Full wwPDB EM Validation Report

Nov 19, 2022 – 08:19 AM EST

PDB ID	:	3J2S
EMDB ID	:	EMD-5559
Title	:	Membrane-bound factor VIII light chain
Authors	:	Stoilova-Mcphie, S.; Lynch, G.C.; Ludtke, S.; Pettitt, B.M.
Deposited on	:	2012-12-21
Resolution	:	15.00 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/EMValidationReportHelp
with specific help available everywhere you see the (i) 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 (i)) were used in the production of this report:

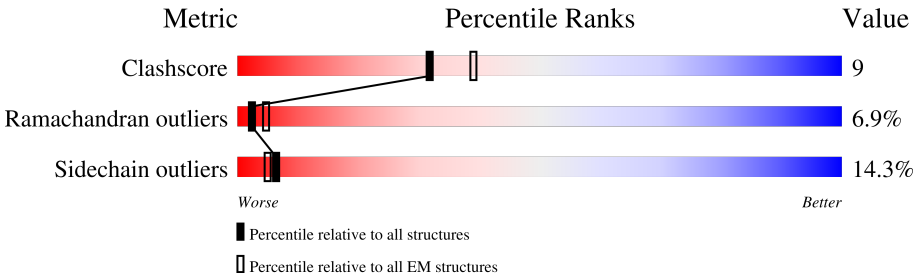
EMDB validation analysis	:	0.0.1.dev43
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

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 15.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	B	642	69% 22% 6%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 5125 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 Coagulation factor VIII light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	B	631	5125	3288	881	924	32	0	0

There is a discrepancy between the modelled and reference sequences:

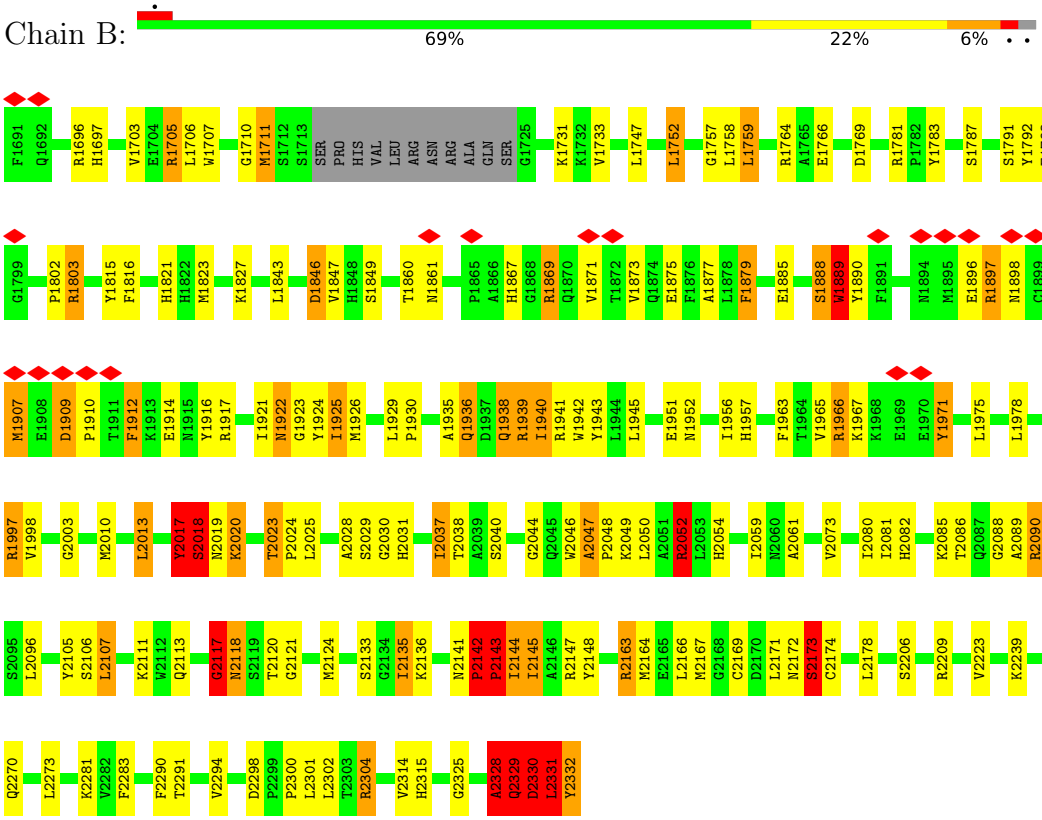
Chain	Residue	Modelled	Actual	Comment	Reference
B	1880	LEU	PHE	SEE REMARK 999	UNP P00451

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: Coagulation factor VIII light chain

Chain B:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=10°, rise=8 Å, axial sym=C1	Depositor
Number of segments used	Not provided	
Resolution determination method	Not provided	
CTF correction method	phase corrected based on first Thon ring	Depositor
Microscope	JEOL 2010F	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/Å^2$)	16	Depositor
Minimum defocus (nm)	-700	Depositor
Maximum defocus (nm)	-4400	Depositor
Magnification	52000	Depositor
Image detector	GATAN ULTRASCAN 4000 (4k x 4k)	Depositor
Maximum map value	0.019	Depositor
Minimum map value	-0.010	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.001	Depositor
Map size (Å)	165.3, 156.6, 81.200005	wwPDB
Map dimensions	57, 54, 28	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	2.9, 2.9, 2.9	Depositor

