Full wwPDB EM Validation Report (i)

Nov 29, 2022 – 07:22 PM JST

PDB ID	:	7WVU
EMDB ID	:	EMD-32858
Title	:	Cryo-EM structure of the human formyl peptide receptor 1 in complex with fMLF and Gi1
Authors	:	Zhu, Y.; Lin, X.; Zong, X.; Han, S.; Zhao, Q.; Wu, B.
Deposited on	:	2022-02-11
Resolution	:	3.30 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:

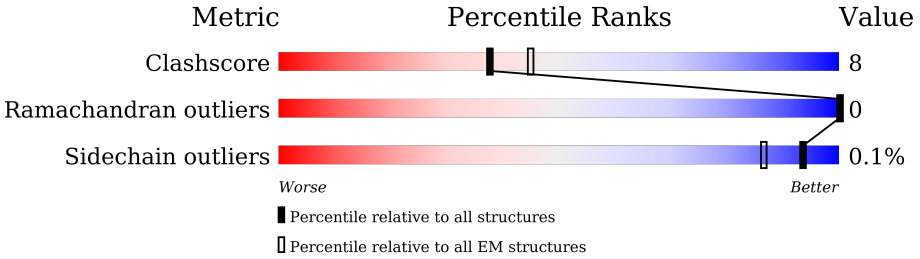
EMDB validation analysis	:	0.0.1.dev43
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (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 3.30 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	R	371	26% 61% 19% 20%
2	A	354	12% 53% 8% 38%
3	B	351	15% 78% 18% •
4	C	71	44% 69% 7% 24%
5	L	3	67% 33% 67%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 6862 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 fMet-Leu-Phe receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	R	298	2279	1520	377	368	14	0	0

There are 51 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	-1	GLY	-	expression tag	UNP P21462
R	0	ALA	-	expression tag	UNP P21462
R	1	PRO	-	expression tag	UNP P21462
R	322	GLU	-	expression tag	UNP P21462
R	323	PHE	-	expression tag	UNP P21462
R	324	LEU	-	expression tag	UNP P21462
R	325	GLU	-	expression tag	UNP P21462
R	326	VAL	-	expression tag	UNP P21462
R	327	LEU	-	expression tag	UNP P21462
R	328	PHE	-	expression tag	UNP P21462
R	329	GLN	-	expression tag	UNP P21462
R	330	GLY	-	expression tag	UNP P21462
R	331	PRO	-	expression tag	UNP P21462
R	332	GLY	-	expression tag	UNP P21462
R	333	SER	-	expression tag	UNP P21462
R	334	TRP	-	expression tag	UNP P21462
R	335	SER	-	expression tag	UNP P21462
R	336	HIS	-	expression tag	UNP P21462
R	337	PRO	-	expression tag	UNP P21462
R	338	GLN	-	expression tag	UNP P21462
R	339	PHE	-	expression tag	UNP P21462
R	340	GLU	-	expression tag	UNP P21462
R	341	LYS	-	expression tag	UNP P21462
R	342	GLY	-	expression tag	UNP P21462
R	343	SER	-	expression tag	UNP P21462
R	344	GLY	-	expression tag	UNP P21462
R	345	ALA	-	expression tag	UNP P21462
R	346	GLY	-	expression tag	UNP P21462

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Chain	Residue	Modelled	Actual	Comment	Reference
R	347	ALA	-	expression tag	UNP P21462
R	348	SER	-	expression tag	UNP P21462
R	349	ALA	-	expression tag	UNP P21462
R	350	GLY	-	expression tag	UNP P21462
R	351	SER	-	expression tag	UNP P21462
R	352	TRP	-	expression tag	UNP P21462
R	353	SER	-	expression tag	UNP P21462
R	354	HIS	-	expression tag	UNP P21462
R	355	PRO	-	expression tag	UNP P21462
R	356	GLN	-	expression tag	UNP P21462
R	357	PHE	-	expression tag	UNP P21462
R	358	GLU	-	expression tag	UNP P21462
R	359	LYS	-	expression tag	UNP P21462
R	360	GLY	-	expression tag	UNP P21462
R	361	SER	-	expression tag	UNP P21462
R	362	ASP	-	expression tag	UNP P21462
R	363	TYR	-	expression tag	UNP P21462
R	364	LYS	-	expression tag	UNP P21462
R	365	ASP	-	expression tag	UNP P21462
R	366	ASP	-	expression tag	UNP P21462
R	367	ASP	-	expression tag	UNP P21462
R	368	ASP	-	expression tag	UNP P21462
R	369	LYS	-	expression tag	UNP P21462

- Molecule 2 is a protein called Guanine nucleotide-binding protein G(i) subunit alpha-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	A	218	1714	1097	283	322	12	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	CYS	SER	engineered mutation	UNP P63096
A	202	THR	GLY	engineered mutation	UNP P63096
A	203	ALA	GLY	engineered mutation	UNP P63096
A	245	ALA	GLU	engineered mutation	UNP P63096
A	326	SER	ALA	engineered mutation	UNP P63096

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	B	336	2457	1526	430	481	20	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	MET	-	expression tag	UNP P62873
B	-9	HIS	-	expression tag	UNP P62873
B	-8	HIS	-	expression tag	UNP P62873
B	-7	HIS	-	expression tag	UNP P62873
B	-6	HIS	-	expression tag	UNP P62873
B	-5	HIS	-	expression tag	UNP P62873
B	-4	HIS	-	expression tag	UNP P62873
B	-3	GLY	-	expression tag	UNP P62873
B	-2	SER	-	expression tag	UNP P62873
B	-1	LEU	-	expression tag	UNP P62873
B	0	LEU	-	expression tag	UNP P62873
B	1	GLN	-	expression tag	UNP P62873

- Molecule 4 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	C	54	382	241	64	74	3	0	0

- Molecule 5 is a protein called FME-LEU-PHE.

