wwPDB EM Validation Summary Report

Sep 24, 2024 – 04:31 PM EDT

PDB ID	:	8U4S
EMDB ID	:	EMD-41893
Title	:	Structure of trimeric CXCR4 in complex with REGN7663 Fab
Authors	:	Saotome, K.; McGoldrick, L.L.; Franklin, M.C.
Deposited on	:	2023-09-11
Resolution	:	3.35 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:

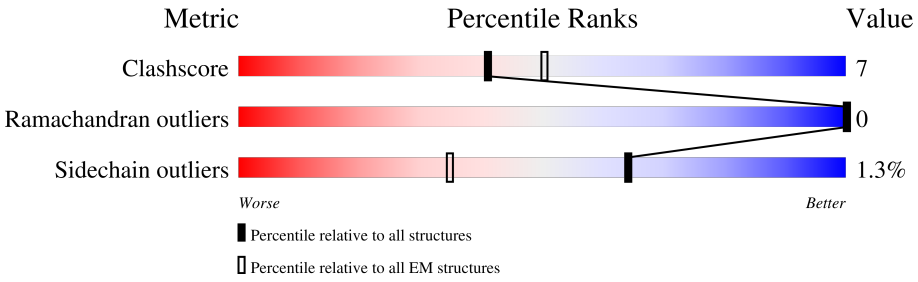
EMDB validation analysis	:	0.0.1.dev112
Mogul	:	2022.3.0, CSD as543be (2022)
MolProbity	:	4.02b-467
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.38.3

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 3.35 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	219	42% 8% • 50%
1	E	219	41% 9% • 50%
1	L	219	43% 6% • 50%
2	B	240	43% 9% 48%
2	F	240	41% 11% 48%
2	H	240	43% 8% 48%
3	C	632	37% 7% 55%
3	G	632	38% 7% 55%

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Mol	Chain	Length	Quality of chain
3	R	632	

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 12684 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 REGN7663 Fab light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	L	110	Total	C	N	O	S	0	0
			856	536	149	167	4		
1	A	110	Total	C	N	O	S	0	0
			856	536	149	167	4		
1	E	110	Total	C	N	O	S	0	0
			856	536	149	167	4		

- Molecule 2 is a protein called REGN7663 Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	125	Total	C	N	O	S	0	0
			976	609	174	187	6		
2	B	125	Total	C	N	O	S	0	0
			976	609	174	187	6		
2	F	125	Total	C	N	O	S	0	0
			976	609	174	187	6		

- Molecule 3 is a protein called C-X-C chemokine receptor type 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	R	284	Total	C	N	O	S	0	0
			2297	1539	365	380	13		
3	C	284	Total	C	N	O	S	0	0
			2297	1539	365	380	13		
3	G	284	Total	C	N	O	S	0	0
			2297	1539	365	380	13		

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	-17	MET	-	initiating methionine	UNP P61073
R	-16	LYS	-	expression tag	UNP P61073
R	-15	THR	-	expression tag	UNP P61073

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Chain	Residue	Modelled	Actual	Comment	Reference
R	-14	ILE	-	expression tag	UNP P61073
R	-13	ILE	-	expression tag	UNP P61073
R	-12	ALA	-	expression tag	UNP P61073
R	-11	LEU	-	expression tag	UNP P61073
R	-10	SER	-	expression tag	UNP P61073
R	-9	TYR	-	expression tag	UNP P61073
R	-8	ILE	-	expression tag	UNP P61073
R	-7	PHE	-	expression tag	UNP P61073
R	-6	CYS	-	expression tag	UNP P61073
R	-5	LEU	-	expression tag	UNP P61073
R	-4	VAL	-	expression tag	UNP P61073
R	-3	PHE	-	expression tag	UNP P61073
R	-2	ALA	-	expression tag	UNP P61073
R	-1	GLY	-	expression tag	UNP P61073
R	0	ALA	-	expression tag	UNP P61073
R	1	PRO	-	expression tag	UNP P61073
R	119	SER	ASN	conflict	UNP P61073
C	-17	MET	-	initiating methionine	UNP P61073
C	-16	LYS	-	expression tag	UNP P61073
C	-15	THR	-	expression tag	UNP P61073
C	-14	ILE	-	expression tag	UNP P61073
C	-13	ILE	-	expression tag	UNP P61073
C	-12	ALA	-	expression tag	UNP P61073
C	-11	LEU	-	expression tag	UNP P61073
C	-10	SER	-	expression tag	UNP P61073
C	-9	TYR	-	expression tag	UNP P61073
C	-8	ILE	-	expression tag	UNP P61073
C	-7	PHE	-	expression tag	UNP P61073
C	-6	CYS	-	expression tag	UNP P61073
C	-5	LEU	-	expression tag	UNP P61073
C	-4	VAL	-	expression tag	UNP P61073
C	-3	PHE	-	expression tag	UNP P61073
C	-2	ALA	-	expression tag	UNP P61073
C	-1	GLY	-	expression tag	UNP P61073
C	0	ALA	-	expression tag	UNP P61073
C	1	PRO	-	expression tag	UNP P61073
C	119	SER	ASN	conflict	UNP P61073
G	-17	MET	-	initiating methionine	UNP P61073
G	-16	LYS	-	expression tag	UNP P61073
G	-15	THR	-	expression tag	UNP P61073
G	-14	ILE	-	expression tag	UNP P61073
G	-13	ILE	-	expression tag	UNP P61073

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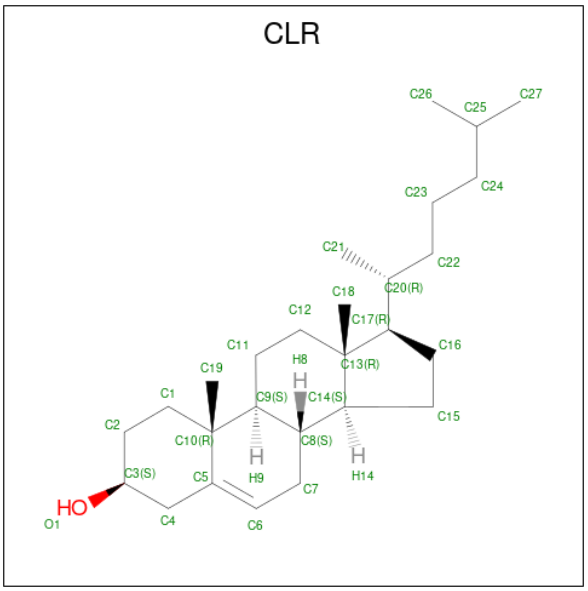
Chain	Residue	Modelled	Actual	Comment	Reference
G	-12	ALA	-	expression tag	UNP P61073
G	-11	LEU	-	expression tag	UNP P61073
G	-10	SER	-	expression tag	UNP P61073
G	-9	TYR	-	expression tag	UNP P61073
G	-8	ILE	-	expression tag	UNP P61073
G	-7	PHE	-	expression tag	UNP P61073
G	-6	CYS	-	expression tag	UNP P61073
G	-5	LEU	-	expression tag	UNP P61073
G	-4	VAL	-	expression tag	UNP P61073
G	-3	PHE	-	expression tag	UNP P61073
G	-2	ALA	-	expression tag	UNP P61073
G	-1	GLY	-	expression tag	UNP P61073
G	0	ALA	-	expression tag	UNP P61073
G	1	PRO	-	expression tag	UNP P61073
G	119	SER	ASN	conflict	UNP P61073

- Molecule 4 is CHOLESTEROL (three-letter code: CLR) (formula: $C_{27}H_{46}O$).



Mol	Chain	Residues	Atoms	AltConf
4	R	1	Total C O 28 27 1	0
4	R	1	Total C O 28 27 1	0
4	C	1	Total C O 28 27 1	0
4	C	1	Total C O 28 27 1	0

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Mol	Chain	Residues	Atoms	AltConf
4	G	1	Total C O 28 27 1	0
4	G	1	Total C O 28 27 1	0

- Molecule 5 is (2R)-1-(hexadecanoyloxy)-3-(phosphonooxy)propan-2-yl (9Z)-octadec-9-enoate (three-letter code: D21) (formula: $C_{37}H_{71}O_8P$).

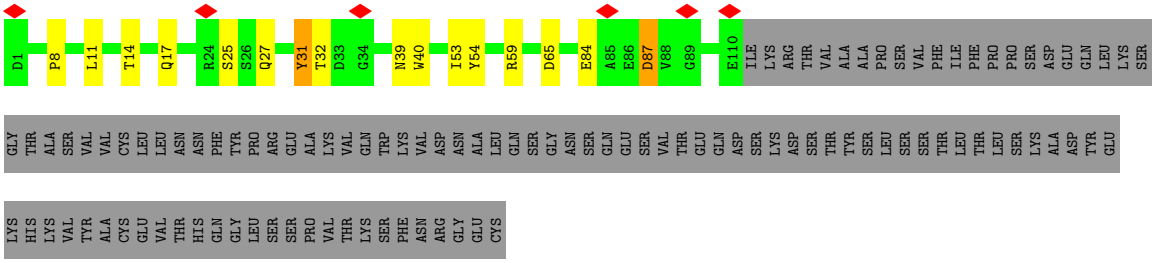
Mol	Chain	Residues	Atoms	AltConf
5	R	1	Total C O P 43 34 8 1	0
5	C	1	Total C O P 43 34 8 1	0
5	G	1	Total C O P 43 34 8 1	0

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.

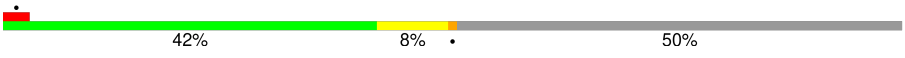
• Molecule 1: REGN7663 Fab light chain

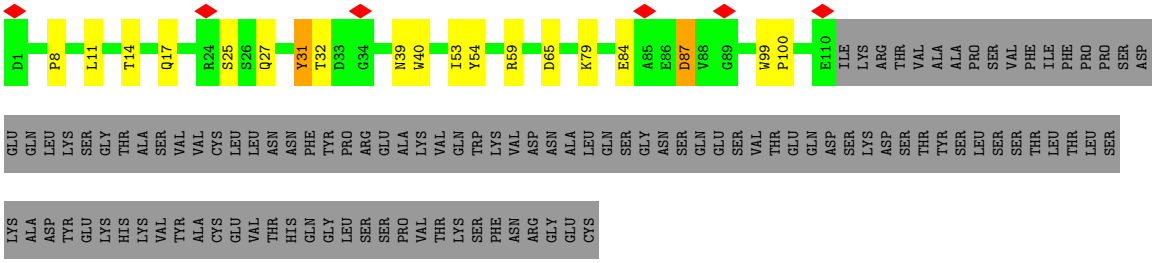
Chain L:



• Molecule 1: REGN7663 Fab light chain

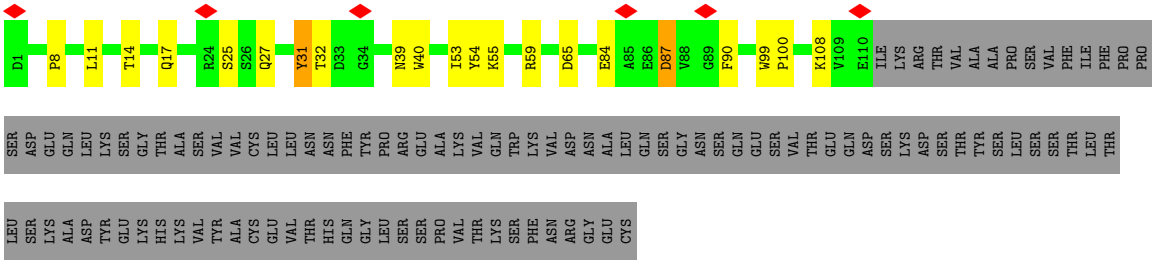
Chain A:





