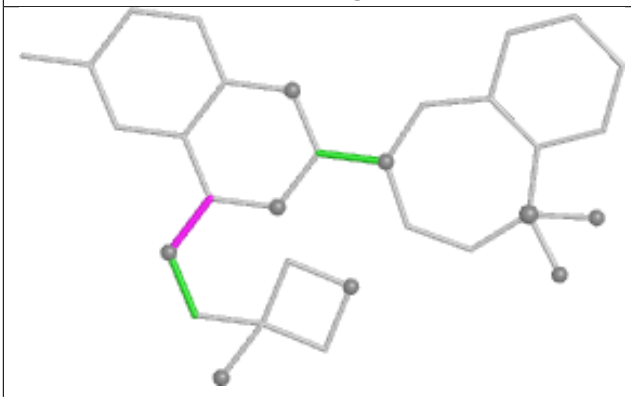
Ligand A1EV1 B 601



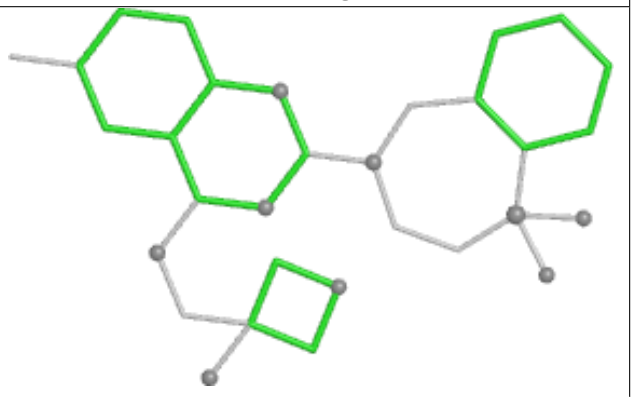
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

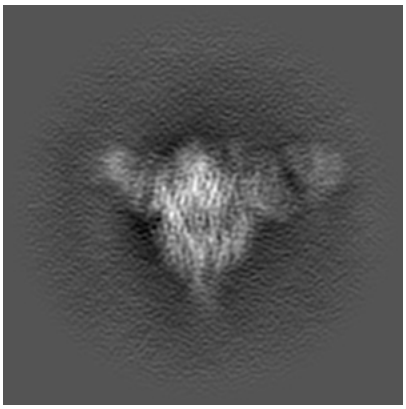
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-65894. These allow visual inspection of the internal detail of the map and identification of artifacts.

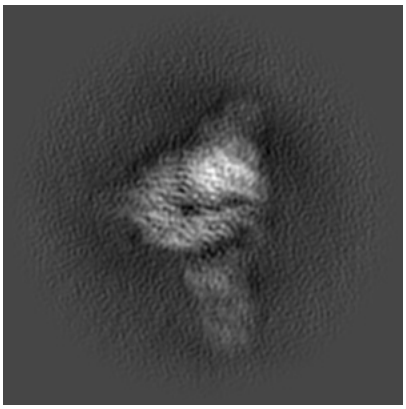
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