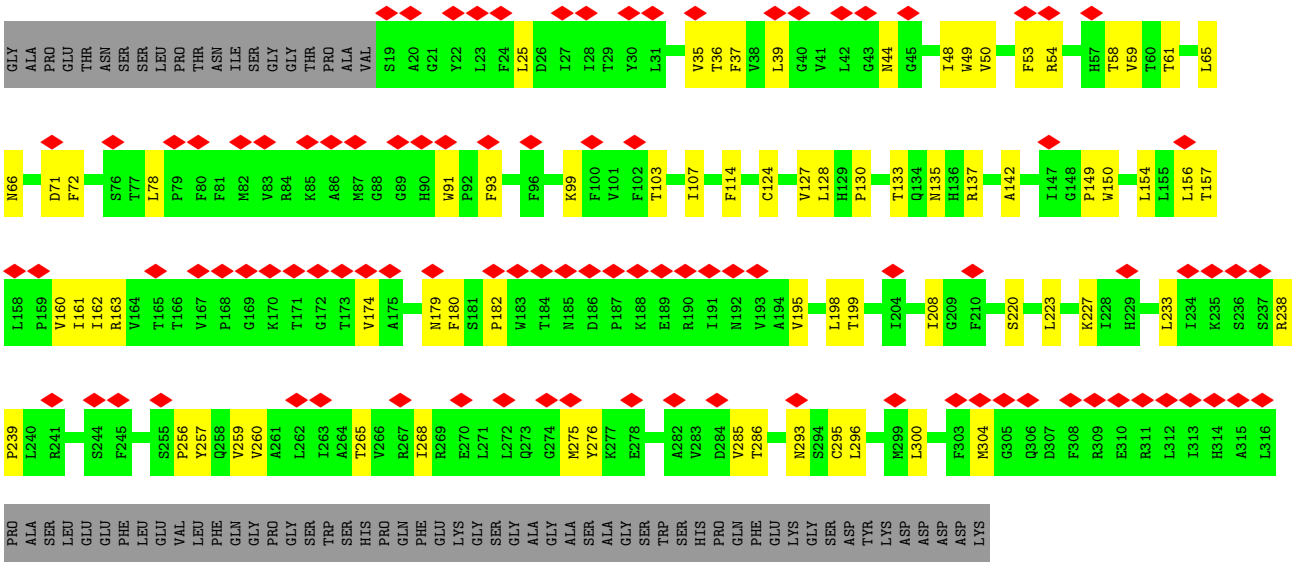
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	L	3	30	21	3	5	1	0	0

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.

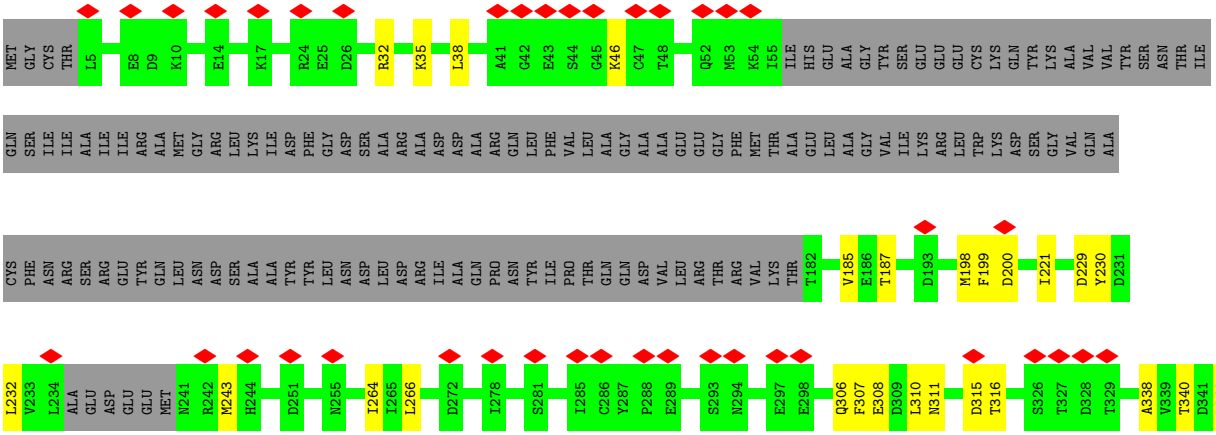
• Molecule 1: fMet-Leu-Phe receptor

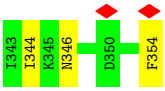
Chain R:



• Molecule 2: Guanine nucleotide-binding protein G(i) subunit alpha-1

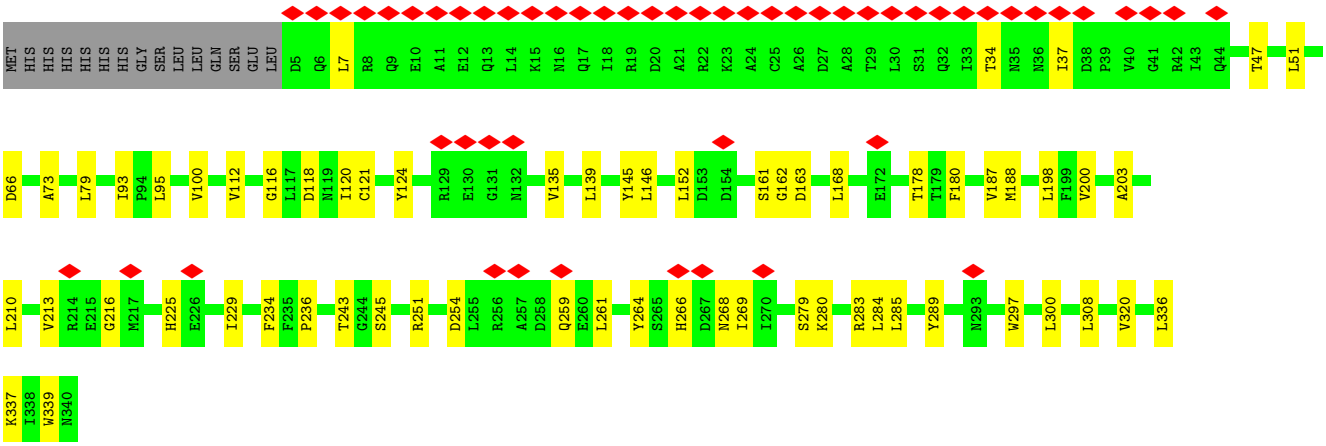
Chain A:





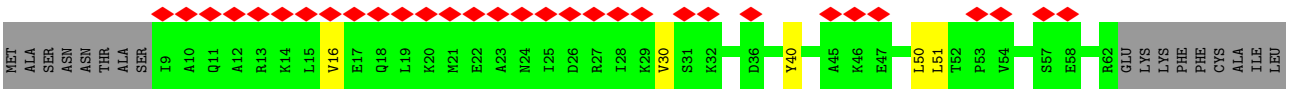
• Molecule 3: Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1

Chain B:



• Molecule 4: Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2

Chain C:



• Molecule 5: FME-LEU-PHE

Chain L:

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	1299041	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	2.1875	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.213	Depositor
Minimum map value	-0.135	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.0298	Depositor
Map size (Å)	267.52, 267.52, 267.52	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.045, 1.045, 1.045	Depositor

5 Model quality (i)

5.1 Standard geometry (i)

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

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	R	0.24	0/2341	0.39	0/3202
2	A	0.24	0/1743	0.39	0/2345
3	B	0.24	0/2504	0.45	0/3412
4	C	0.24	0/388	0.36	0/529
5	L	0.26	0/20	0.45	0/24
All	All	0.24	0/6996	0.41	0/9512

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	R	2279	0	2319	46	0
2	A	1714	0	1673	20	0
3	B	2457	0	2262	46	0
4	C	382	0	357	5	0
5	L	30	0	30	3	0
All	All	6862	0	6641	109	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 8.