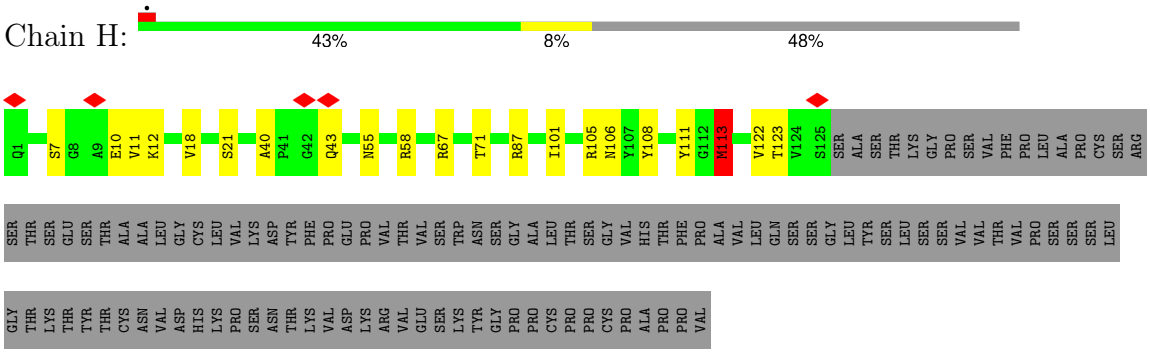
• Molecule 1: REGN7663 Fab light chain

Chain E:



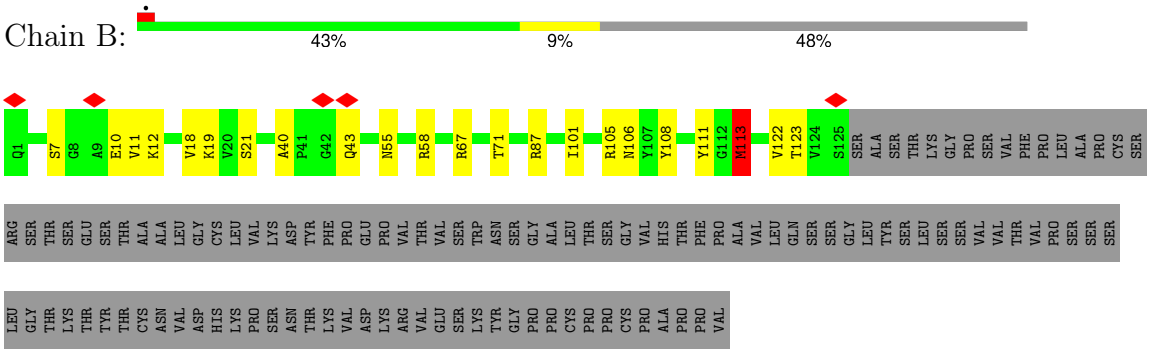
• Molecule 2: REGN7663 Fab heavy chain

Chain H:



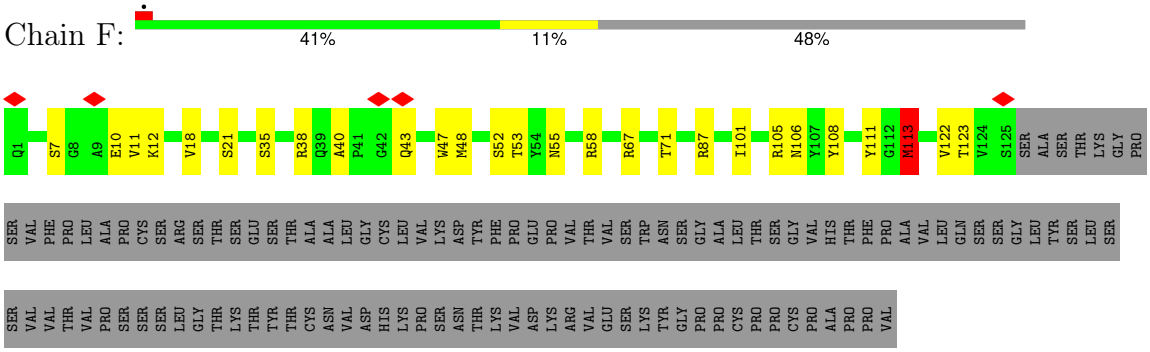
• Molecule 2: REGN7663 Fab heavy chain

Chain B:



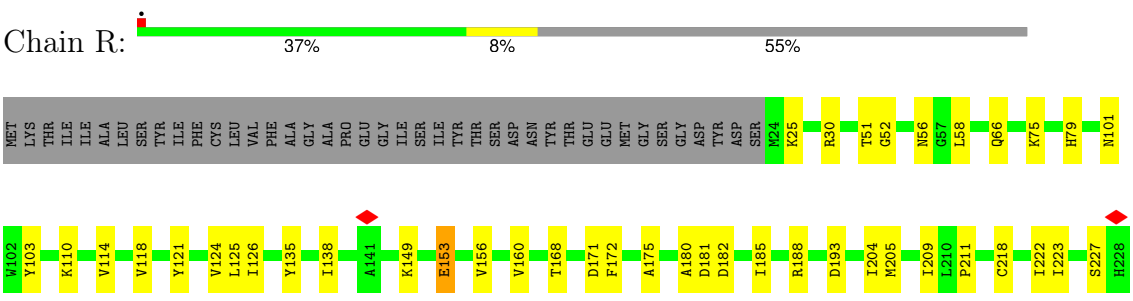
• Molecule 2: REGN7663 Fab heavy chain

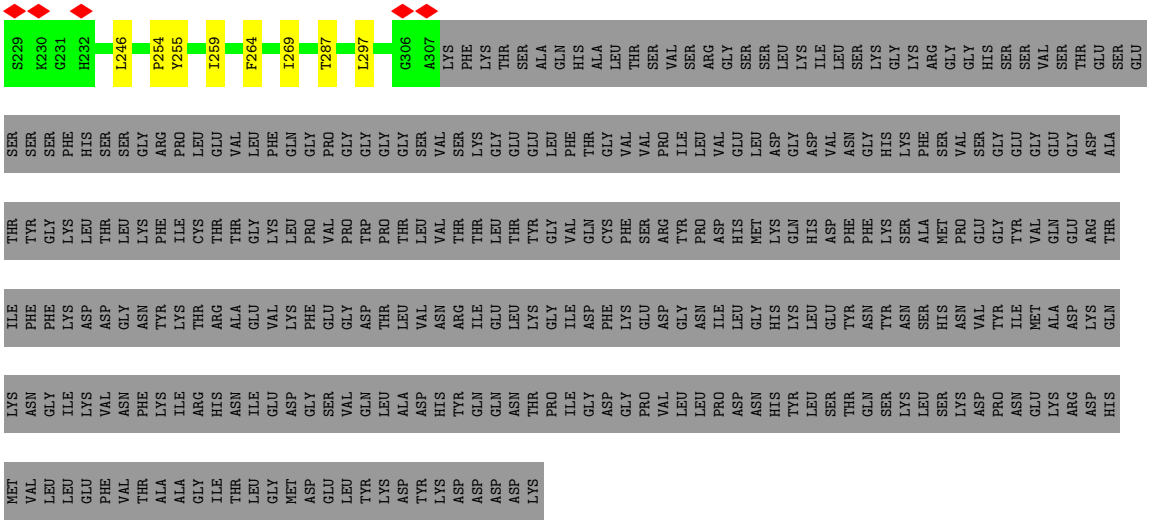
Chain F:



• Molecule 3: C-X-C chemokine receptor type 4

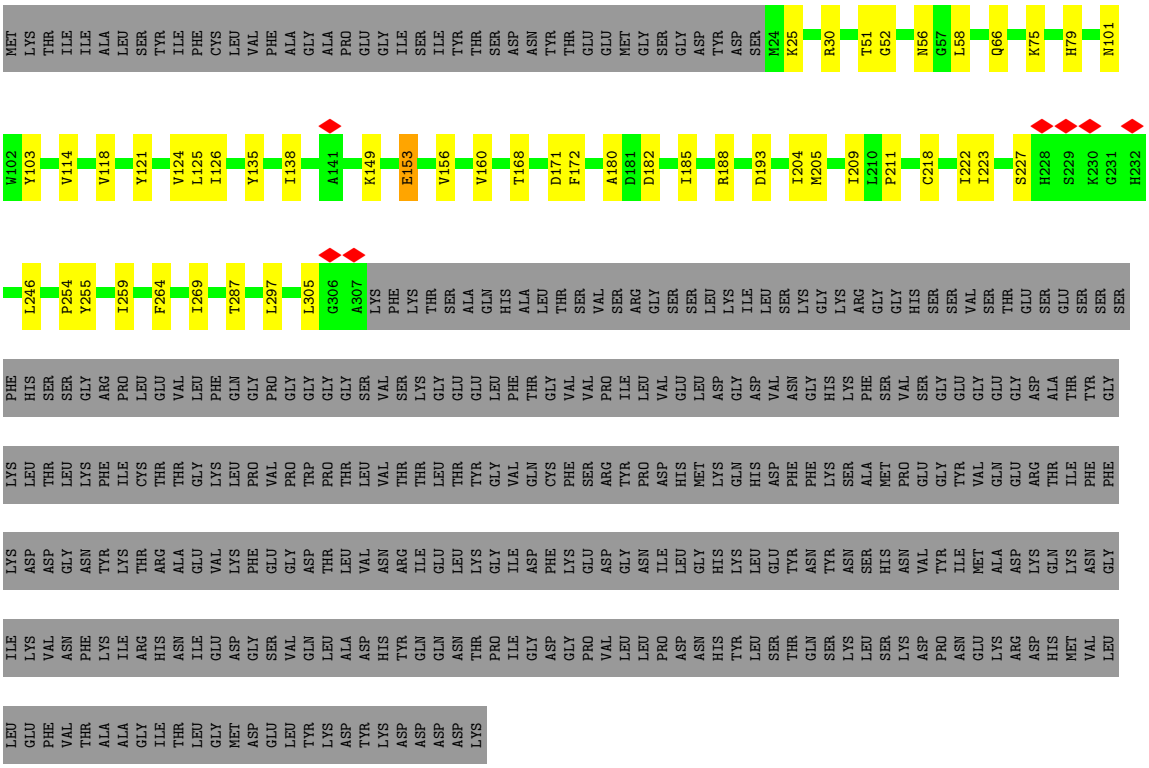
Chain R:





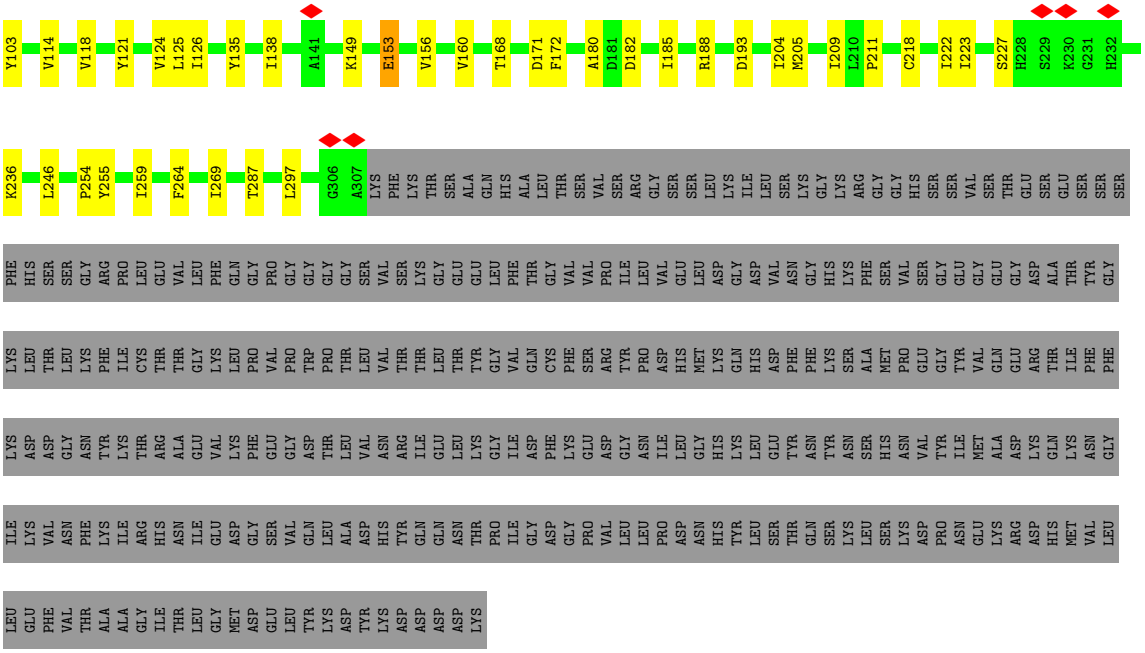
• Molecule 3: C-X-C chemokine receptor type 4

Chain C:



• Molecule 3: C-X-C chemokine receptor type 4

Chain G:


Y103 V114 V118 Y121 V124 L125 I126 Y135 I138 A141 K149 E153 V156 V160 T168 D171 F172 A180 D181 D182 I185 R188 D193 I204 M205 I209 L210 P211 C218 I222 I223 S227 H228 S229 K230 G231 H232
K236 L246 P254 Y255 I259 F264 I269 T287 L297 G306 A307 LYS PHE LYS THR SER ALA GLN HIS ALA LEU THR SER VAL SER ARG GLY SER SER LEU LYS ILE LEU SER LYS GLY LYS ARG GLY GLY HIS SER SER VAL SER THR GLU SER GLU SER SER SER
PHE HIS SER SER GLY ARG PRO LEU GLU VAL LEU PHE GLN GLY PRO GLY GLY GLY GLY GLY SER VAL SER LYS GLY GLU GLU LEU PHE THR GLY VAL VAL PRO ILE LEU VAL GLU LEU ASP GLY ASP VAL ASN GLY HIS LYS PHE SER VAL SER GLY GLU GLY GLU GLY ASP ALA THR TYR GLY
LYS LEU THR LEU LYS PHE ILE CYS THR THR GLY LYS LEU PRO VAL PRO TRP PRO THR LEU VAL THR THR LEU THR TYR GLY VAL GLN CYS PHE SER ARG TYR PRO ASP HIS MET LYS GLN HIS ASP PHE PHE LYS SER ALA MET PRO GLU GLY TYR VAL GLN GLU ARG THR ILE PHE PHE
LYS ASP ASP GLY ASN TYR LYS THR ARG ALA GLU VAL LYS PHE GLU GLY ASP THR LEU VAL ASN ARG ILE GLU LEU LYS GLY ILE ASP PHE LYS GLU ASP GLY ASN ILE LEU GLY HIS LYS LEU GLU TYR ASN TYR ASN SER HIS ASN VAL TYR ILE MET ALA ASP LYS GLN LYS ASN GLY
ILE LYS VAL ASN PHE LYS ILE ARG HIS ASN ILE GLU ASP GLY SER VAL GLN LEU ALA ASP HIS TYR GLN GLN ASN THR PRO ILE GLY ASP GLY PRO VAL LEU LEU PRO ASP ASN HIS TYR LEU SER THR GLN SER LYS LEU SER LYS ASP PRO ASN GLU LYS ARG ASP HIS MET VAL LEU
LEU GLU PHE VAL THR ALA ALA GLY ILE THR LEU GLY MET ASP GLU LEU TYR LYS ASP TYR LYS ASP ASP ASP ASP LYS

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	27104	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 ($e^-/Å^2$)	40	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.975	Depositor
Minimum map value	-0.647	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.029	Depositor
Recommended contour level	0.15	Depositor
Map size (Å)	306.0, 306.0, 306.0	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.85, 0.85, 0.85	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: D21, CLR

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	0.29	0/877	0.59	0/1191
1	E	0.29	0/877	0.59	0/1191
1	L	0.29	0/877	0.59	0/1191
2	B	0.28	0/999	0.63	2/1355 (0.1%)
2	F	0.27	0/999	0.63	2/1355 (0.1%)
2	H	0.27	0/999	0.63	2/1355 (0.1%)
3	C	0.29	0/2361	0.48	0/3214
3	G	0.29	0/2361	0.48	0/3214
3	R	0.29	0/2361	0.48	0/3214
All	All	0.28	0/12711	0.54	6/17280 (0.0%)

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	113	MET	CB-CG-SD	7.02	133.45	112.40
2	H	113	MET	CB-CG-SD	7.01	133.43	112.40
2	B	113	MET	CB-CG-SD	7.00	133.40	112.40
2	B	113	MET	CA-CB-CG	5.82	123.20	113.30
2	H	113	MET	CA-CB-CG	5.80	123.16	113.30

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	856	0	824	13	0
1	E	856	0	824	15	0
1	L	856	0	824	12	0
2	B	976	0	934	13	0
2	F	976	0	934	16	0
2	H	976	0	934	12	0
3	C	2297	0	2365	37	0
3	G	2297	0	2365	36	0
3	R	2297	0	2365	38	0
4	C	56	0	92	4	0
4	G	56	0	92	5	0
4	R	56	0	92	5	0
5	C	43	0	0	0	0
5	G	43	0	0	0	0
5	R	43	0	0	0	0
All	All	12684	0	12645	170	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 7.

The worst 5 of 170 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:L:31:TYR:HD1	1:L:32:THR:H	1.13	0.97
1:E:31:TYR:HD1	1:E:32:THR:H	1.13	0.93
1:A:31:TYR:HD1	1:A:32:THR:H	1.13	0.92
4:R:703:CLR:H112	4:G:703:CLR:H12	1.73	0.71
4:C:703:CLR:H12	4:G:703:CLR:H112	1.73	0.71

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	108/219 (49%)	105 (97%)	3 (3%)	0	100	100
1	E	108/219 (49%)	105 (97%)	3 (3%)	0	100	100
1	L	108/219 (49%)	105 (97%)	3 (3%)	0	100	100
2	B	123/240 (51%)	117 (95%)	6 (5%)	0	100	100
2	F	123/240 (51%)	117 (95%)	6 (5%)	0	100	100
2	H	123/240 (51%)	117 (95%)	6 (5%)	0	100	100
3	C	282/632 (45%)	276 (98%)	6 (2%)	0	100	100
3	G	282/632 (45%)	276 (98%)	6 (2%)	0	100	100
3	R	282/632 (45%)	276 (98%)	6 (2%)	0	100	100
All	All	1539/3273 (47%)	1494 (97%)	45 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	97/195 (50%)	95 (98%)	2 (2%)	48	69
1	E	97/195 (50%)	95 (98%)	2 (2%)	48	69
1	L	97/195 (50%)	95 (98%)	2 (2%)	48	69
2	B	102/203 (50%)	99 (97%)	3 (3%)	37	62
2	F	102/203 (50%)	99 (97%)	3 (3%)	37	62
2	H	102/203 (50%)	99 (97%)	3 (3%)	37	62
3	C	252/552 (46%)	251 (100%)	1 (0%)	89	94
3	G	252/552 (46%)	251 (100%)	1 (0%)	89	94
3	R	252/552 (46%)	251 (100%)	1 (0%)	89	94
All	All	1353/2850 (48%)	1335 (99%)	18 (1%)	64	79

5 of 18 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	F	58	ARG
3	G	153	GLU
2	F	113	MET
2	B	58	ARG
1	E	87	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	66	GLN
2	F	57	ASN
3	G	66	GLN
3	R	66	GLN
2	H	57	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)

9 ligands are 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	CLR	G	701	-	31,31,31	1.08	2 (6%)	48,48,48	1.33	7 (14%)
4	CLR	C	701	-	31,31,31	1.07	2 (6%)	48,48,48	1.33	7 (14%)
5	D21	C	702	-	42,42,45	1.47	4 (9%)	45,47,50	1.10	4 (8%)
4	CLR	C	703	-	31,31,31	1.10	2 (6%)	48,48,48	1.36	7 (14%)
5	D21	G	702	-	42,42,45	1.47	4 (9%)	45,47,50	1.10	4 (8%)
4	CLR	R	703	-	31,31,31	1.09	2 (6%)	48,48,48	1.36	6 (12%)
5	D21	R	702	-	42,42,45	1.47	4 (9%)	45,47,50	1.10	4 (8%)
4	CLR	G	703	-	31,31,31	1.10	2 (6%)	48,48,48	1.36	6 (12%)
4	CLR	R	701	-	31,31,31	1.08	2 (6%)	48,48,48	1.33	7 (14%)

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	CLR	G	701	-	-	7/10/68/68	0/4/4/4
4	CLR	C	701	-	-	7/10/68/68	0/4/4/4
5	D21	C	702	-	-	19/44/44/47	-
4	CLR	C	703	-	-	4/10/68/68	0/4/4/4
5	D21	G	702	-	-	19/44/44/47	-
4	CLR	R	703	-	-	4/10/68/68	0/4/4/4
5	D21	R	702	-	-	19/44/44/47	-
4	CLR	G	703	-	-	4/10/68/68	0/4/4/4
4	CLR	R	701	-	-	7/10/68/68	0/4/4/4

The worst 5 of 24 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	R	702	D21	P23-O22	8.07	1.85	1.60
5	G	702	D21	P23-O22	8.05	1.85	1.60
5	C	702	D21	P23-O22	8.04	1.85	1.60
5	R	702	D21	O22-C21	-2.89	1.33	1.44
5	G	702	D21	O22-C21	-2.88	1.33	1.44

The worst 5 of 52 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	G	702	D21	O25-P23-O22	-3.02	98.81	106.67
5	R	702	D21	O25-P23-O22	-3.01	98.81	106.67
5	C	702	D21	O25-P23-O22	-3.01	98.83	106.67
4	R	701	CLR	C13-C17-C20	-2.98	114.89	119.50
4	G	701	CLR	C13-C17-C20	-2.98	114.89	119.50

There are no chirality outliers.

5 of 90 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	R	701	CLR	C21-C20-C22-C23
4	C	701	CLR	C21-C20-C22-C23
4	G	701	CLR	C21-C20-C22-C23
4	R	703	CLR	C21-C20-C22-C23
4	C	703	CLR	C21-C20-C22-C23

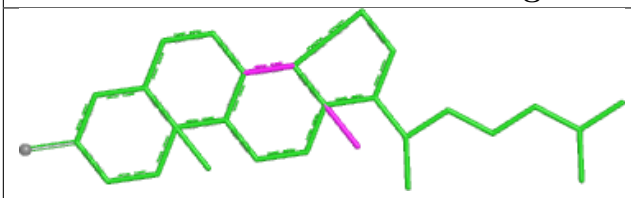
There are no ring outliers.