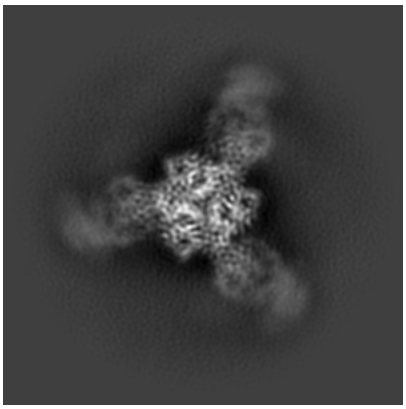
6.1.1 Primary map



X

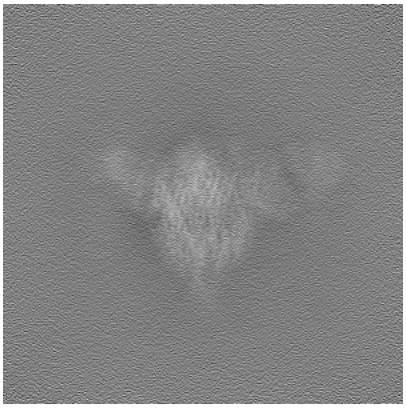


Y

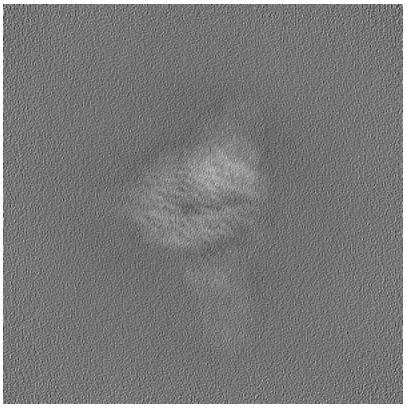


Z

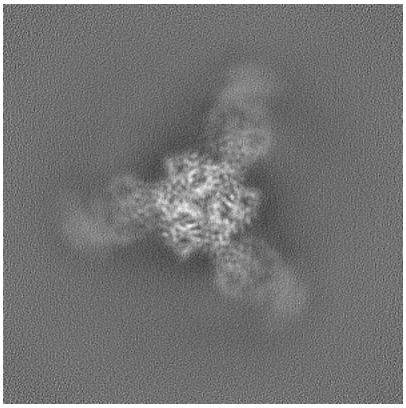
6.1.2 Raw map



X



Y

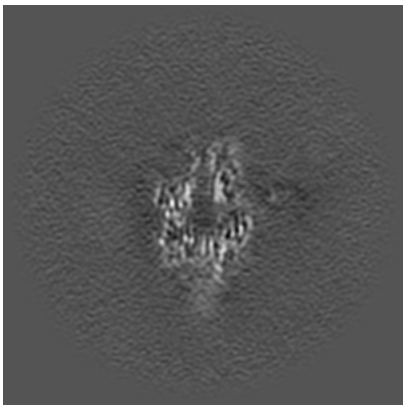


Z

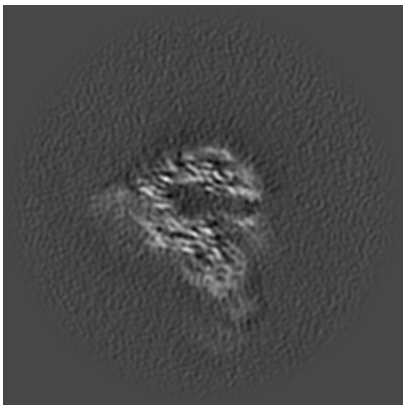
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

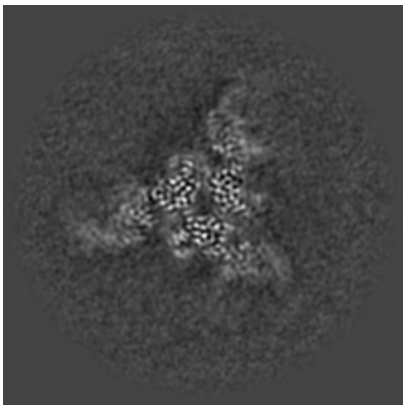
6.2.1 Primary map



X Index: 180

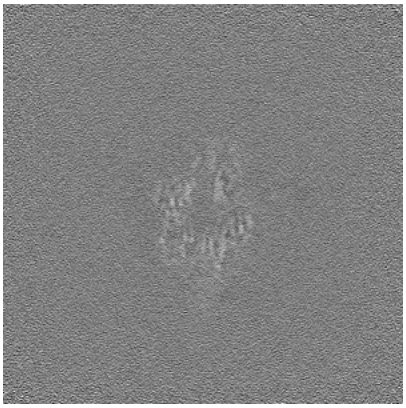


Y Index: 180

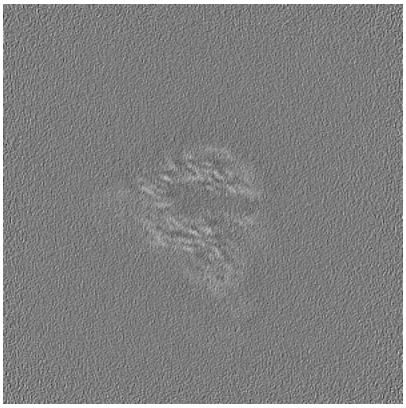


Z Index: 180

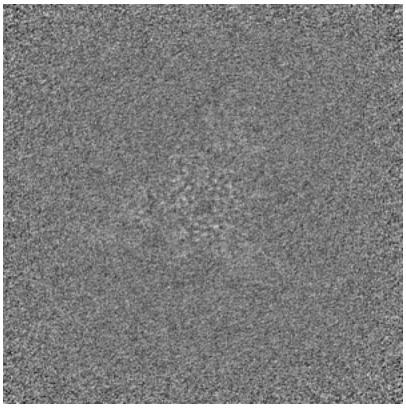
6.2.2 Raw map



X Index: 180



Y Index: 180

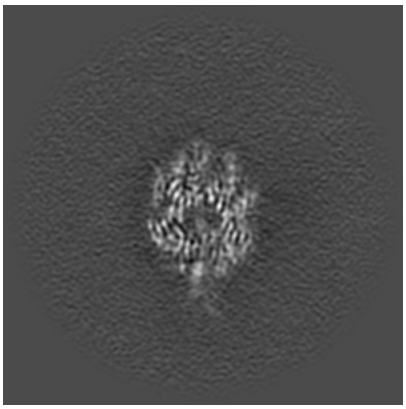


Z Index: 180

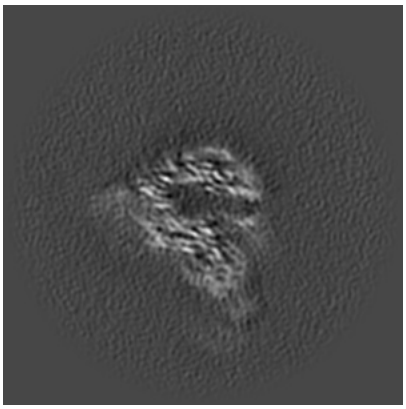
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