All (109) 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)
3:B:254:ASP:HB2	3:B:261:LEU:HD11	1.70	0.73
1:R:162:ILE:HG22	1:R:163:ARG:HG3	1.71	0.72
2:A:187:THR:HB	2:A:198:MET:HB2	1.75	0.68
3:B:34:THR:HG21	3:B:300:LEU:HD12	1.76	0.68
1:R:130:PRO:HG2	2:A:340:THR:HG23	1.77	0.67
2:A:229:ASP:HA	2:A:232:LEU:HD13	1.79	0.65
1:R:163:ARG:HE	1:R:182:PRO:HD3	1.62	0.64
1:R:58:THR:H	1:R:61:THR:HB	1.64	0.63
1:R:135:ASN:HA	2:A:32:ARG:HG3	1.82	0.61
1:R:107:ILE:HD12	1:R:154:LEU:HD12	1.82	0.60
1:R:49:TRP:HD1	1:R:53:PHE:HD2	1.48	0.60
1:R:99:LYS:NZ	1:R:161:ILE:O	2.36	0.58
3:B:264:TYR:HE2	3:B:285:LEU:HD13	1.69	0.58
1:R:37:PHE:HB2	1:R:78:LEU:HD23	1.87	0.57
3:B:152:LEU:HD21	3:B:213:VAL:HG21	1.87	0.57
3:B:161:SER:OG	3:B:163:ASP:OD1	2.22	0.57
2:A:342:VAL:O	2:A:346:ASN:ND2	2.38	0.56
3:B:225:HIS:NE2	3:B:243:THR:OG1	2.39	0.56
3:B:279:SER:HA	3:B:320:VAL:HG11	1.88	0.56
1:R:199:THR:HA	1:R:265:THR:HG23	1.88	0.56
3:B:145:TYR:OH	3:B:188:MET:SD	2.57	0.55
2:A:306:GLN:NE2	2:A:306:GLN:O	2.40	0.54
3:B:47:THR:HA	3:B:339:TRP:CD1	2.42	0.54
2:A:315:ASP:OD1	2:A:316:THR:N	2.40	0.54
2:A:342:VAL:HG12	2:A:346:ASN:HD21	1.72	0.54
1:R:50:VAL:HA	1:R:54:ARG:HG2	1.89	0.53
1:R:114:PHE:HB3	1:R:149:PRO:HB3	1.90	0.53
1:R:257:TYR:OH	5:L:2:LEU:O	2.27	0.53
1:R:293:ASN:HA	1:R:296:LEU:HG	1.90	0.53
1:R:66:ASN:HD21	1:R:150:TRP:HE1	1.57	0.52
3:B:187:VAL:HA	3:B:203:ALA:HA	1.90	0.52
1:R:35:VAL:HG22	1:R:39:LEU:HD23	1.92	0.52
1:R:128:LEU:HD13	1:R:227:LYS:HD2	1.91	0.52
1:R:195:VAL:HG13	1:R:268:ILE:HG23	1.91	0.52
3:B:146:LEU:HD23	3:B:161:SER:HB3	1.92	0.51
3:B:200:VAL:HG12	3:B:210:LEU:HA	1.93	0.51
1:R:124:CYS:HB2	1:R:220:SER:HB3	1.93	0.51
3:B:198:LEU:HD12	3:B:210:LEU:HD21	1.93	0.51
1:R:239:PRO:HG3	2:A:354:PHE:HA	1.93	0.50
2:A:230:TYR:HA	2:A:243:MET:HB2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:308:LEU:HB3	3:B:339:TRP:HZ3	1.76	0.50
1:R:130:PRO:HG3	2:A:344:ILE:HD11	1.93	0.49
1:R:103:THR:O	1:R:107:ILE:HG12	2.13	0.49
3:B:168:LEU:HB3	3:B:178:THR:HB	1.94	0.49
3:B:266:HIS:ND1	3:B:268:ASN:OD1	2.46	0.48
3:B:251:ARG:HH11	3:B:251:ARG:HG3	1.77	0.48
3:B:259:GLN:HE22	4:C:30:VAL:HG21	1.76	0.48
2:A:185:VAL:HB	2:A:200:ASP:HB2	1.95	0.48
2:A:221:ILE:HB	2:A:264:ILE:HG22	1.96	0.48
3:B:95:LEU:HD13	3:B:100:VAL:HG11	1.96	0.48
3:B:145:TYR:O	3:B:162:GLY:N	2.47	0.47
3:B:279:SER:HB2	4:C:50:LEU:HD12	1.96	0.47
3:B:279:SER:OG	3:B:280:LYS:N	2.48	0.47
3:B:121:CYS:HB3	3:B:139:LEU:HB2	1.95	0.47
1:R:54:ARG:HD2	1:R:54:ARG:N	2.30	0.47
1:R:36:THR:HG23	1:R:295:CYS:HB3	1.97	0.47
1:R:156:LEU:HD11	1:R:208:ILE:HG21	1.97	0.47
1:R:25:LEU:HD11	1:R:285:VAL:HG11	1.97	0.46
3:B:79:LEU:HD22	3:B:93:ILE:HD12	1.95	0.46
3:B:47:THR:HG23	3:B:339:TRP:HE1	1.80	0.46
1:R:179:ASN:OD1	1:R:180:PHE:N	2.48	0.46
1:R:233:LEU:HD13	2:A:344:ILE:HG21	1.97	0.46
3:B:229:ILE:HD13	3:B:245:SER:HB3	1.96	0.46
1:R:128:LEU:HD21	1:R:223:LEU:HB3	1.97	0.46
3:B:100:VAL:HA	3:B:116:GLY:HA3	1.98	0.46
3:B:200:VAL:HG22	3:B:234:PHE:CE1	2.51	0.45
1:R:257:TYR:HA	1:R:286:THR:HB	1.98	0.45
3:B:289:TYR:HH	3:B:297:TRP:HE1	1.63	0.45
1:R:44:ASN:HD21	1:R:71:ASP:HB2	1.82	0.44
1:R:127:VAL:HG12	1:R:227:LYS:HD3	1.98	0.44
5:L:2:LEU:HD23	5:L:2:LEU:HA	1.87	0.44
3:B:180:PHE:HE1	3:B:216:GLY:HA2	1.83	0.44
3:B:146:LEU:HA	3:B:161:SER:HA	2.00	0.44
1:R:91:TRP:CE2	1:R:93:PHE:HB2	2.52	0.44
1:R:238:ARG:NH2	1:R:304:MET:O	2.49	0.44
3:B:118:ASP:O	3:B:120:ILE:HG13	2.17	0.43
5:L:2:LEU:HB3	5:L:3:PHE:H	1.58	0.43
1:R:48:ILE:HD13	1:R:65:LEU:HD12	1.99	0.43
1:R:133:THR:OG1	1:R:137:ARG:HD3	2.18	0.43
1:R:256:PRO:HA	1:R:259:VAL:HG12	1.99	0.43
3:B:266:HIS:HB3	3:B:269:ILE:HG12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:300:LEU:HA	1:R:304:MET:HG3	2.01	0.43
1:R:59:VAL:HG13	1:R:142:ALA:HB2	1.98	0.43
3:B:283:ARG:HB3	4:C:51:LEU:HD11	1.99	0.43
3:B:79:LEU:HD21	3:B:112:VAL:HG21	2.01	0.43
2:A:338:ALA:O	2:A:342:VAL:HG23	2.18	0.43
1:R:275:MET:HG2	1:R:276:TYR:CD1	2.55	0.42
2:A:308:GLU:O	2:A:311:ASN:ND2	2.52	0.42
3:B:34:THR:O	3:B:37:ILE:HG22	2.19	0.42
1:R:198:LEU:HD23	1:R:198:LEU:HA	1.87	0.42
1:R:260:VAL:HG21	1:R:286:THR:HG21	2.01	0.42
2:A:38:LEU:HB2	2:A:199:PHE:O	2.19	0.42
1:R:44:ASN:HB3	1:R:72:PHE:HB2	2.01	0.42
3:B:73:ALA:HB1	3:B:100:VAL:HG21	2.00	0.42
3:B:7:LEU:HG	4:C:16:VAL:HG21	2.02	0.42
3:B:308:LEU:HD23	3:B:339:TRP:HE3	1.85	0.41
1:R:157:THR:O	1:R:160:VAL:N	2.53	0.41
3:B:283:ARG:HD2	3:B:300:LEU:HD23	2.03	0.41
3:B:124:TYR:CE2	3:B:135:VAL:HG22	2.56	0.41
3:B:236:PRO:HB2	4:C:40:TYR:CE2	2.55	0.41
1:R:35:VAL:O	1:R:39:LEU:HG	2.21	0.41
3:B:66:ASP:OD1	3:B:66:ASP:N	2.52	0.41
3:B:284:LEU:HD23	3:B:284:LEU:H	1.86	0.41
3:B:337:LYS:HB3	3:B:339:TRP:NE1	2.36	0.41
2:A:266:LEU:HD12	2:A:266:LEU:HA	1.92	0.40
2:A:35:LYS:HB3	2:A:199:PHE:HE2	1.85	0.40
2:A:307:PHE:HA	2:A:310:LEU:HD12	2.03	0.40
3:B:51:LEU:HD22	3:B:336:LEU:HD22	2.02	0.40
1:R:91:TRP:N	1:R:174:VAL:O	2.55	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	R	296/371 (80%)	288 (97%)	8 (3%)	0	100	100
2	A	212/354 (60%)	211 (100%)	1 (0%)	0	100	100
3	B	334/351 (95%)	324 (97%)	10 (3%)	0	100	100
4	C	52/71 (73%)	50 (96%)	2 (4%)	0	100	100
5	L	1/3 (33%)	0	1 (100%)	0	100	100
All	All	895/1150 (78%)	873 (98%)	22 (2%)	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	R	237/311 (76%)	237 (100%)	0	100	100
2	A	182/306 (60%)	181 (100%)	1 (0%)	88	93
3	B	246/293 (84%)	246 (100%)	0	100	100
4	C	35/58 (60%)	35 (100%)	0	100	100
5	L	2/2 (100%)	2 (100%)	0	100	100
All	All	702/970 (72%)	701 (100%)	1 (0%)	93	97