5 Model quality (i)

5.1 Standard geometry (i)

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	B	0.80	3/5270 (0.1%)	1.32	40/7136 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	35

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	2173	SER	C-N	5.49	1.46	1.34
1	B	2331	LEU	C-N	5.11	1.45	1.34
1	B	2329	GLN	N-CA	5.00	1.56	1.46

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	2328	ALA	O-C-N	-15.44	97.99	122.70
1	B	2329	GLN	N-CA-CB	10.57	129.63	110.60
1	B	2018	SER	N-CA-CB	-9.36	96.46	110.50
1	B	2017	TYR	O-C-N	-8.72	108.74	122.70
1	B	2052	ARG	NE-CZ-NH1	8.56	124.58	120.30
1	B	2330	ASP	CA-C-N	8.29	135.44	117.20
1	B	2331	LEU	O-C-N	-7.84	110.16	122.70
1	B	2331	LEU	C-N-CA	7.82	141.25	121.70
1	B	2018	SER	CA-CB-OG	7.41	131.20	111.20
1	B	2290	PHE	C-N-CA	7.34	140.04	121.70
1	B	2173	SER	O-C-N	-7.28	111.06	122.70
1	B	2328	ALA	CA-C-N	-7.26	101.24	117.20
1	B	2330	ASP	O-C-N	-7.25	111.09	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	1888	SER	C-N-CA	7.03	139.26	121.70
1	B	2133	SER	CB-CA-C	-7.01	96.78	110.10
1	B	2142	PRO	CA-C-N	6.76	136.02	117.10
1	B	1802	PRO	C-N-CA	6.63	138.27	121.70
1	B	1936	GLN	CB-CG-CD	6.50	128.50	111.60
1	B	2020	LYS	N-CA-C	6.43	128.37	111.00
1	B	2330	ASP	C-N-CA	6.39	137.68	121.70
1	B	2019	ASN	C-N-CA	6.24	137.31	121.70
1	B	2143	PRO	CA-N-CD	-6.07	103.00	111.50
1	B	2328	ALA	N-CA-CB	-5.91	101.83	110.10
1	B	2142	PRO	N-CA-CB	-5.88	96.13	102.60
1	B	1869	ARG	NE-CZ-NH2	5.87	123.23	120.30
1	B	1993	ALA	N-CA-CB	-5.77	102.03	110.10
1	B	1924	TYR	CB-CG-CD1	-5.76	117.55	121.00
1	B	2018	SER	N-CA-C	5.68	126.33	111.00
1	B	2164	MET	CG-SD-CE	-5.60	91.23	100.20
1	B	2172	ASN	C-N-CA	5.58	135.65	121.70
1	B	1963	PHE	CB-CG-CD1	5.58	124.70	120.80
1	B	1711	MET	CG-SD-CE	5.54	109.06	100.20
1	B	2147	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	B	1889	TRP	C-N-CA	5.49	135.42	121.70
1	B	2331	LEU	CA-C-N	5.36	128.99	117.20
1	B	1941	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	B	2018	SER	C-N-CA	5.24	134.79	121.70
1	B	2209	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	B	1963	PHE	CB-CG-CD2	-5.05	117.27	120.80
1	B	2047	ALA	N-CA-CB	-5.00	103.10	110.10

There are no chirality outliers.

All (35) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	1705	ARG	Peptide
1	B	1803	ARG	Mainchain
1	B	1846	ASP	Mainchain
1	B	1869	ARG	Sidechain
1	B	1885	GLU	Peptide
1	B	1889	TRP	Mainchain
1	B	1909	ASP	Peptide
1	B	1917	ARG	Sidechain,Peptide
1	B	1925	ILE	Peptide
1	B	1939	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	B	1971	TYR	Sidechain
1	B	2017	TYR	Mainchain
1	B	2018	SER	Peptide
1	B	2044	GLY	Peptide
1	B	2048	PRO	Peptide
1	B	2052	ARG	Sidechain
1	B	2090	ARG	Sidechain
1	B	2117	GLY	Peptide
1	B	2118	ASN	Mainchain
1	B	2141	ASN	Peptide
1	B	2142	PRO	Peptide
1	B	2148	TYR	Sidechain
1	B	2171	LEU	Peptide
1	B	2173	SER	Peptide,Mainchain
1	B	2174	CYS	Peptide
1	B	2298	ASP	Peptide
1	B	2304	ARG	Sidechain
1	B	2325	GLY	Peptide
1	B	2328	ALA	Peptide,Mainchain
1	B	2330	ASP	Peptide,Mainchain
1	B	2331	LEU	Peptide

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	B	5125	0	4995	96	0
All	All	5125	0	4995	96	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 9.