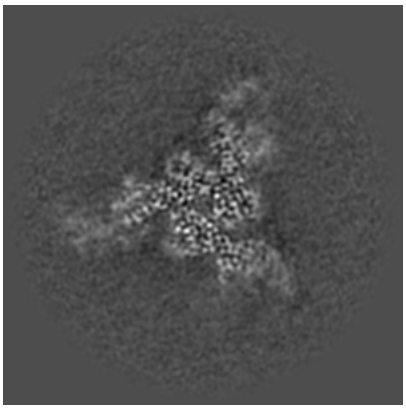
6.3.1 Primary map



X Index: 167

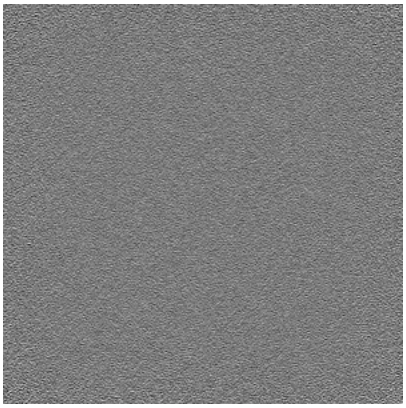


Y Index: 180

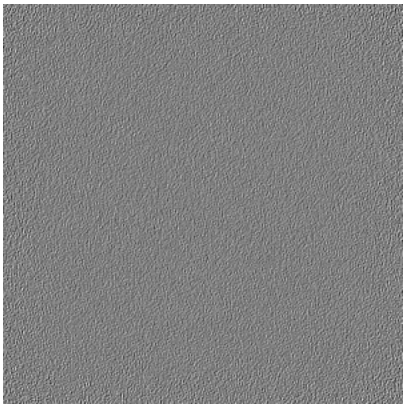


Z Index: 191

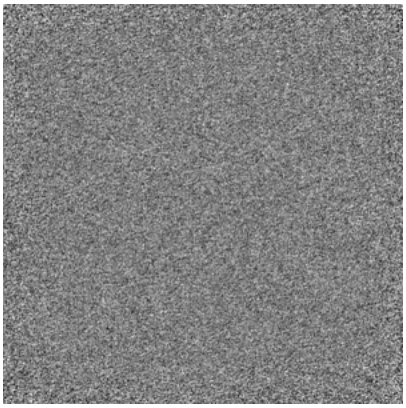
6.3.2 Raw map



X Index: 0



Y Index: 0

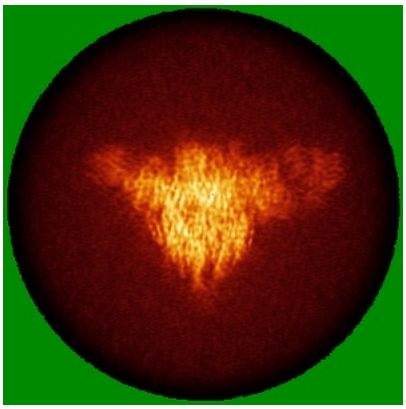


Z Index: 0

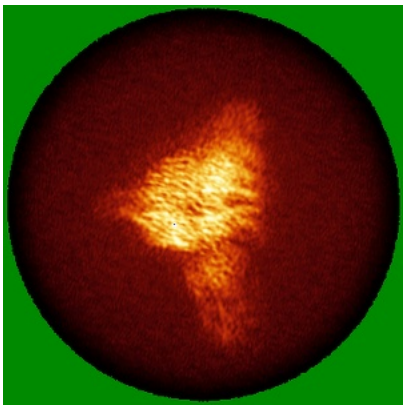
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

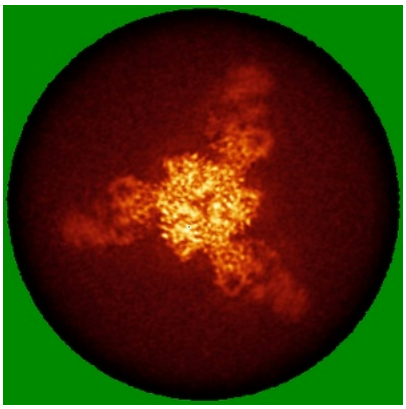
6.4.1 Primary map



X

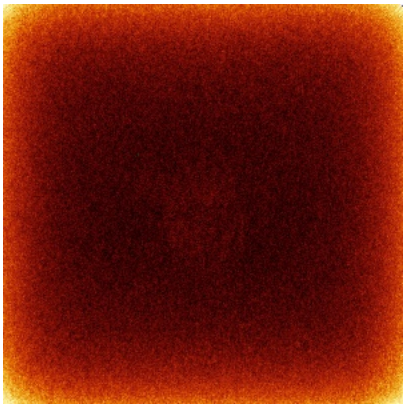


Y

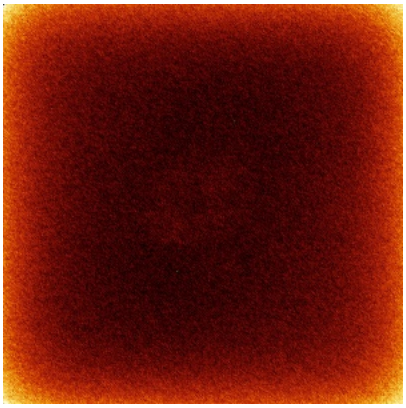


Z

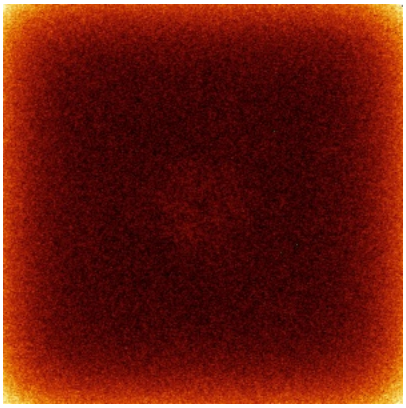
6.4.2 Raw map



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)