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	46	LYS

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

Mol	Chain	Res	Type
1	R	44	ASN
2	A	204	GLN
2	A	346	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)

1 non-standard protein/DNA/RNA residue 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	$\#\|Z\| > 2$
5	FME	L	1	5	8,9,10	0.92	0	7,9,11	0.89	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	FME	L	1	5	-	5/7/9/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	L	1	FME	O1-CN-N-CA
5	L	1	FME	CB-CA-N-CN
5	L	1	FME	C-CA-CB-CG
5	L	1	FME	N-CA-CB-CG
5	L	1	FME	CA-CB-CG-SD

There are no ring outliers.

No monomer is involved in short contacts.

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.

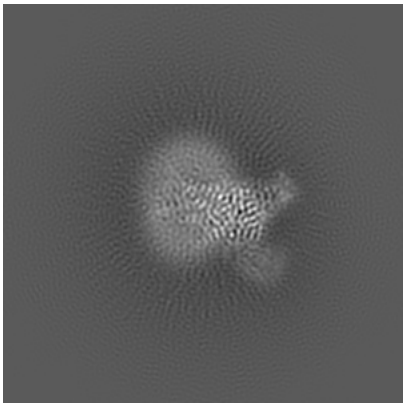
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-32858. These allow visual inspection of the internal detail of the map and identification of artifacts.