All (96) 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:2013:LEU:HG	1:B:2054:HIS:CE1	2.02	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2304:ARG:HH11	1:B:2331:LEU:HD11	1.31	0.93
1:B:1731:LYS:NZ	1:B:1889:TRP:CZ2	2.36	0.92
1:B:2018:SER:HB2	1:B:2023:THR:OG1	1.68	0.92
1:B:1792:TYR:HB3	1:B:1803:ARG:NH1	1.87	0.89
1:B:2107:LEU:HD23	1:B:2302:LEU:HD13	1.58	0.86
1:B:2304:ARG:NH1	1:B:2331:LEU:HD11	1.92	0.84
1:B:1781:ARG:NH2	1:B:1888:SER:HA	1.92	0.83
1:B:2304:ARG:HH11	1:B:2331:LEU:CD1	1.92	0.82
1:B:1781:ARG:HH22	1:B:1888:SER:HA	1.42	0.81
1:B:2118:ASN:HB3	1:B:2332:TYR:CE1	2.17	0.79
1:B:1936:GLN:HG2	1:B:2018:SER:OG	1.84	0.78
1:B:2013:LEU:CD2	1:B:2054:HIS:NE2	2.49	0.76
1:B:2018:SER:HB2	1:B:2023:THR:CB	2.15	0.75
1:B:2107:LEU:HD23	1:B:2302:LEU:CD1	2.16	0.74
1:B:2018:SER:CB	1:B:2023:THR:OG1	2.36	0.73
1:B:2013:LEU:CG	1:B:2054:HIS:CE1	2.72	0.72
1:B:1936:GLN:CG	1:B:2018:SER:OG	2.39	0.71
1:B:2013:LEU:HD21	1:B:2054:HIS:CD2	2.26	0.70
1:B:1792:TYR:CB	1:B:1803:ARG:NH1	2.54	0.70
1:B:1792:TYR:HB3	1:B:1803:ARG:HH12	1.56	0.69
1:B:1936:GLN:HB3	1:B:2018:SER:OG	1.92	0.69
1:B:2017:TYR:C	1:B:2018:SER:OG	2.27	0.67
1:B:2013:LEU:HD23	1:B:2054:HIS:NE2	2.13	0.63
1:B:1781:ARG:HH11	1:B:1783:TYR:HE1	1.48	0.62
1:B:1897:ARG:HE	1:B:1898:ASN:CG	2.03	0.62
1:B:2018:SER:HB2	1:B:2023:THR:HG21	1.81	0.62
1:B:2089:ALA:O	1:B:2096:LEU:HD23	2.00	0.62
1:B:1781:ARG:HD2	1:B:1783:TYR:CE1	2.34	0.62
1:B:1995:ILE:HB	1:B:2085:LYS:HZ3	1.64	0.61
1:B:2013:LEU:CD2	1:B:2054:HIS:CE1	2.84	0.61
1:B:1764:ARG:NH2	1:B:1875:GLU:OE1	2.33	0.60
1:B:2018:SER:HB2	1:B:2023:THR:CG2	2.30	0.60
1:B:2142:PRO:HB2	1:B:2330:ASP:C	2.22	0.60
1:B:1792:TYR:CB	1:B:1803:ARG:HH11	2.14	0.59
1:B:2107:LEU:HD22	1:B:2300:PRO:O	2.02	0.59
1:B:1936:GLN:CB	1:B:2018:SER:OG	2.51	0.59
1:B:2120:THR:HG23	1:B:2121:GLY:H	1.68	0.58
1:B:2013:LEU:CD2	1:B:2054:HIS:CD2	2.86	0.58
1:B:2118:ASN:CB	1:B:2332:TYR:CE1	2.86	0.57
1:B:1781:ARG:NH1	1:B:1846:ASP:OD1	2.38	0.56
1:B:1781:ARG:HH22	1:B:1888:SER:CA	2.15	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1781:ARG:HD2	1:B:1783:TYR:HE1	1.73	0.52
1:B:1846:ASP:HA	1:B:1849:SER:OG	2.09	0.52
1:B:1936:GLN:HG2	1:B:2018:SER:CB	2.39	0.52
1:B:2330:ASP:O	1:B:2331:LEU:CD2	2.59	0.51
1:B:1752:LEU:HD13	1:B:1752:LEU:H	1.76	0.51
1:B:2145:ILE:HD12	1:B:2302:LEU:HG	1.92	0.51
1:B:2082:HIS:HA	1:B:2329:GLN:CD	2.31	0.50
1:B:2145:ILE:CG1	1:B:2328:ALA:HB2	2.41	0.50
1:B:1897:ARG:NE	1:B:1898:ASN:OD1	2.43	0.49
1:B:2143:PRO:HA	1:B:2329:GLN:HB2	1.94	0.49
1:B:2018:SER:OG	1:B:2023:THR:OG1	2.26	0.49
1:B:2017:TYR:CE2	1:B:2029:SER:HB2	2.47	0.49
1:B:2018:SER:C	1:B:2023:THR:HG21	2.33	0.49
1:B:2082:HIS:HA	1:B:2329:GLN:NE2	2.28	0.48
1:B:2025:LEU:CD1	1:B:2081:ILE:HG22	2.42	0.48
1:B:1759:LEU:HD11	1:B:1879:PHE:CE2	2.48	0.48
1:B:2281:LYS:HG2	1:B:2283:PHE:CZ	2.48	0.47
1:B:2025:LEU:HD13	1:B:2081:ILE:HG22	1.96	0.47
1:B:2145:ILE:HG23	1:B:2328:ALA:CB	2.45	0.47
1:B:2145:ILE:HG23	1:B:2328:ALA:HB2	1.95	0.47
1:B:1731:LYS:O	1:B:1889:TRP:HZ2	1.98	0.46
1:B:2145:ILE:CD1	1:B:2302:LEU:HG	2.45	0.46
1:B:1696:ARG:NH2	1:B:1766:GLU:OE1	2.26	0.46
1:B:2052:ARG:HH11	1:B:2052:ARG:HG3	1.80	0.46
1:B:2013:LEU:HG	1:B:2054:HIS:HE1	1.74	0.46
1:B:2314:VAL:HG13	1:B:2315:HIS:N	2.31	0.45
1:B:2117:GLY:H	1:B:2124:MET:HB2	1.81	0.45
1:B:1935:ALA:HA	1:B:2017:TYR:O	2.17	0.45
1:B:2013:LEU:HD21	1:B:2054:HIS:CG	2.51	0.45
1:B:2080:ILE:HG12	1:B:2328:ALA:HA	1.99	0.45
1:B:2143:PRO:CG	1:B:2144:ILE:H	2.30	0.45
1:B:2088:GLY:O	1:B:2163:ARG:HD3	2.17	0.44
1:B:2167:MET:HA	1:B:2167:MET:HE2	1.99	0.44
1:B:2023:THR:HG23	1:B:2024:PRO:O	2.18	0.44
1:B:1731:LYS:NZ	1:B:1889:TRP:HZ2	2.06	0.43
1:B:2143:PRO:HG2	1:B:2144:ILE:H	1.82	0.43
1:B:2013:LEU:HD21	1:B:2054:HIS:CE1	2.51	0.43
1:B:1936:GLN:HG2	1:B:2018:SER:HB3	2.00	0.43
1:B:1966:ARG:H	1:B:1966:ARG:HD3	1.84	0.43
1:B:2143:PRO:O	1:B:2329:GLN:NE2	2.52	0.43
1:B:1940:ILE:HD11	1:B:1942:TRP:CE2	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1997:ARG:HB2	1:B:2013:LEU:HA	2.01	0.42
1:B:1877:ALA:HA	1:B:1943:TYR:HB2	2.02	0.42
1:B:1923:GLY:H	1:B:1925:ILE:HG12	1.84	0.42
1:B:2142:PRO:HG2	1:B:2331:LEU:HB2	2.01	0.42
1:B:1936:GLN:HA	1:B:1990:PRO:HG2	2.01	0.42
1:B:1897:ARG:NH1	1:B:1897:ARG:HB3	2.33	0.42
1:B:2314:VAL:HG13	1:B:2315:HIS:H	1.85	0.41
1:B:1897:ARG:NH2	1:B:1898:ASN:OD1	2.52	0.41
1:B:2105:TYR:CG	1:B:2144:ILE:HG23	2.55	0.41
1:B:2330:ASP:O	1:B:2331:LEU:HD23	2.20	0.41
1:B:1921:ILE:O	1:B:1922:ASN:HB2	2.21	0.41
1:B:1935:ALA:HB3	1:B:1938:GLN:HB2	2.03	0.40
1:B:2145:ILE:HD13	1:B:2239:LYS:HB2	2.03	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	B	627/642 (98%)	471 (75%)	113 (18%)	43 (7%)	1 15