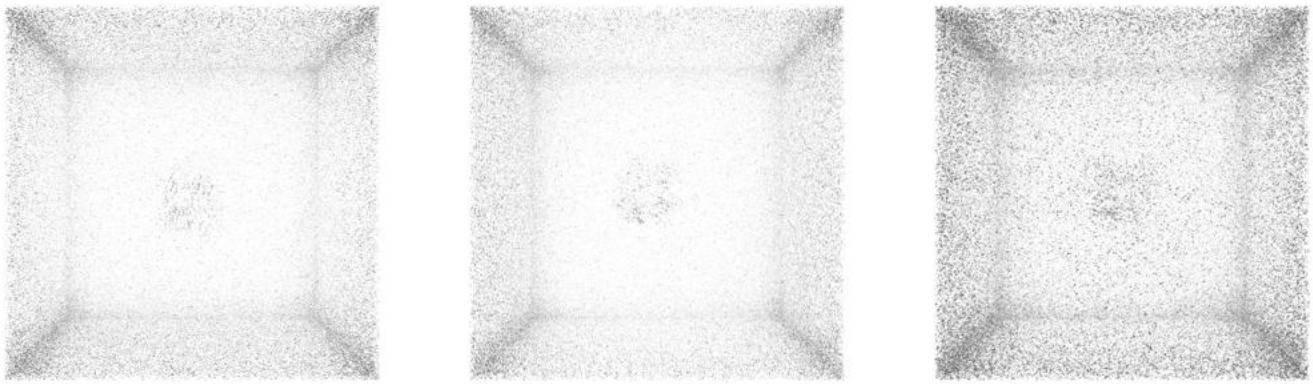
6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.054. 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



X Y Z

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.

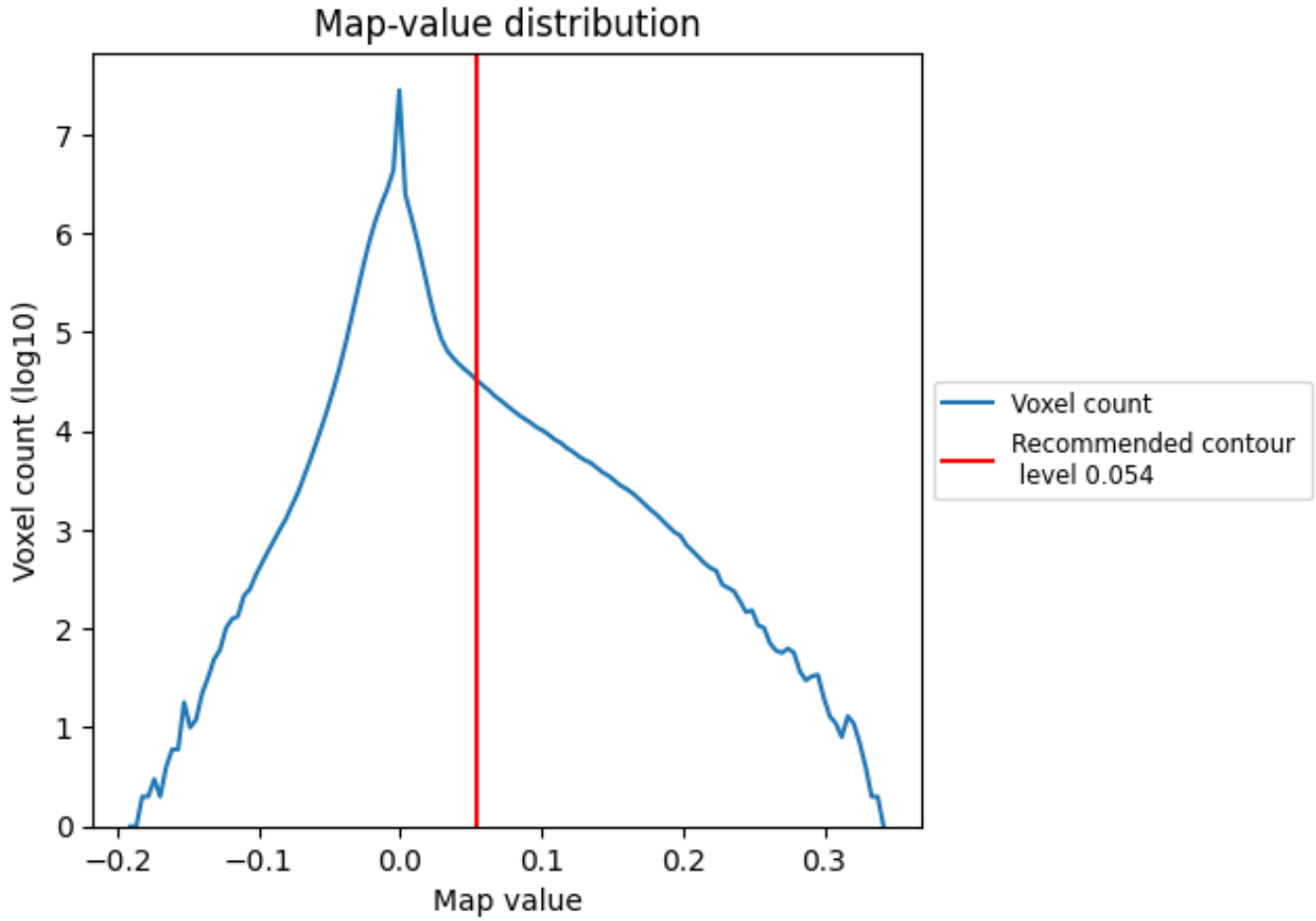
6.6 Mask visualisation (i)