5 monomers are involved in 11 short contacts:

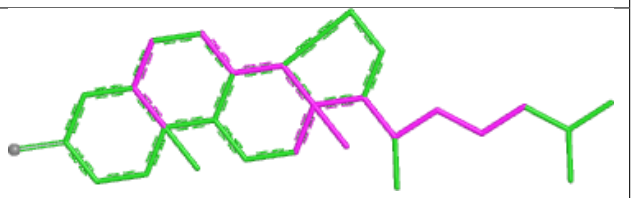
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	701	CLR	1	0
4	C	703	CLR	4	0
4	R	703	CLR	4	0
4	G	703	CLR	4	0
4	R	701	CLR	1	0

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.

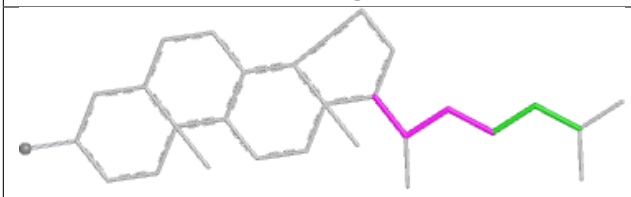
Ligand CLR G 701



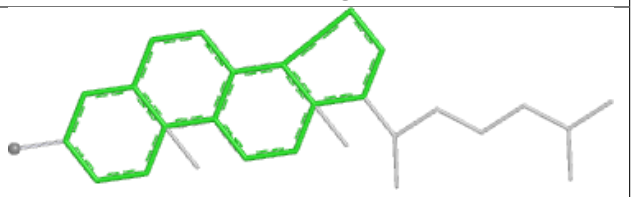
Bond lengths



Bond angles

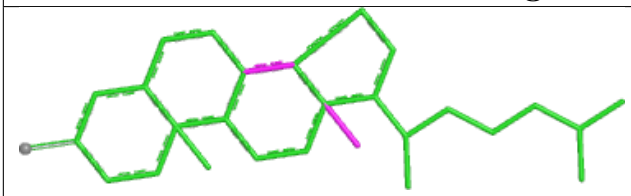


Torsions

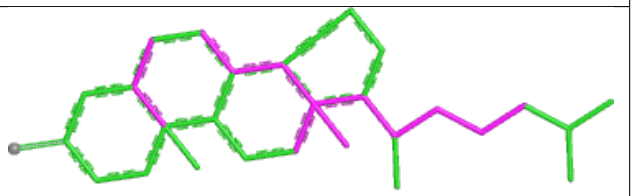


Rings

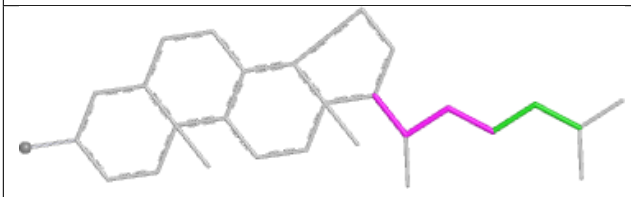
Ligand CLR C 701



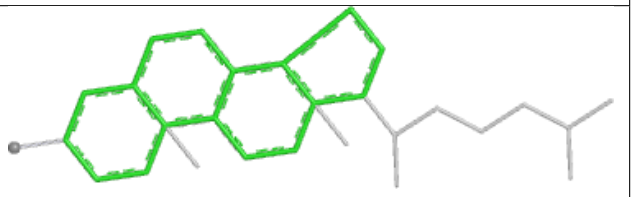
Bond lengths



Bond angles

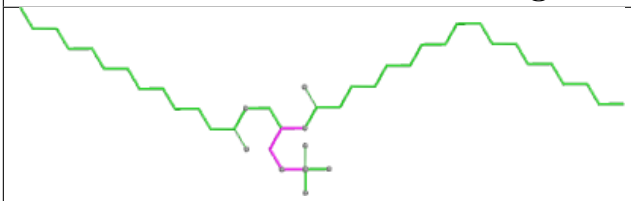


Torsions

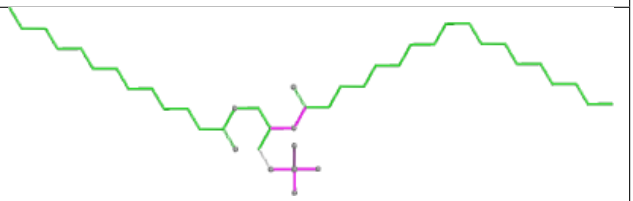


Rings

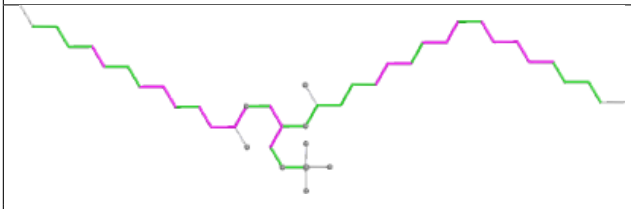
Ligand D21 C 702



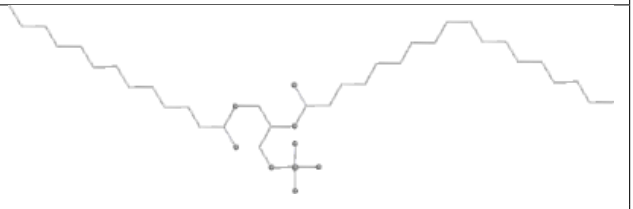
Bond lengths



Bond angles

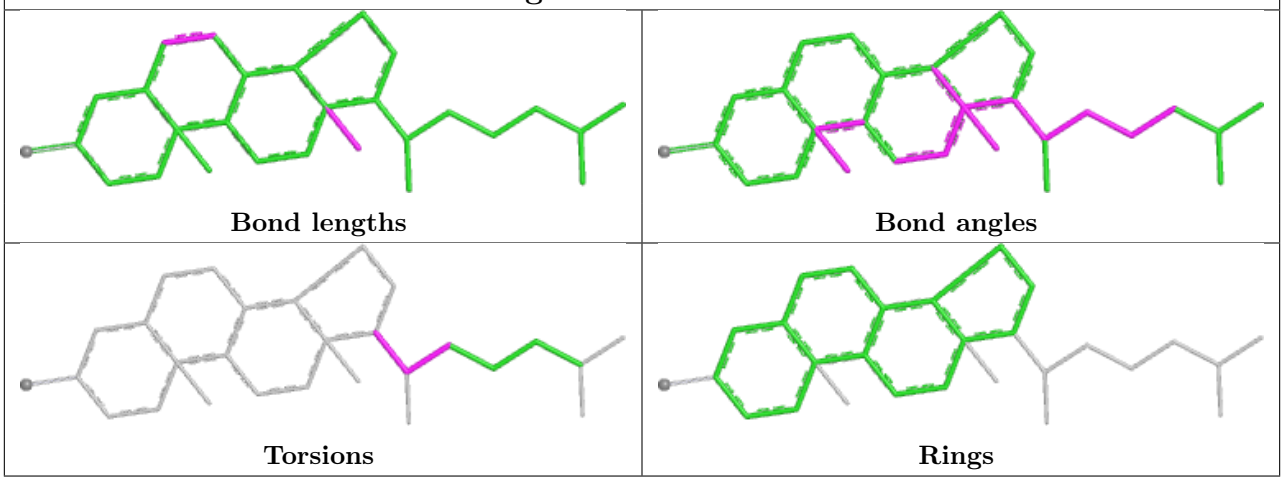


Torsions

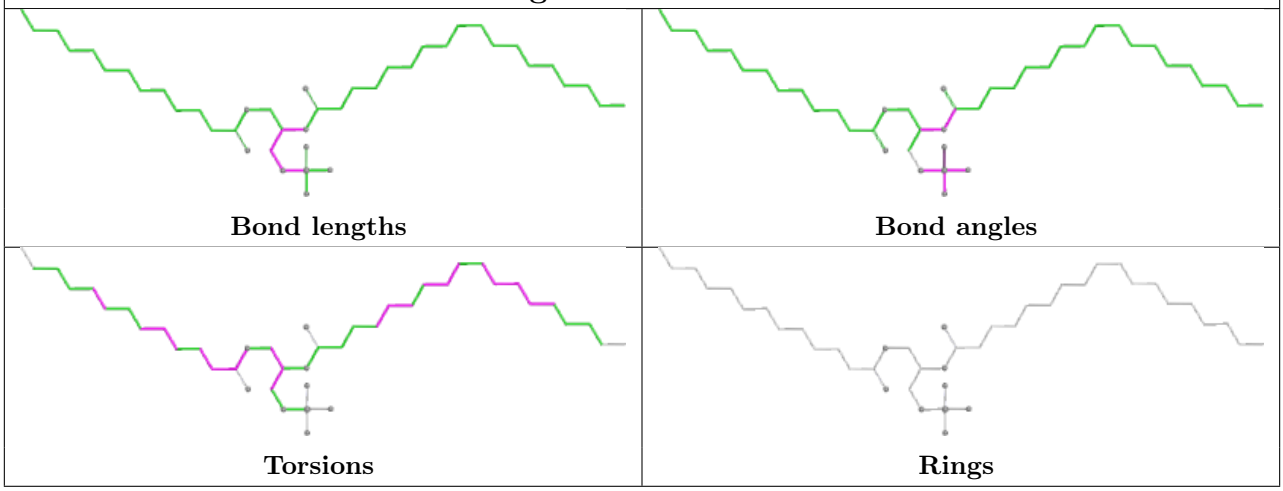


Rings