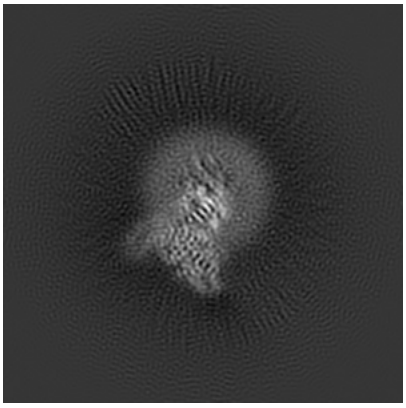
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections (i)

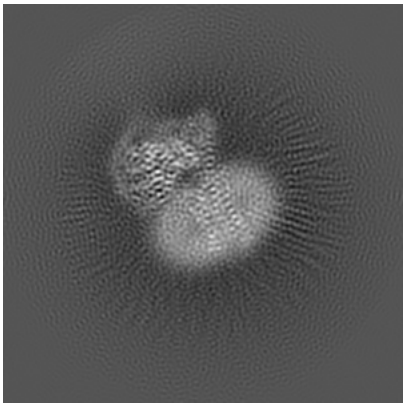
6.1.1 Primary map



X



Y

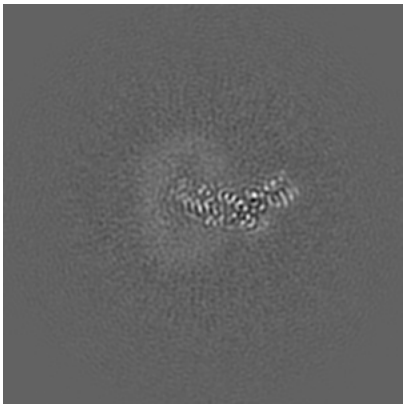


Z

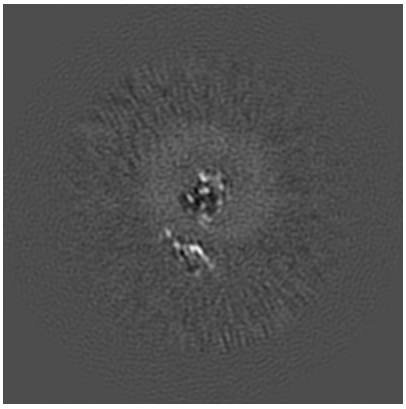
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

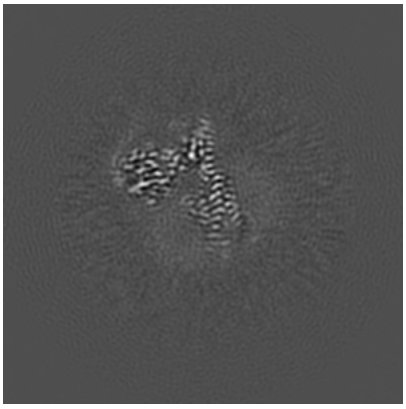
6.2.1 Primary map



X Index: 128



Y Index: 128

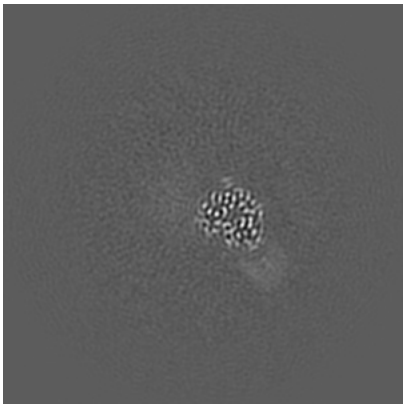


Z Index: 128

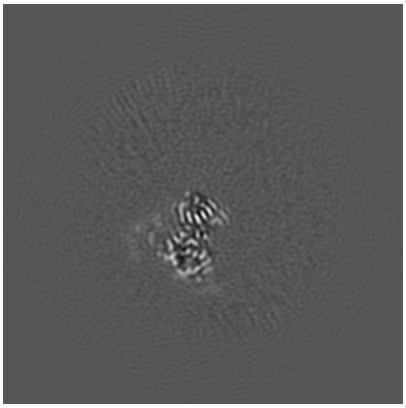
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

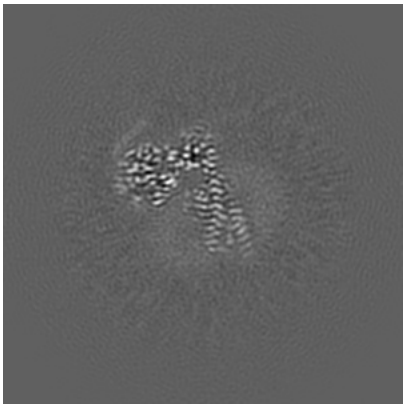
6.3.1 Primary map



X Index: 96



Y Index: 159



Z Index: 125

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 0.0298. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