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

7 Map analysis

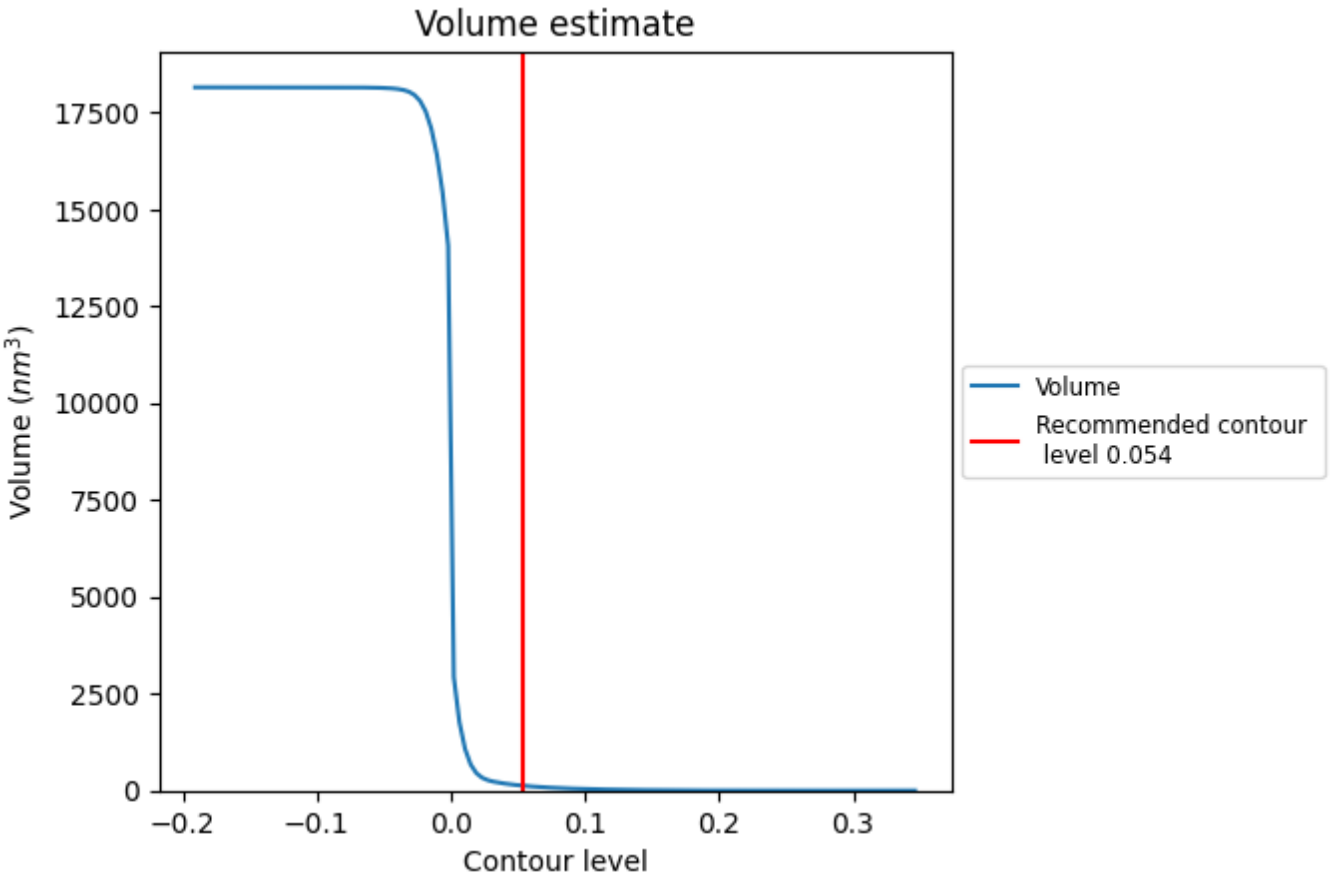
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