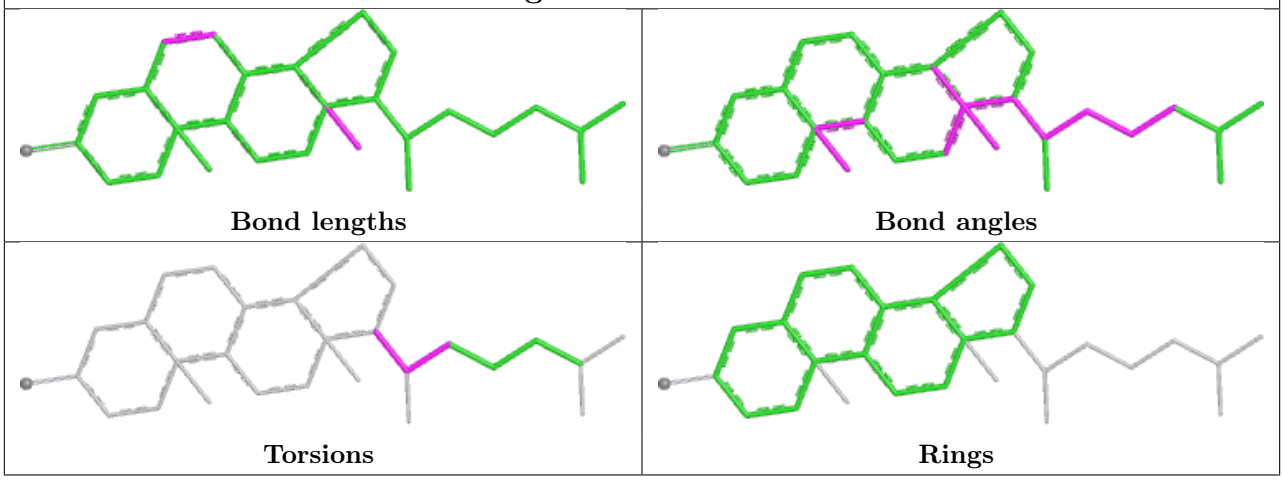
Ligand CLR C 703



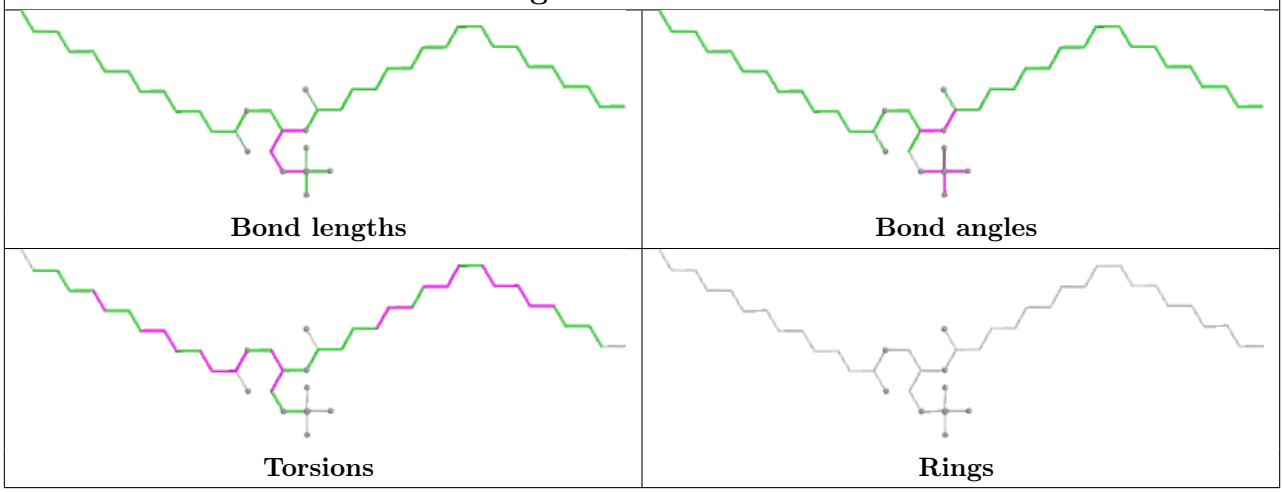
Ligand D21 G 702



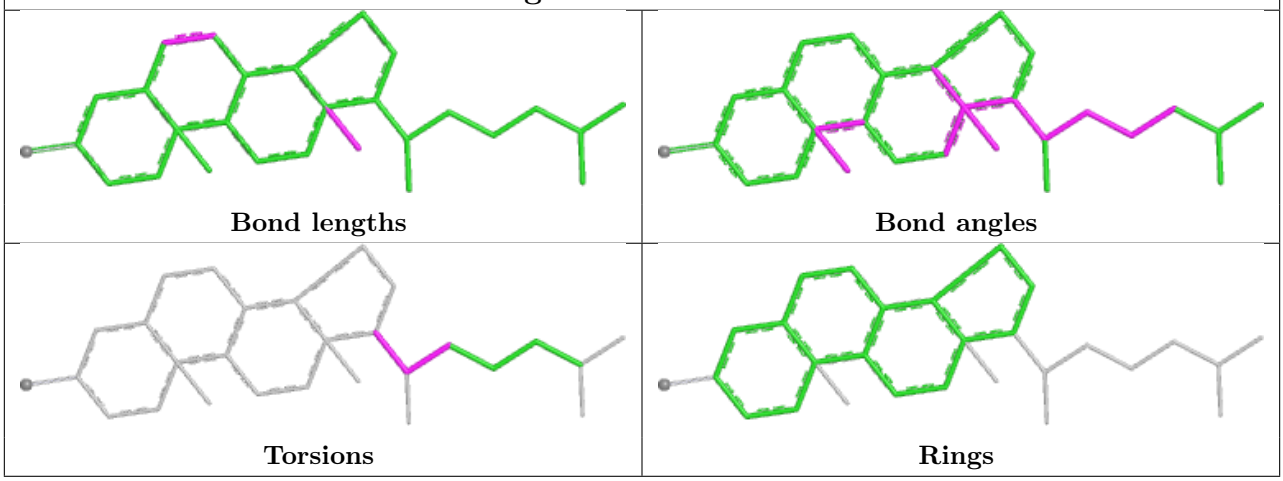
Ligand CLR R 703



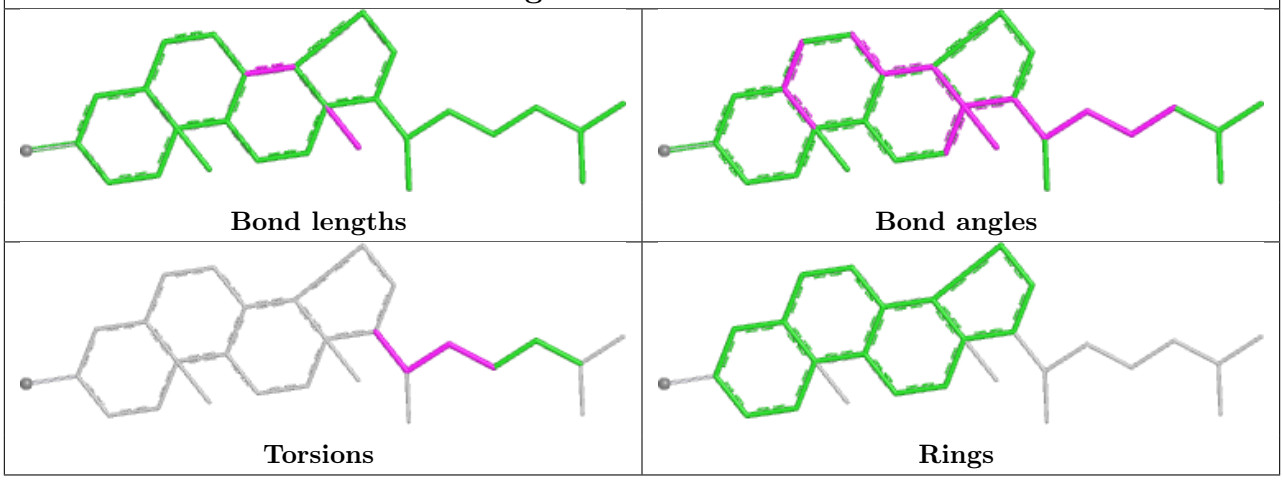
Ligand D21 R 702



Ligand CLR G 703



Ligand CLR R 701



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.

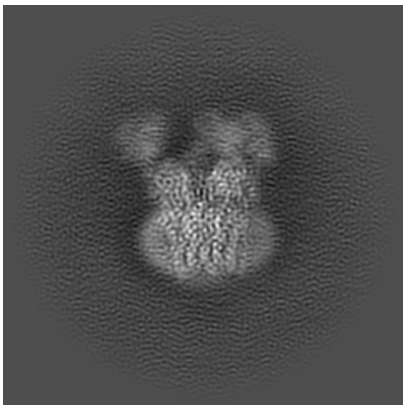
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-41893. These allow visual inspection of the internal detail of the map and identification of artifacts.

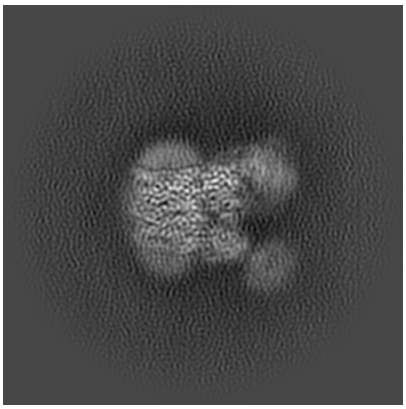
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

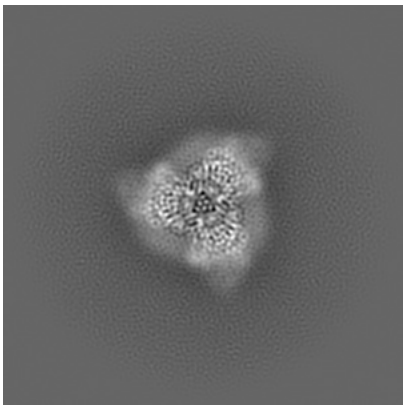
6.1.1 Primary map



X

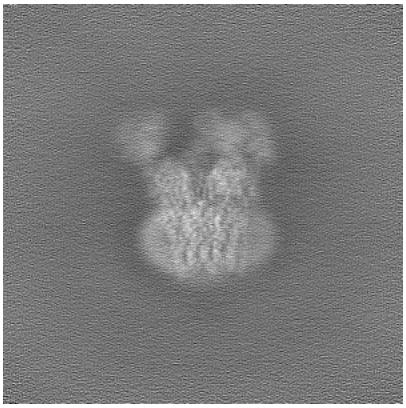


Y

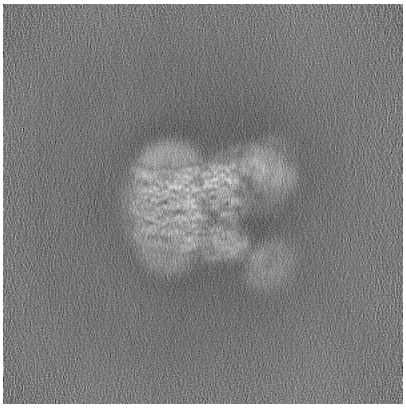


Z

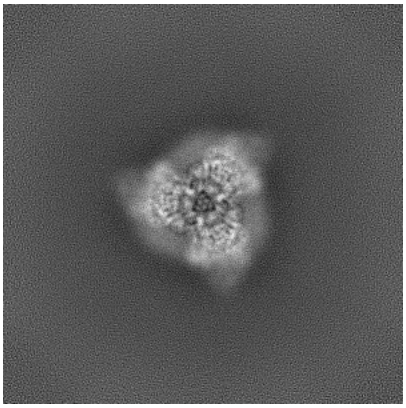
6.1.2 Raw map



X



Y

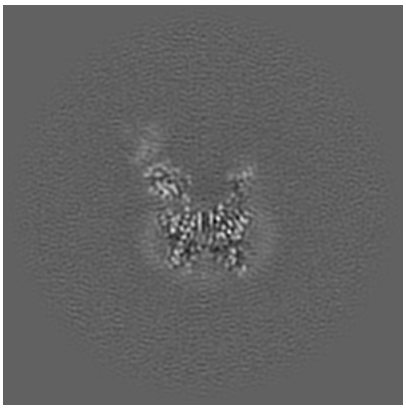


Z

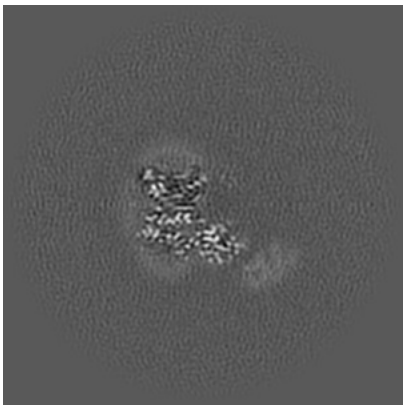
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