All (43) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1705	ARG
1	B	1706	LEU
1	B	1710	GLY
1	B	1757	GLY
1	B	1871	VAL
1	B	1889	TRP
1	B	1890	TYR
1	B	1910	PRO

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Mol	Chain	Res	Type
1	B	1912	PHE
1	B	1922	ASN
1	B	1930	PRO
1	B	2003	GLY
1	B	2018	SER
1	B	2049	LYS
1	B	2061	ALA
1	B	2092	LYS
1	B	2135	ILE
1	B	2142	PRO
1	B	2143	PRO
1	B	2291	THR
1	B	1794	GLU
1	B	1952	ASN
1	B	2020	LYS
1	B	2037	ILE
1	B	2038	THR
1	B	2117	GLY
1	B	2169	CYS
1	B	2331	LEU
1	B	1791	SER
1	B	1821	HIS
1	B	1907	MET
1	B	2252	LEU
1	B	2028	ALA
1	B	2047	ALA
1	B	1867	HIS
1	B	1906	GLN
1	B	1967	LYS
1	B	2040	SER
1	B	2144	ILE
1	B	2173	SER
1	B	1873	VAL
1	B	2030	GLY
1	B	2059	ILE

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	B	560/570 (98%)	480 (86%)	80 (14%)	3 16

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	1697	HIS
1	B	1703	VAL
1	B	1707	TRP
1	B	1711	MET
1	B	1733	VAL
1	B	1747	LEU
1	B	1752	LEU
1	B	1758	LEU
1	B	1759	LEU
1	B	1769	ASP
1	B	1787	SER
1	B	1793	GLU
1	B	1798	GLN
1	B	1815	TYR
1	B	1816	PHE
1	B	1823	MET
1	B	1827	LYS
1	B	1843	LEU
1	B	1847	VAL
1	B	1860	THR
1	B	1861	ASN
1	B	1879	PHE
1	B	1889	TRP
1	B	1896	GLU
1	B	1897	ARG
1	B	1900	ARG
1	B	1907	MET
1	B	1909	ASP
1	B	1912	PHE
1	B	1914	GLU
1	B	1916	TYR
1	B	1926	MET
1	B	1929	LEU
1	B	1938	GLN
1	B	1939	ARG
1	B	1940	ILE