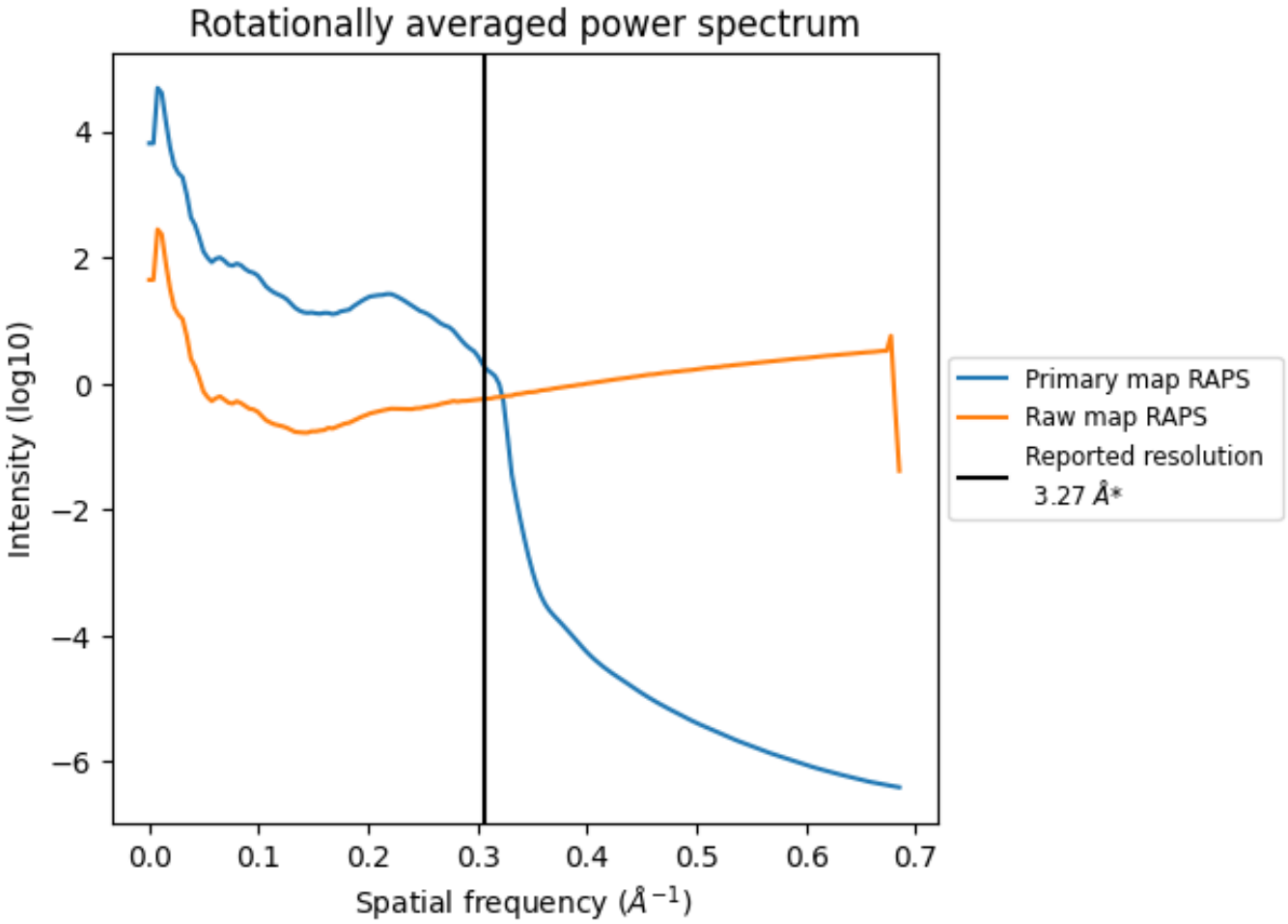
7.2 Volume estimate (i)



The volume at the recommended contour level is 125 nm^3; this corresponds to an approximate mass of 113 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 (i)