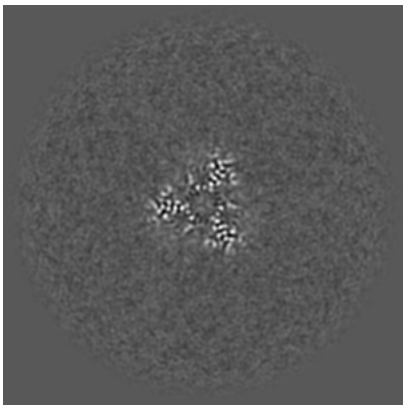
6.2.1 Primary map



X Index: 180

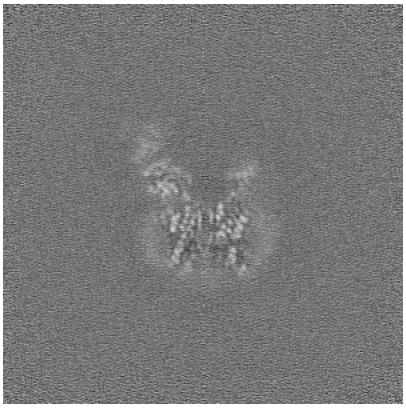


Y Index: 180

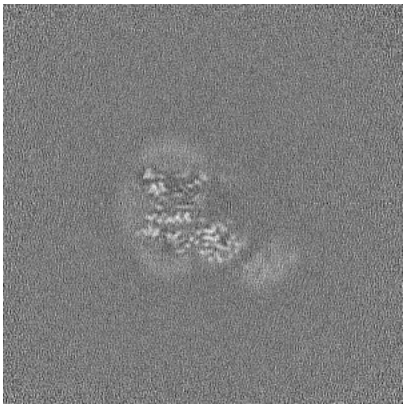


Z Index: 180

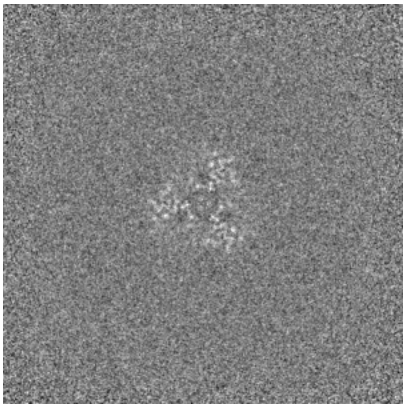
6.2.2 Raw map



X Index: 180



Y Index: 180

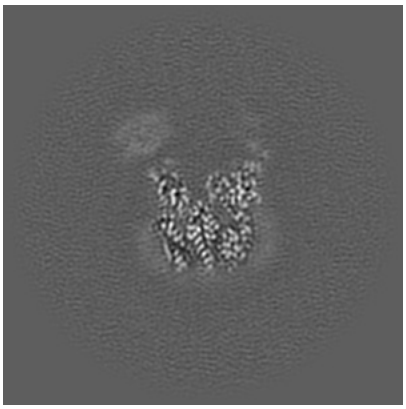


Z Index: 180

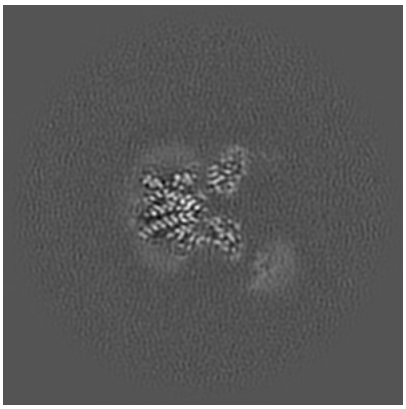
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

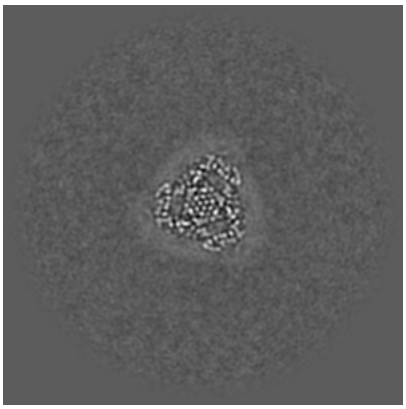
6.3.1 Primary map



X Index: 195

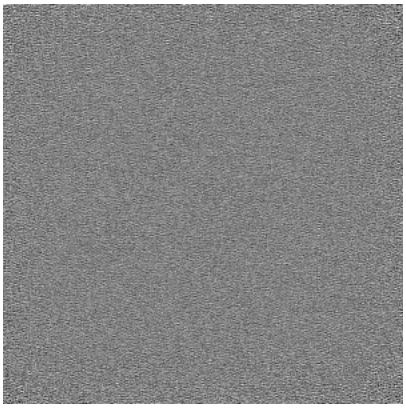


Y Index: 193

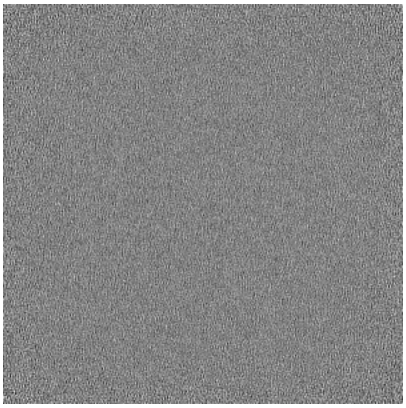


Z Index: 168

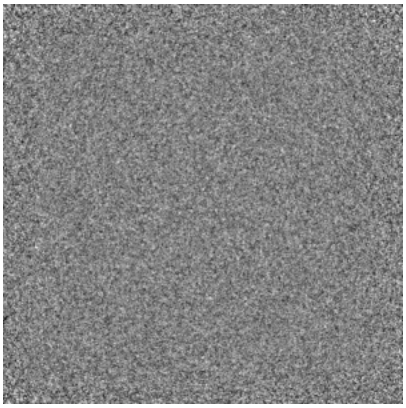
6.3.2 Raw map



X Index: 0



Y Index: 0

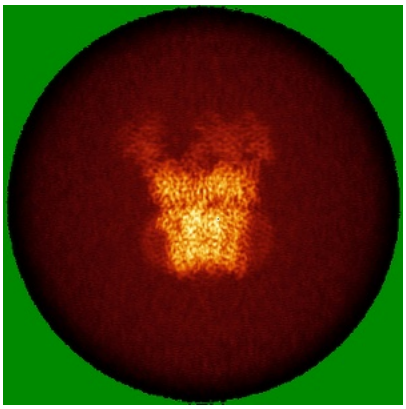


Z Index: 0

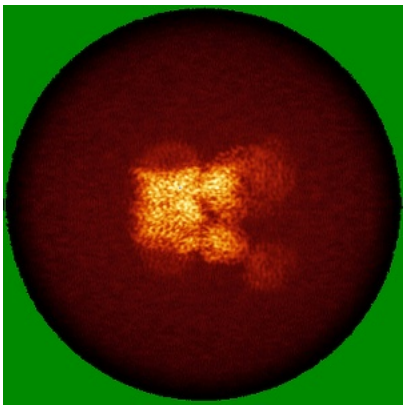
The images above show the largest variance slices of the map in three orthogonal directions.

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

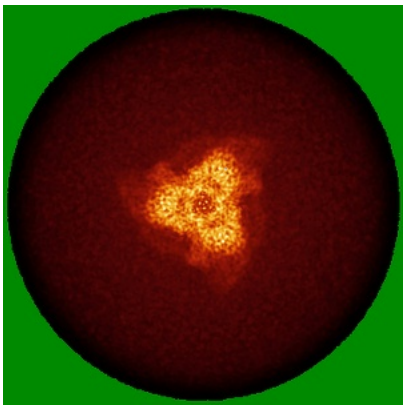
6.4.1 Primary map



X

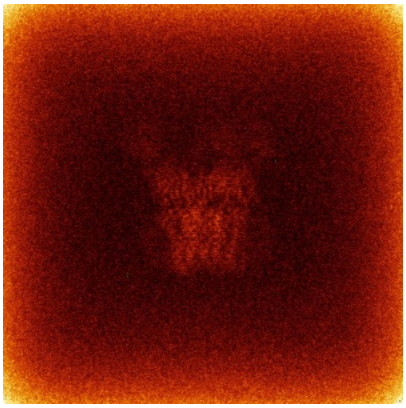


Y

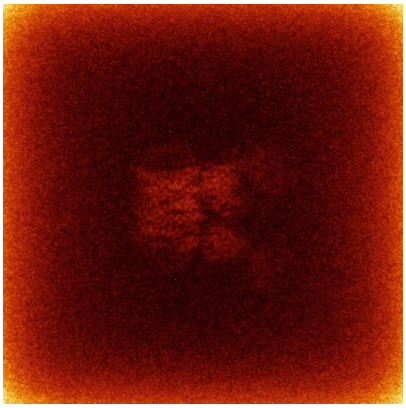


Z

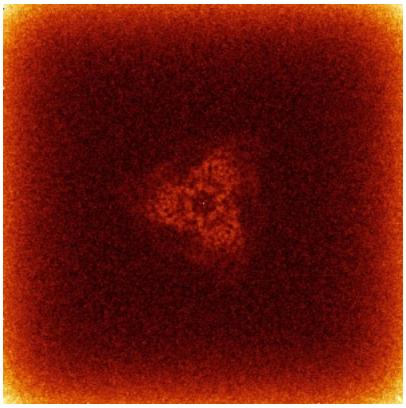
6.4.2 Raw map



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 0.15. 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.

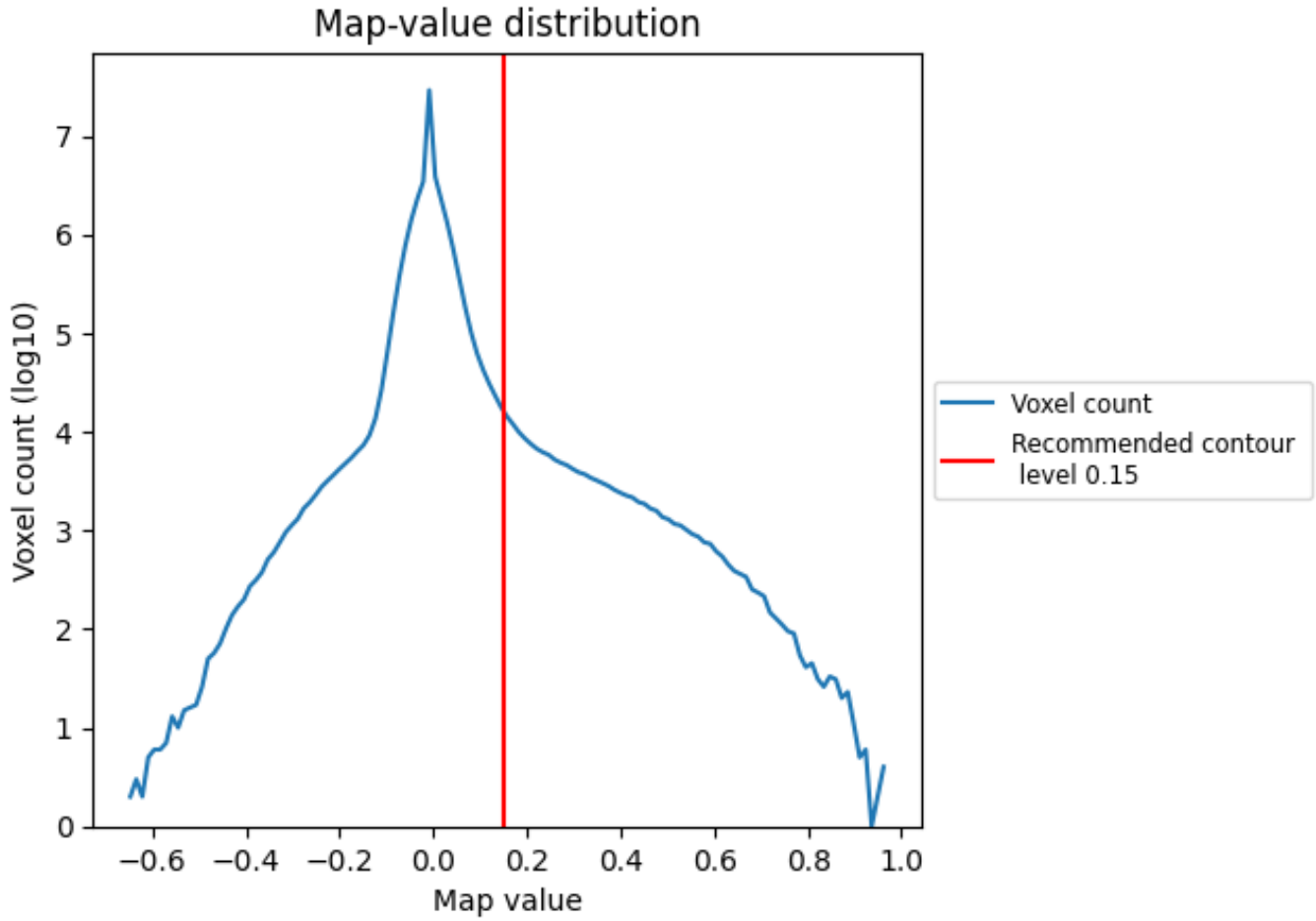
6.6 Mask visualisation (i)