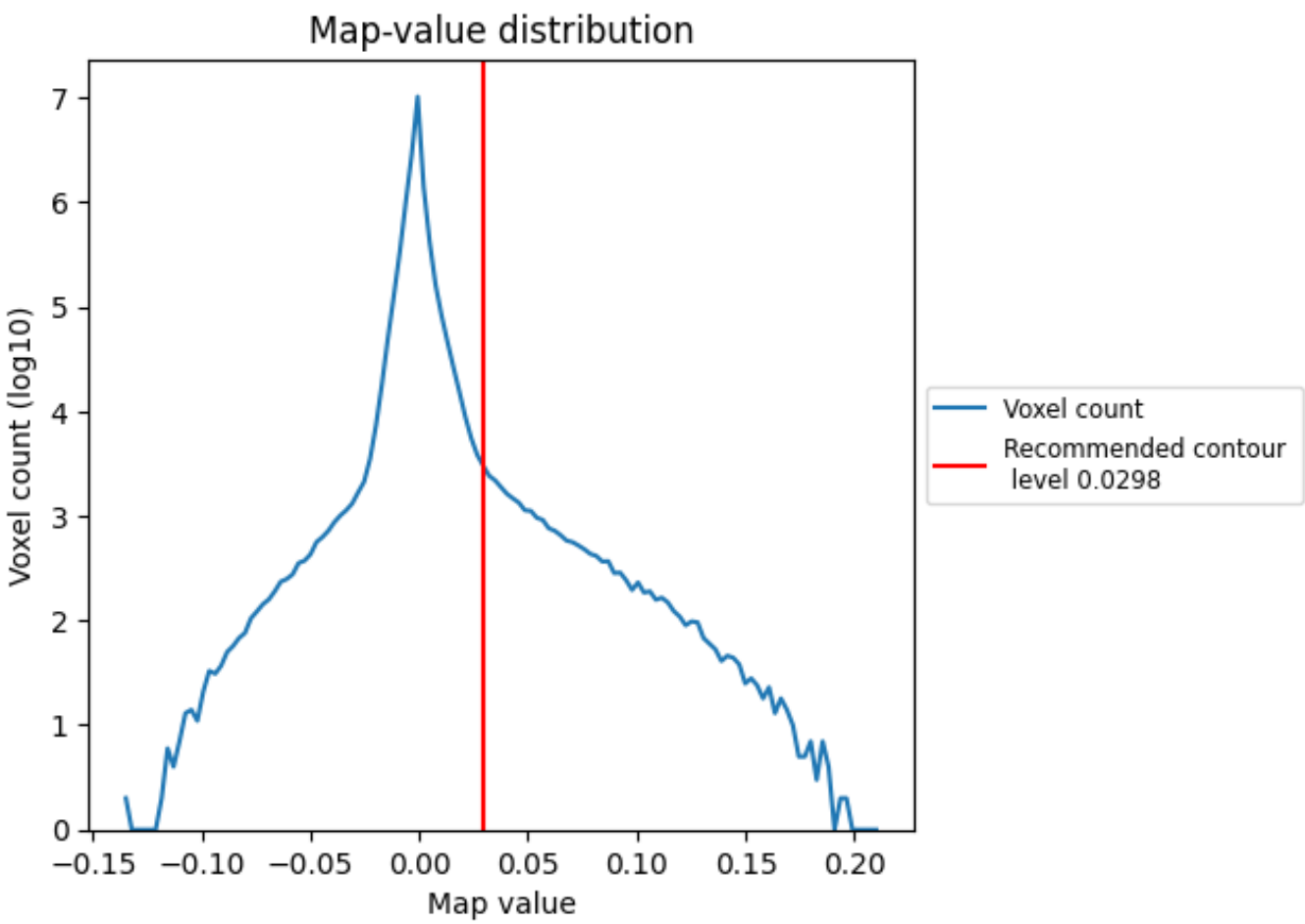
6.5 Mask visualisation (i)

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

7 Map analysis (i)

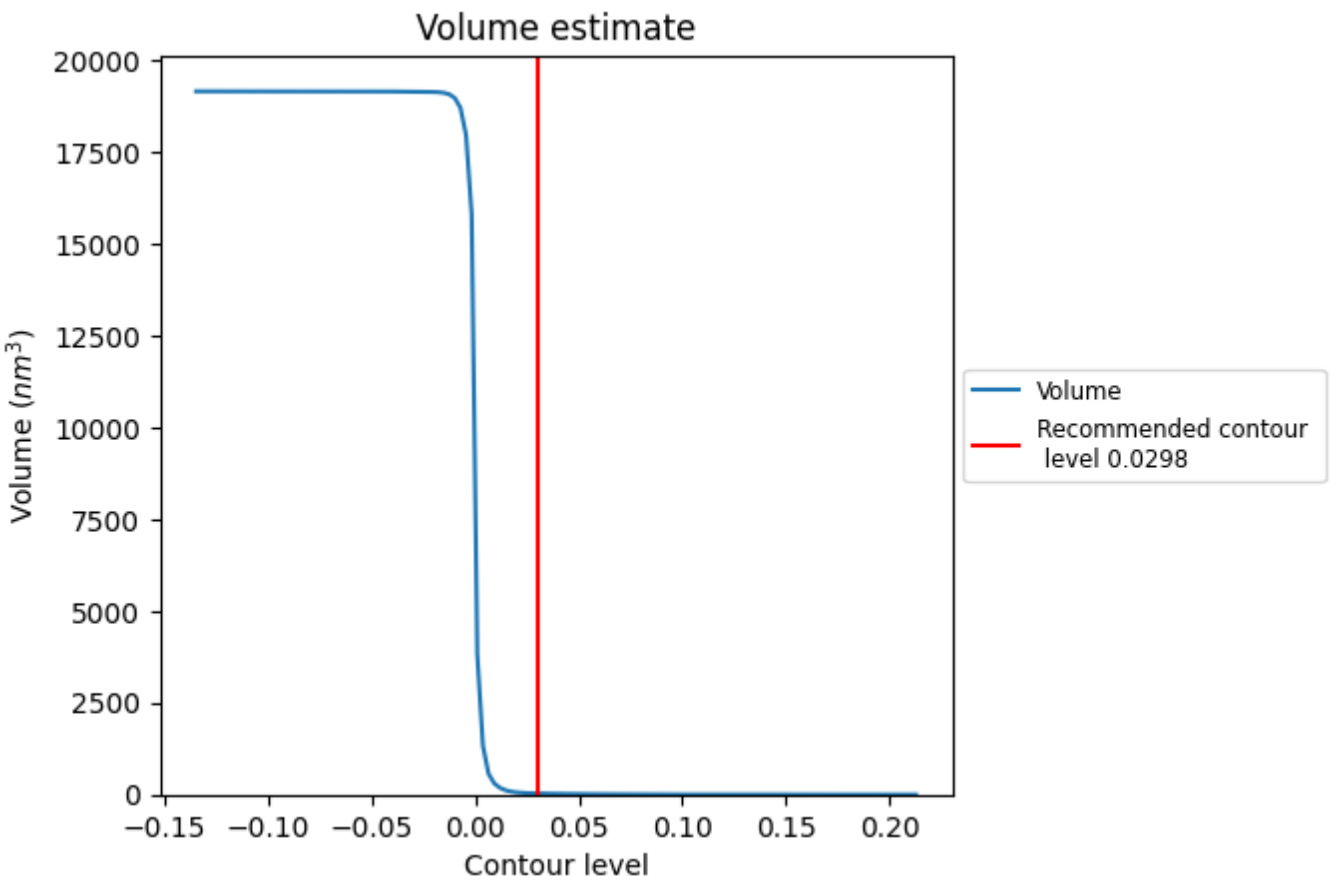
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution (i)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