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Mol	Chain	Res	Type
1	B	1945	LEU
1	B	1951	GLU
1	B	1956	ILE
1	B	1957	HIS
1	B	1965	VAL
1	B	1966	ARG
1	B	1971	TYR
1	B	1975	LEU
1	B	1978	LEU
1	B	1997	ARG
1	B	1998	VAL
1	B	2010	MET
1	B	2013	LEU
1	B	2023	THR
1	B	2031	HIS
1	B	2037	ILE
1	B	2046	TRP
1	B	2050	LEU
1	B	2052	ARG
1	B	2073	VAL
1	B	2086	THR
1	B	2090	ARG
1	B	2091	GLN
1	B	2092	LYS
1	B	2094	SER
1	B	2106	SER
1	B	2107	LEU
1	B	2111	LYS
1	B	2113	GLN
1	B	2135	ILE
1	B	2136	LYS
1	B	2145	ILE
1	B	2163	ARG
1	B	2166	LEU
1	B	2178	LEU
1	B	2206	SER
1	B	2223	VAL
1	B	2249	LYS
1	B	2270	GLN
1	B	2273	LEU
1	B	2294	VAL
1	B	2301	LEU

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Mol	Chain	Res	Type
1	B	2329	GLN
1	B	2332	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	1729	GLN
1	B	1778	GLN
1	B	1867	HIS
1	B	1874	GLN
1	B	2036	GLN
1	B	2129	ASN
1	B	2315	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

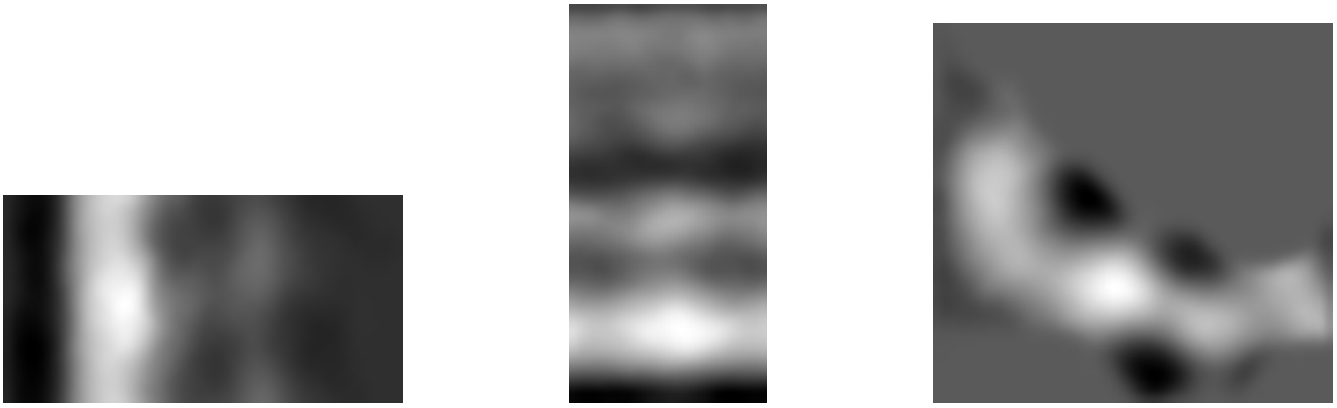
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-5559. 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

6.1.1 Primary map



X Y Z

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

6.2 Central slices

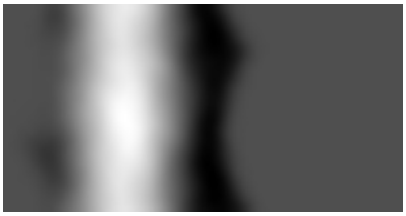
6.2.1 Primary map

X Index: 28 Y Index: 27 Z Index: 14

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

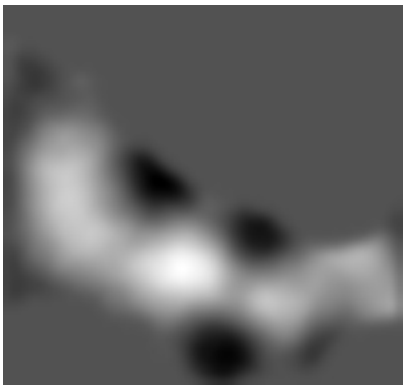
6.3 Largest variance slices (i)

6.3.1 Primary map



X Index: 24

Y Index: 18

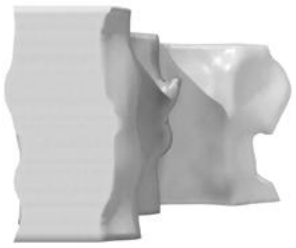


Z Index: 12

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

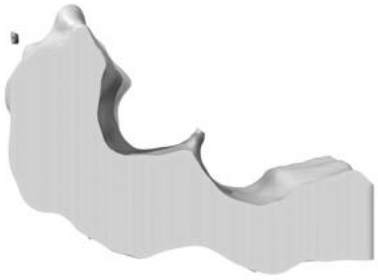
6.4 Orthogonal surface views (i)

6.4.1 Primary map



X

Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.001. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