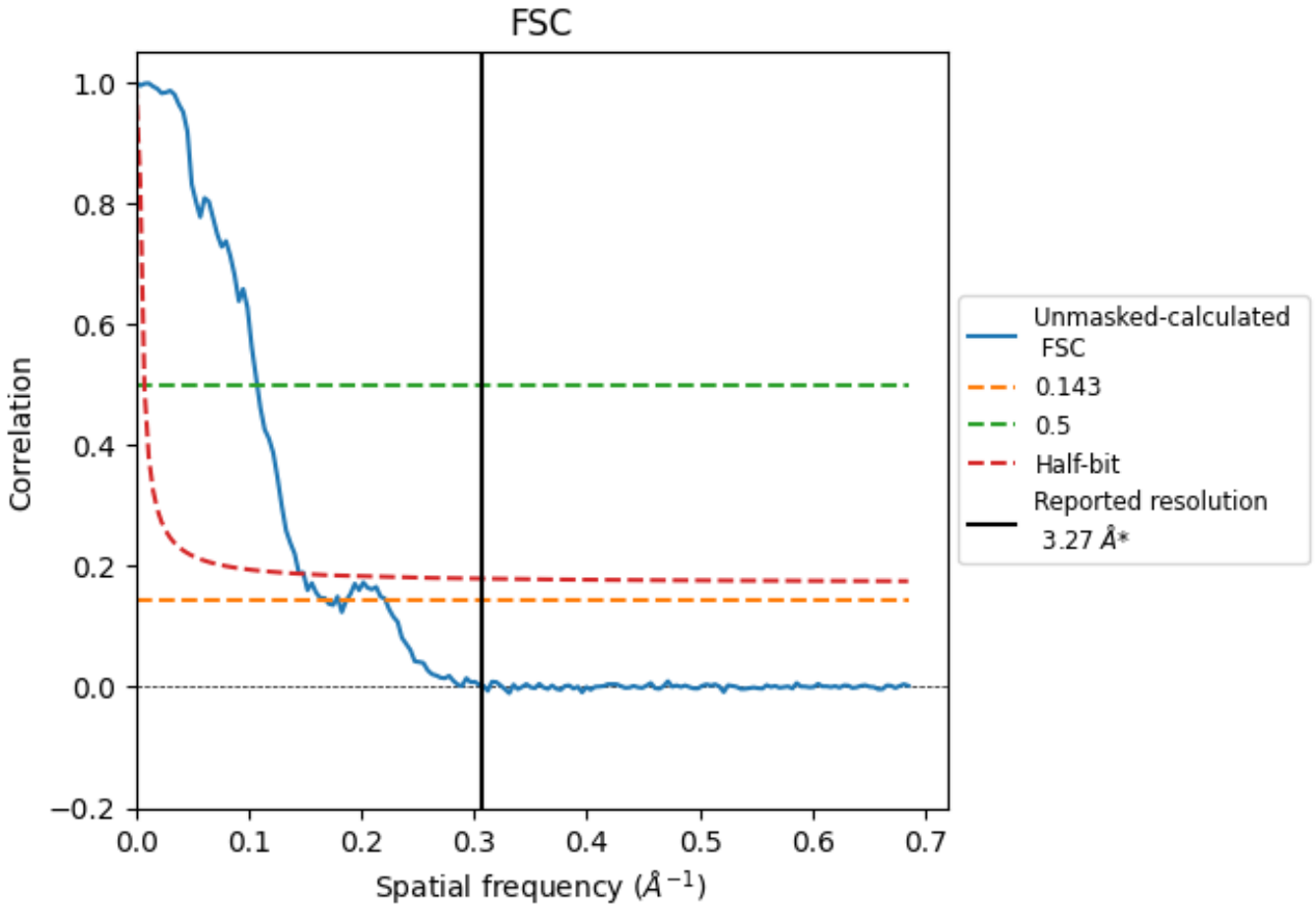


*Reported resolution corresponds to spatial frequency of 0.306 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.306 Å^{-1}

8.2 Resolution estimates (i)

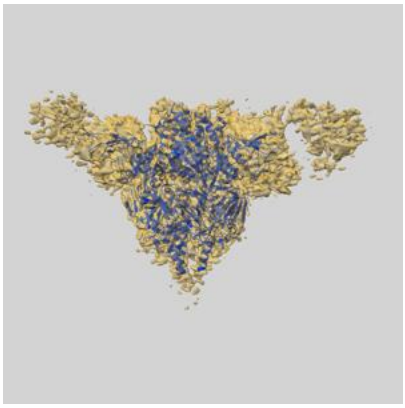
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.27	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	5.92	9.30	6.72

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 5.92 differs from the reported value 3.27 by more than 10 %

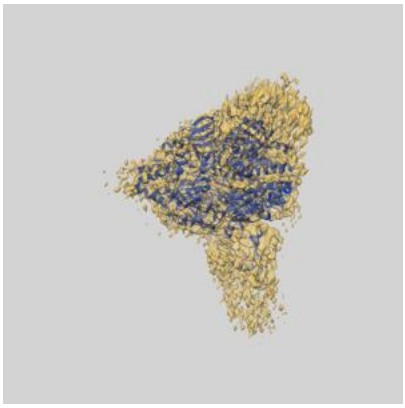
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-65894 and PDB model 9WDP. Per-residue inclusion information can be found in section 3 on page 10.

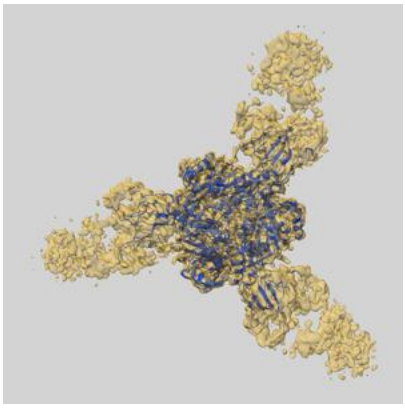
9.1 Map-model overlay



X



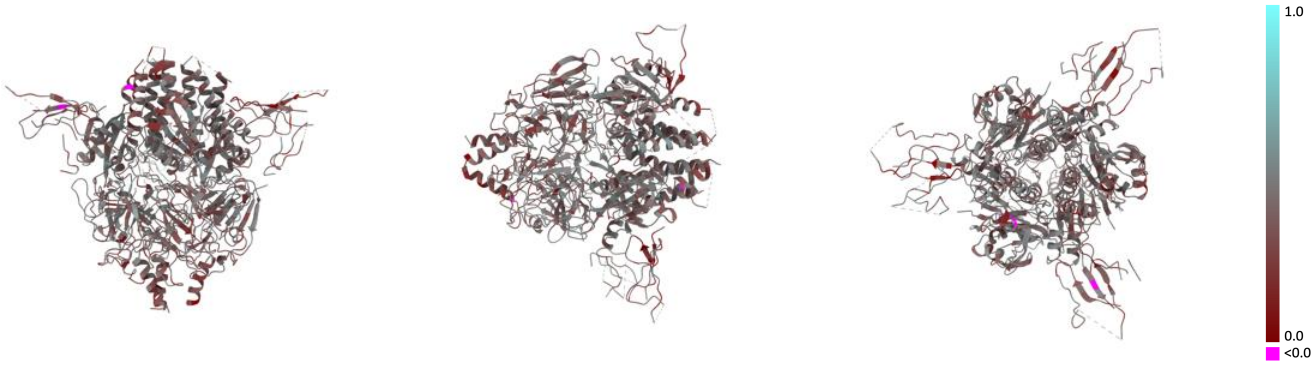
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.054 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.