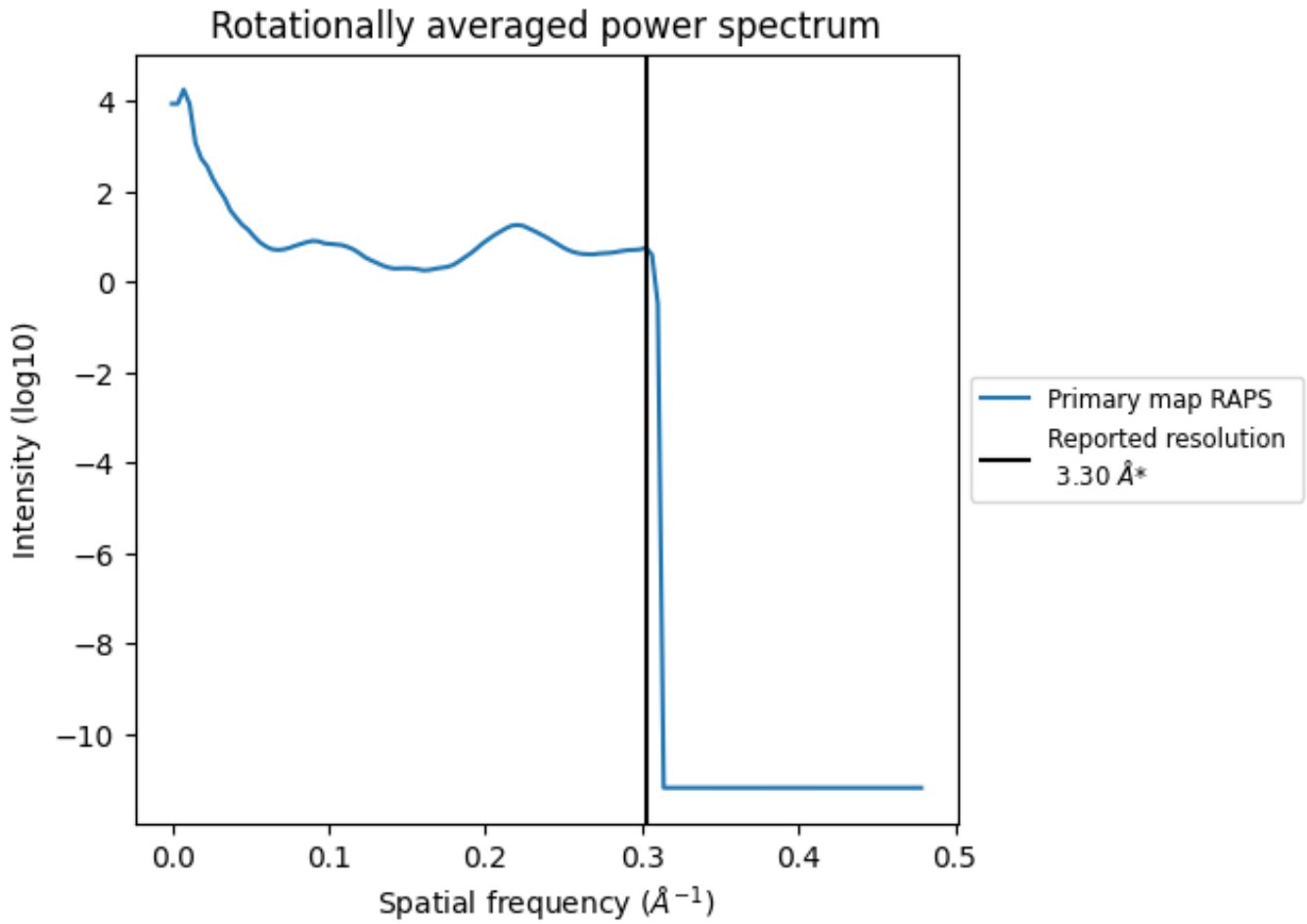
7.2 Volume estimate (i)



The volume at the recommended contour level is 31 nm^3; this corresponds to an approximate mass of 28 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 0.303 Å^{-1}

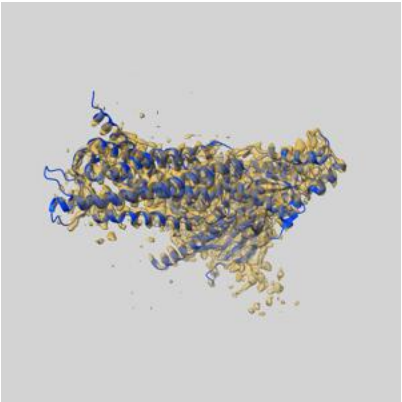
8 Fourier-Shell correlation (i)

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

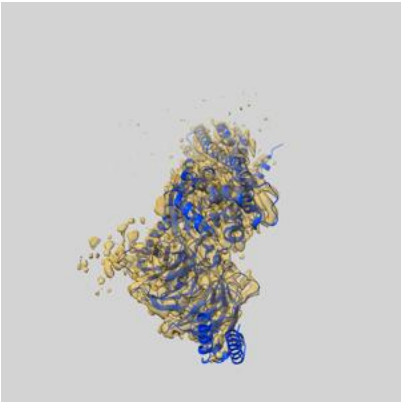
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-32858 and PDB model 7WVU. Per-residue inclusion information can be found in section 3 on page 6.

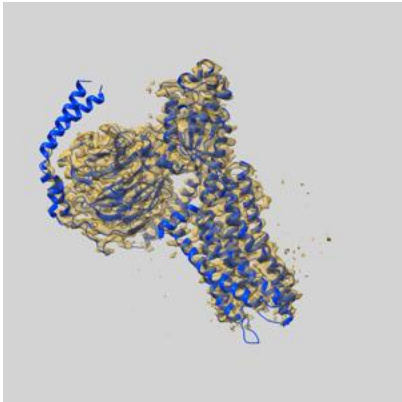
9.1 Map-model overlay (i)



X



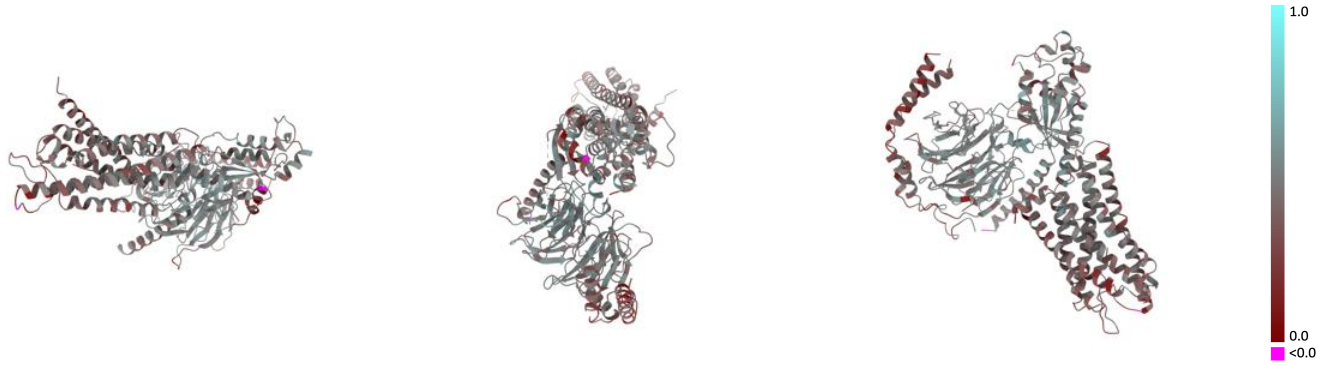
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0298 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