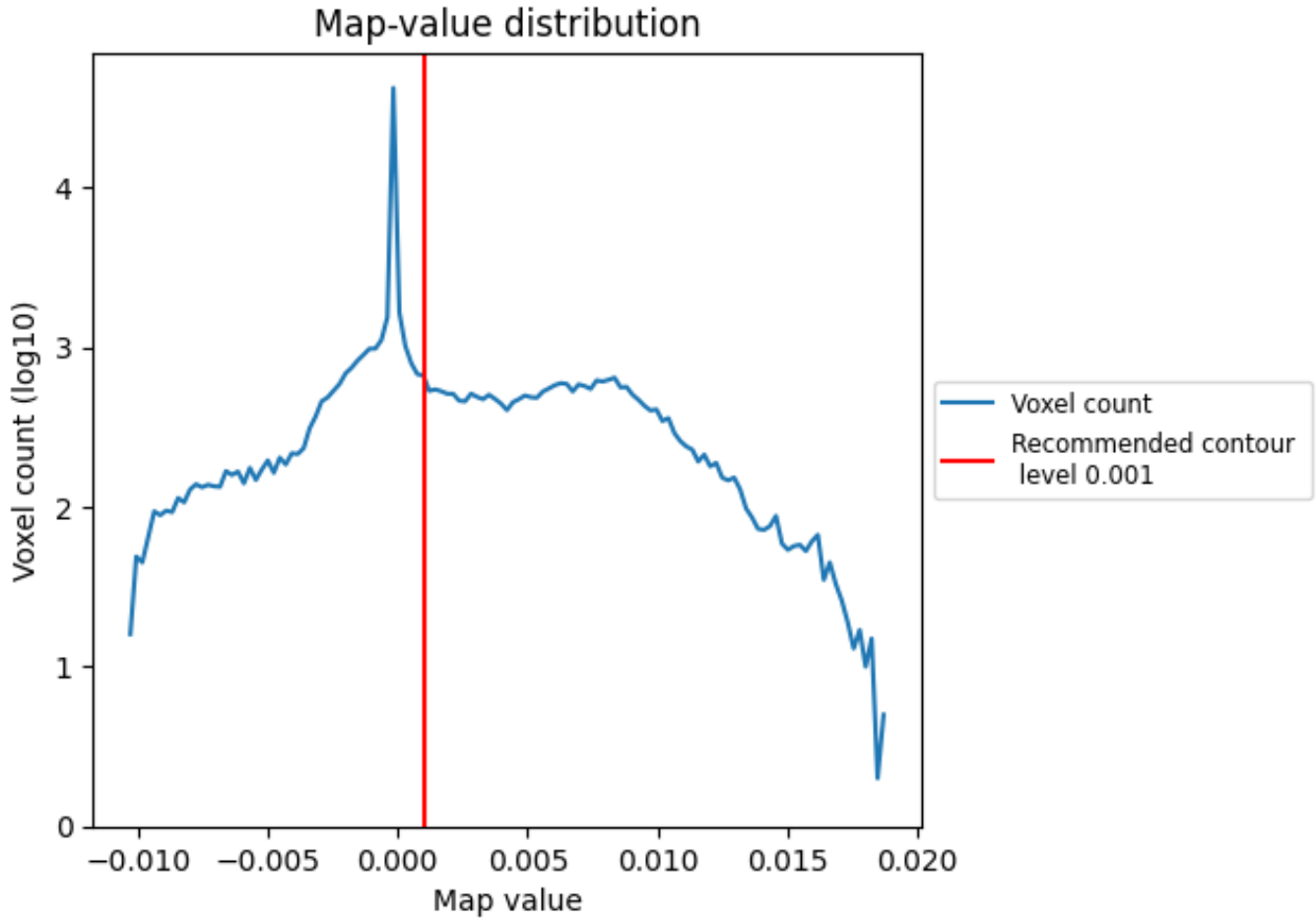
6.5 Mask visualisation (i)

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

7 Map analysis

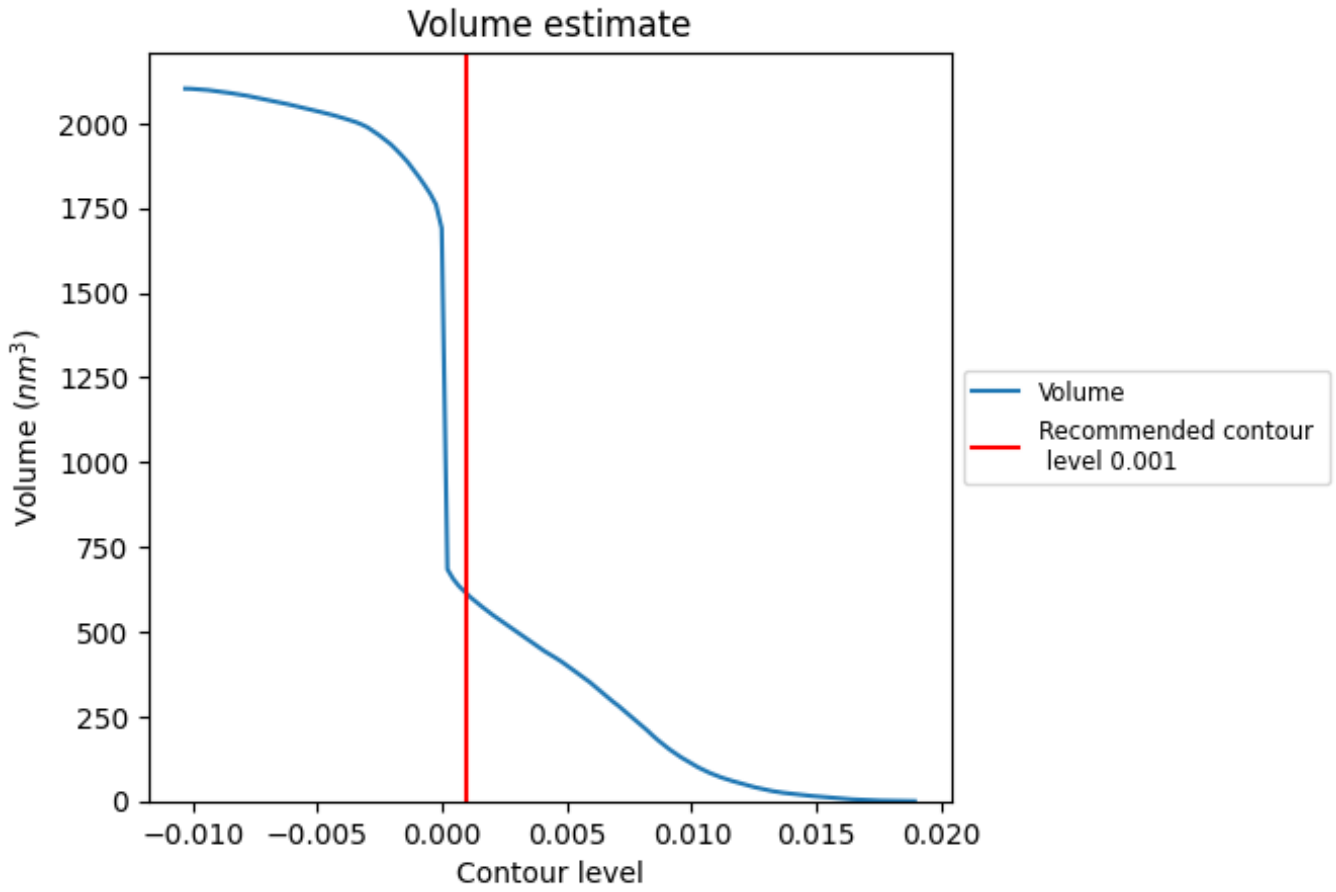
This section contains the results of statistical analysis of the map.