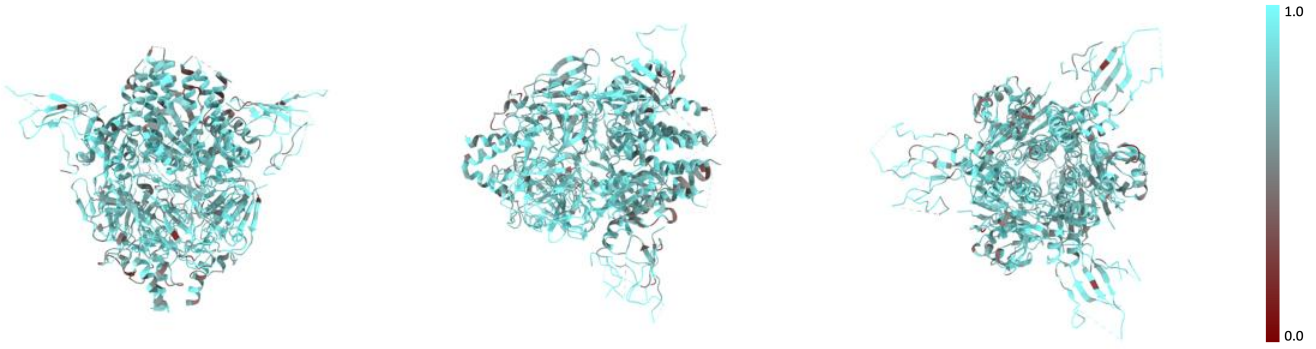
9.2 Q-score mapped to coordinate model



X Y Z

The images above show the model with each residue coloured according 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.

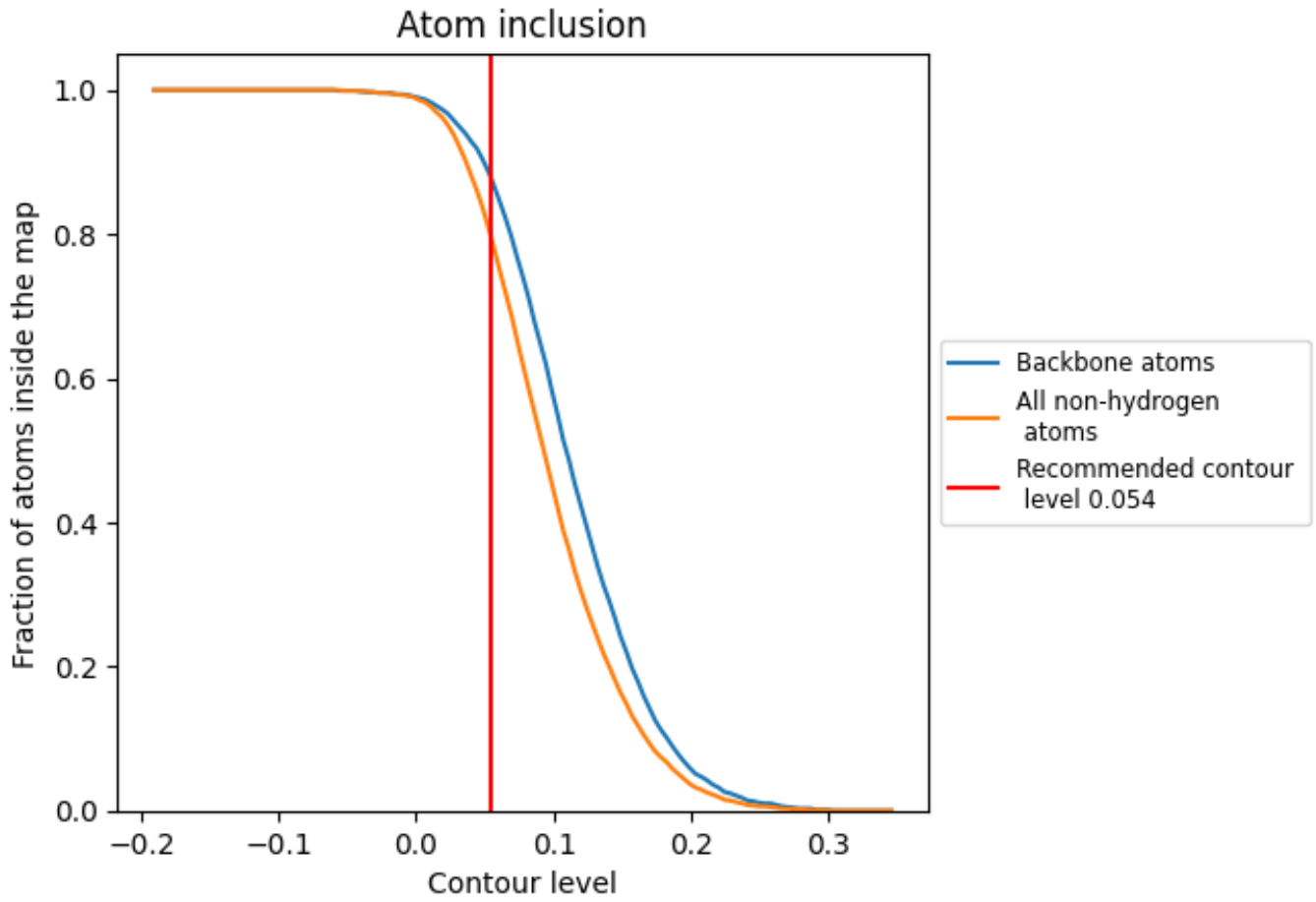
9.3 Atom inclusion mapped to coordinate model



X Y Z

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 (0.054).

9.4 Atom inclusion (i)



At the recommended contour level, 88% of all backbone atoms, 80% 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 (0.054) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8000	0.3990
A	0.7880	0.4100
B	0.8040	0.4060
C	0.8000	0.4000
D	0.7870	0.3130
E	0.8740	0.4370
F	0.8540	0.3720
G	0.7950	0.3500
H	0.8000	0.3430
L	0.7530	0.3700