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

7 Map analysis

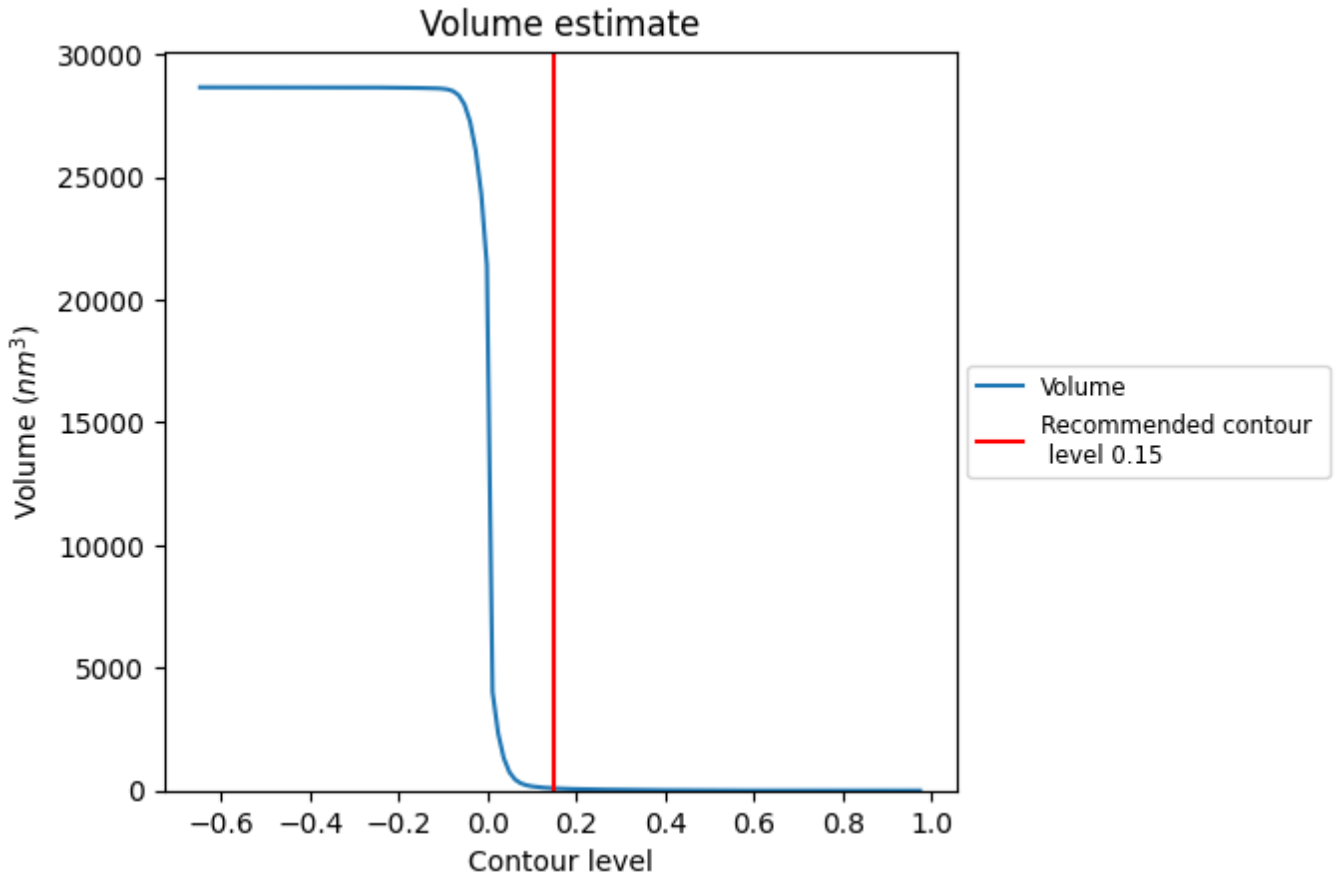
This section contains the results of statistical analysis of the map.

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.

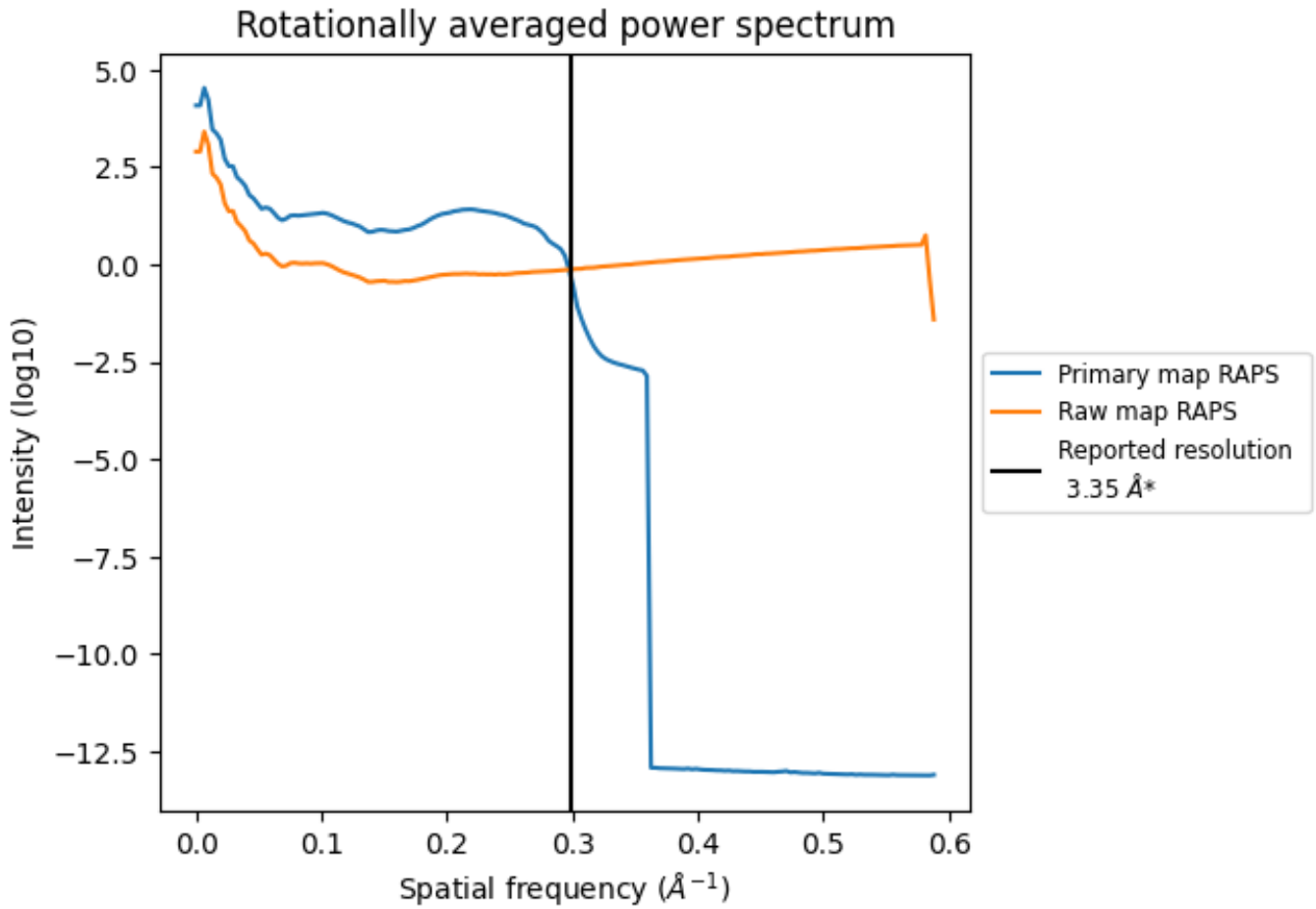
7.2 Volume estimate (i)



The volume at the recommended contour level is 92 nm^3; this corresponds to an approximate mass of 83 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)