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



The volume at the recommended contour level is 611 nm^3; this corresponds to an approximate mass of 552 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

This section was not generated. The rotationally averaged power spectrum is only generated for cubic maps.

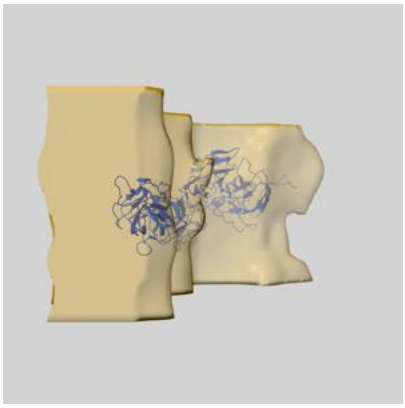
8 Fourier-Shell correlation (i)

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

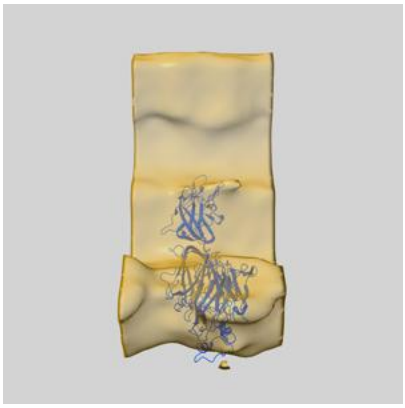
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-5559 and PDB model 3J2S. Per-residue inclusion information can be found in section 3 on page 4.

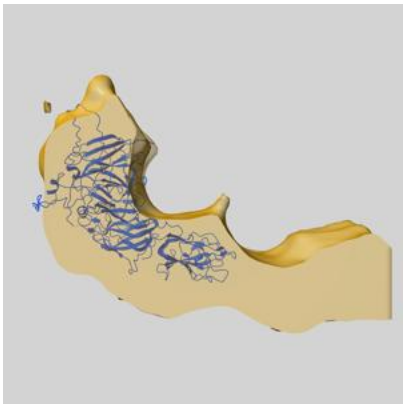
9.1 Map-model overlay



X



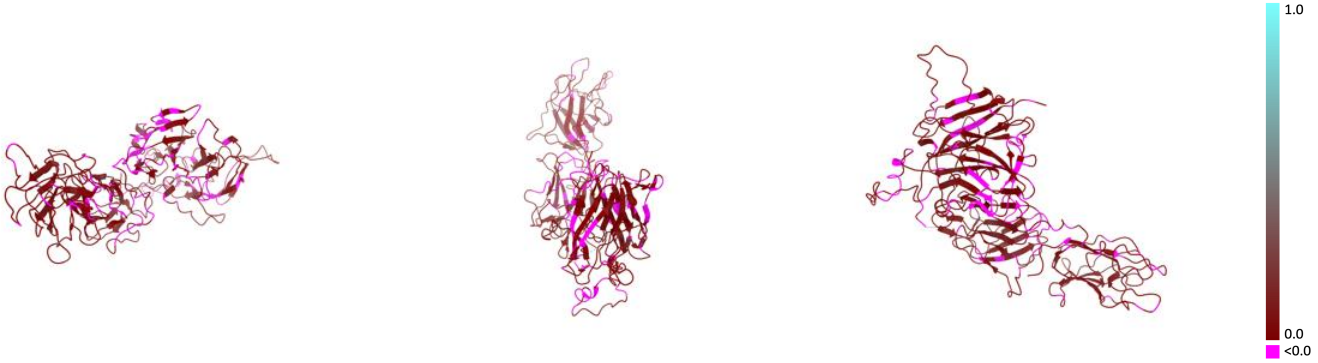
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.001 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.