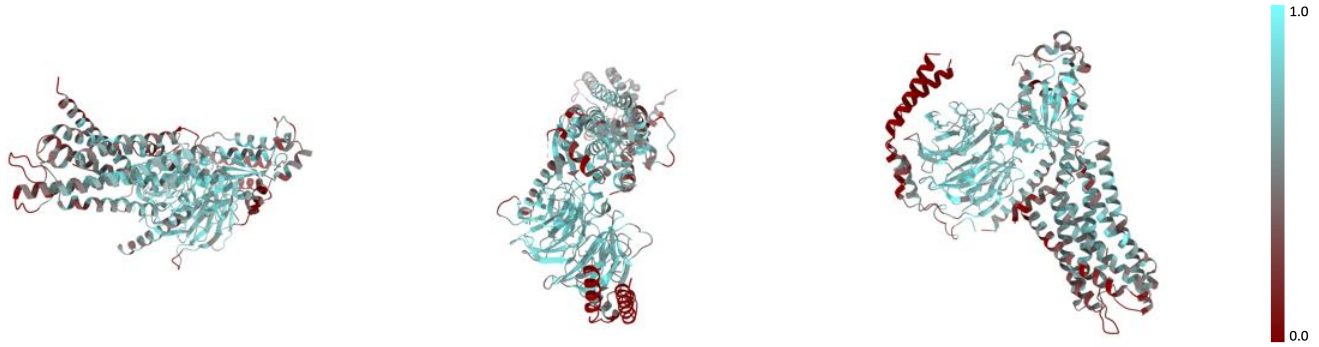
9.2 Q-score mapped to coordinate model (i)



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.

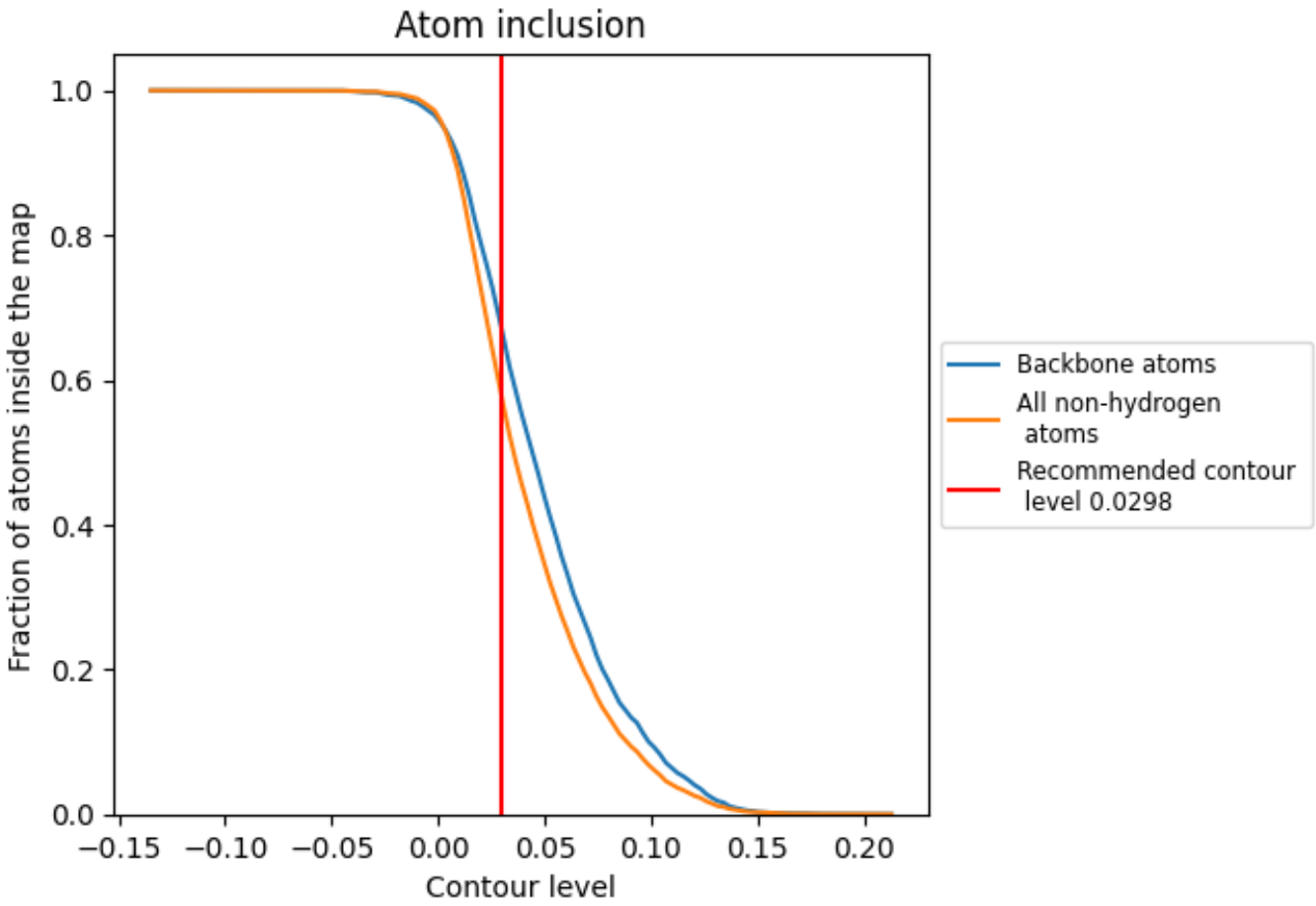
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.0298).

9.4 Atom inclusion (i)



At the recommended contour level, 67% of all backbone atoms, 58% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.0298) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5763	0.4340
A	0.6128	0.4570
B	0.6581	0.4560
C	0.3298	0.3840
L	0.2333	0.4130
R	0.5069	0.4000