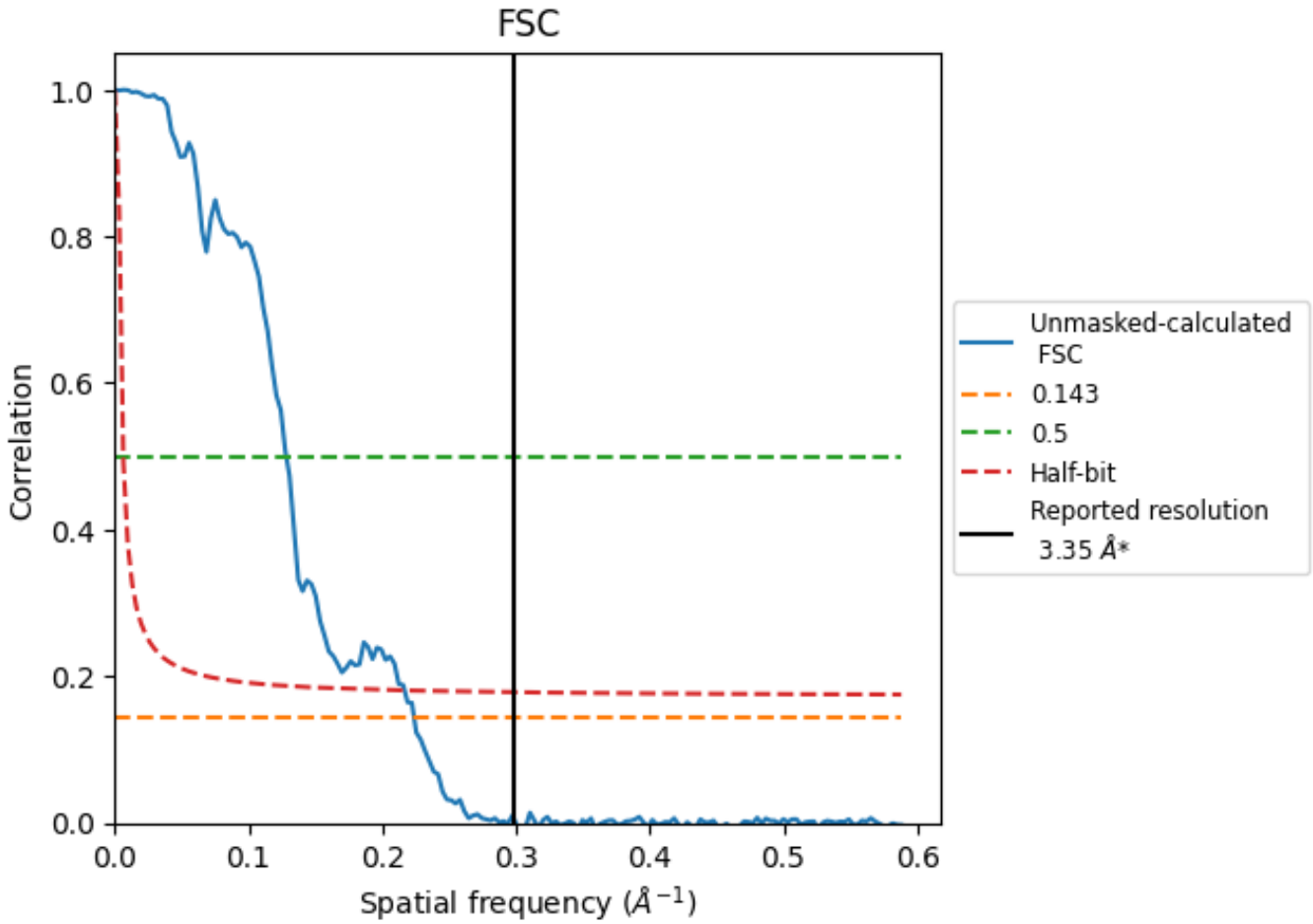


*Reported resolution corresponds to spatial frequency of 0.299 $Å^{-1}$

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 0.299 Å^{-1}

8.2 Resolution estimates (i)

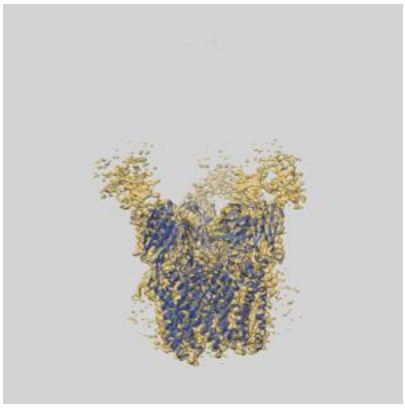
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.35	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.47	7.80	4.61

*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 4.47 differs from the reported value 3.35 by more than 10 %

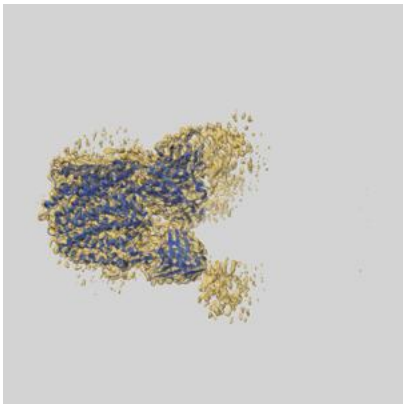
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-41893 and PDB model 8U4S. Per-residue inclusion information can be found in section 3 on page 8.

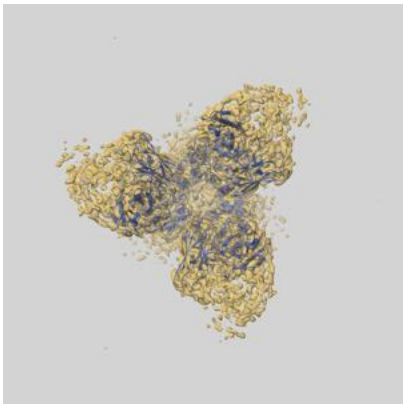
9.1 Map-model overlay



X



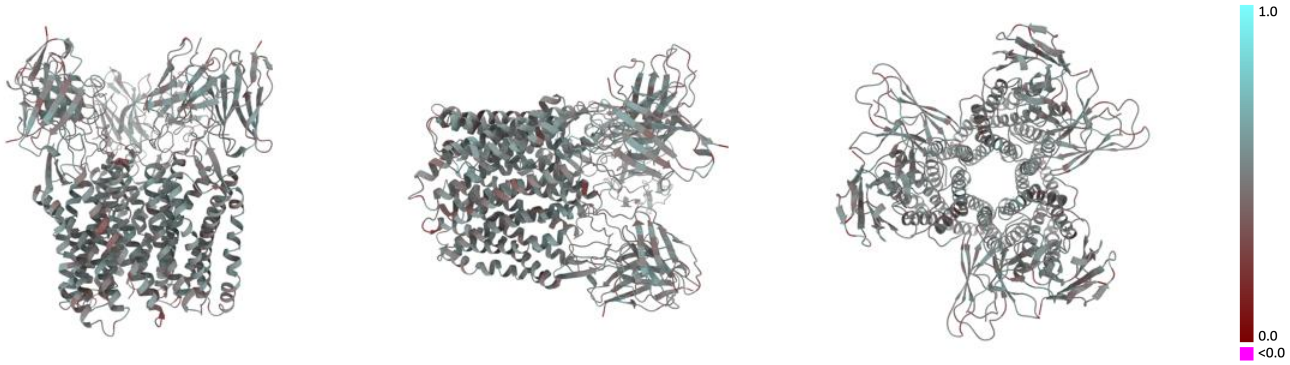
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.15 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.

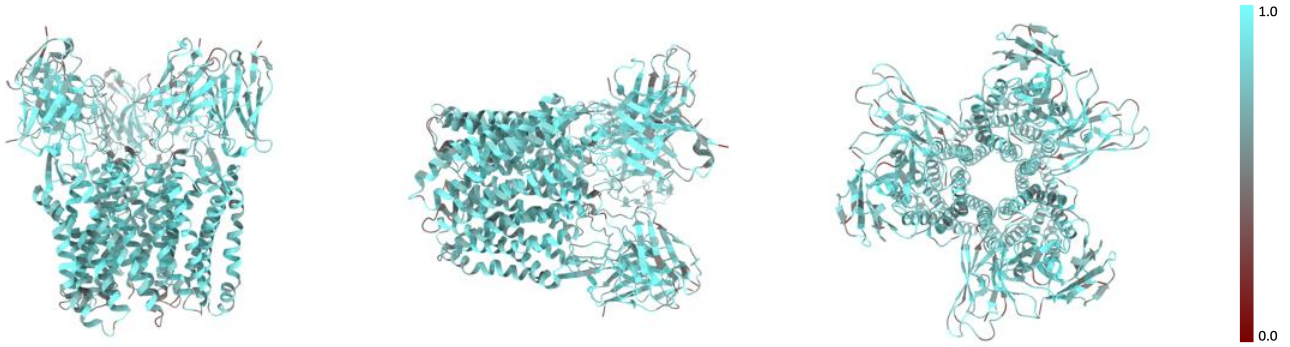
9.2 Q-score mapped to coordinate model



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.

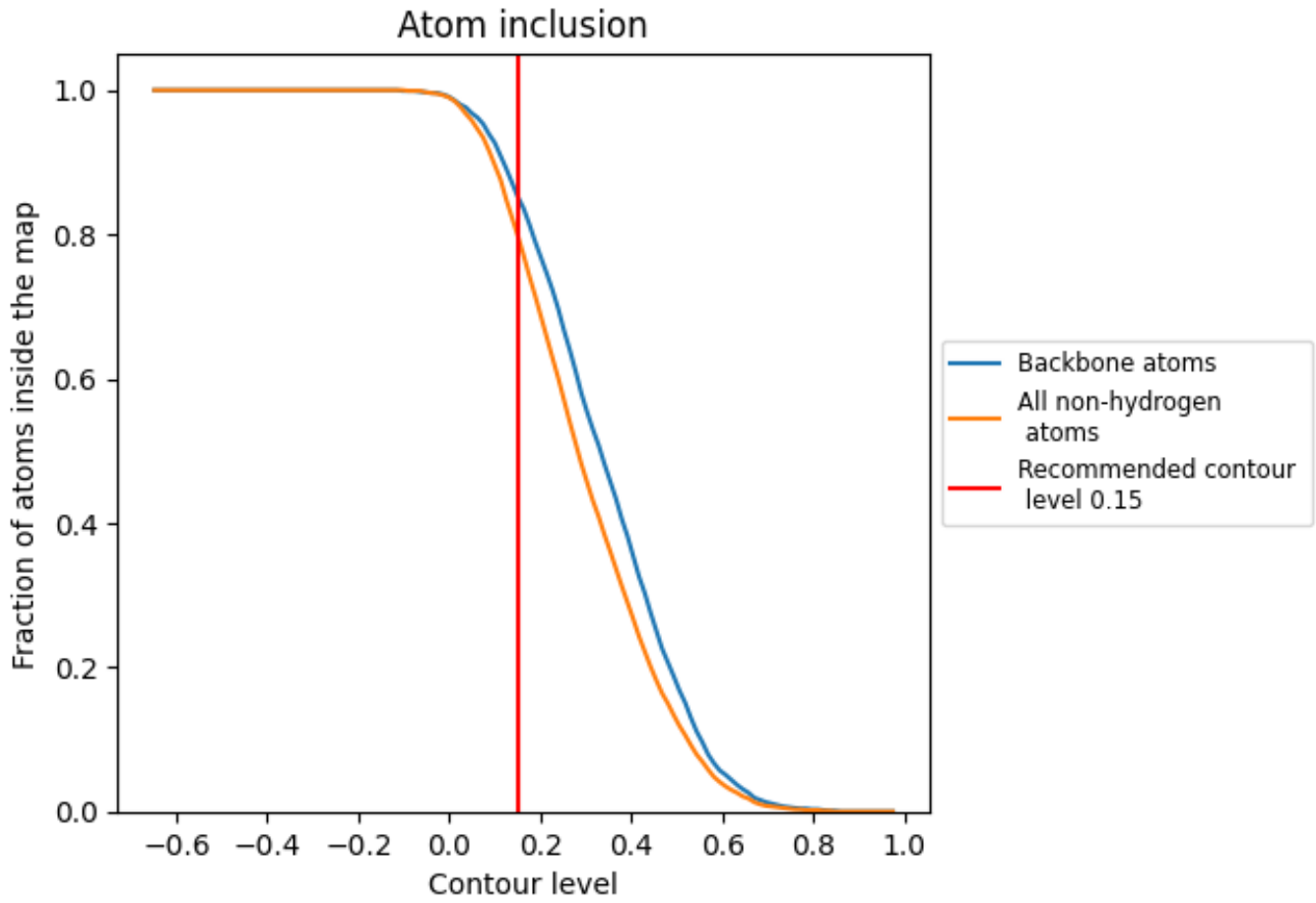
9.3 Atom inclusion mapped to coordinate model



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 (0.15).

9.4 Atom inclusion (i)



At the recommended contour level, 85% 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 (0.15) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8000	0.4900
A	0.7710	0.4820
B	0.7990	0.4870
C	0.8100	0.4930
E	0.7720	0.4810
F	0.7960	0.4870
G	0.8110	0.4930
H	0.7960	0.4900
L	0.7740	0.4820
R	0.8120	0.4950