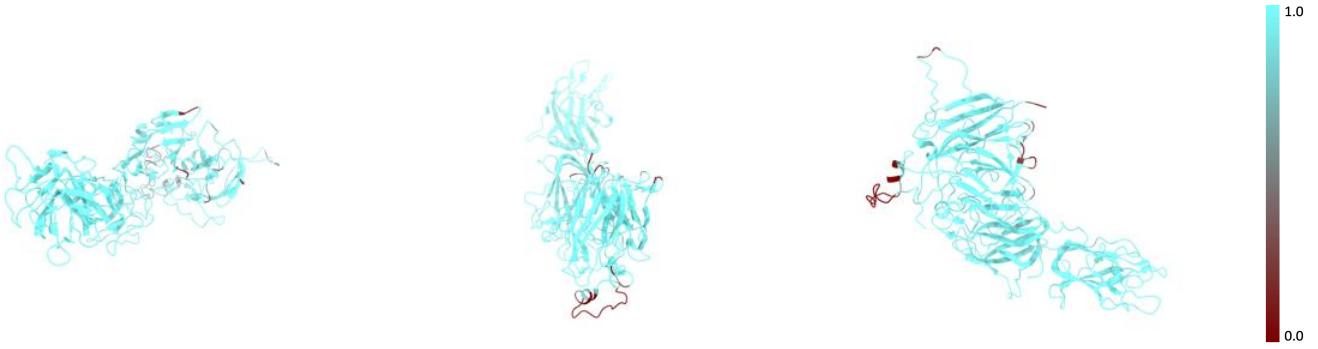
9.2 Q-score mapped to coordinate model



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.

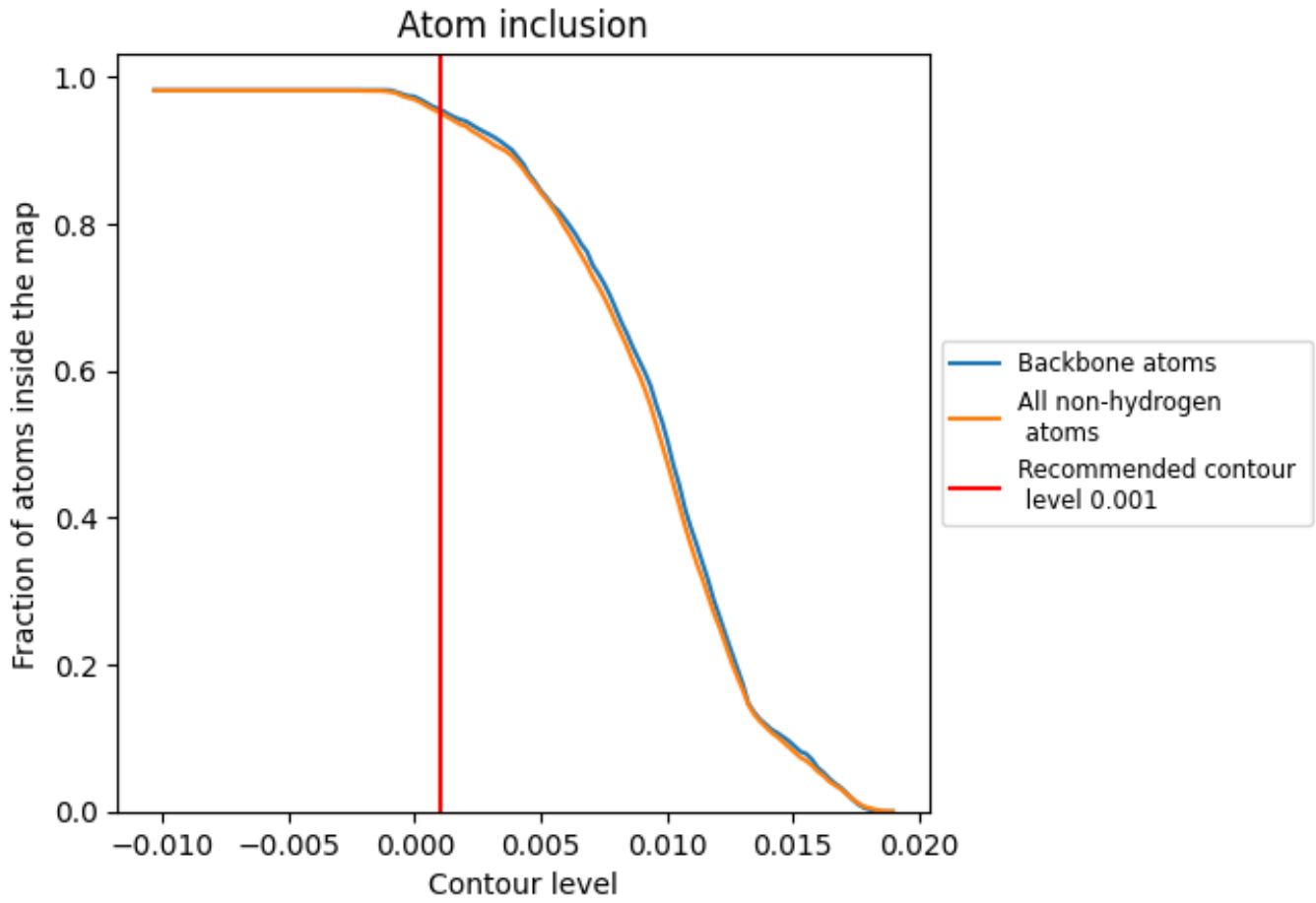
9.3 Atom inclusion mapped to coordinate model



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 (0.001).

9.4 Atom inclusion (i)



At the recommended contour level, 96% of all backbone atoms, 95% 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 (0.001) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9512	0.0430
B	0.9512	0.0430