



Full wwPDB EM Validation Report ⓘ

Mar 26, 2026 – 02:09 AM UTC

PDB ID : 7MPE / pdb_00007mpe
EMDB ID : EMD-23932
Title : Cryo-EM structure of the yeast cadmium factor 1 protein (Ycf1p)
Authors : Bickers, S.C.; Benlekbir, S.; Rubinstein, J.L.; Kanelis, V.
Deposited on : 2021-05-04
Resolution : 3.20 Å(reported)

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<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.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

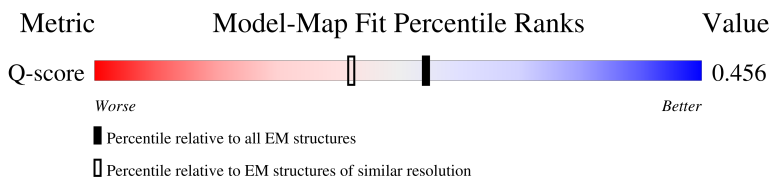
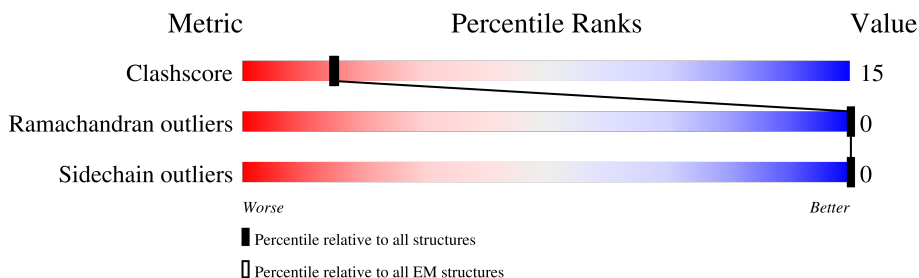
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.20 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	15020 (2.70 - 3.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1537	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 11214 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 Metal resistance protein YCF1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1428	Total	C	N	O	S	0	0
			11214	7251	1873	2039	51		

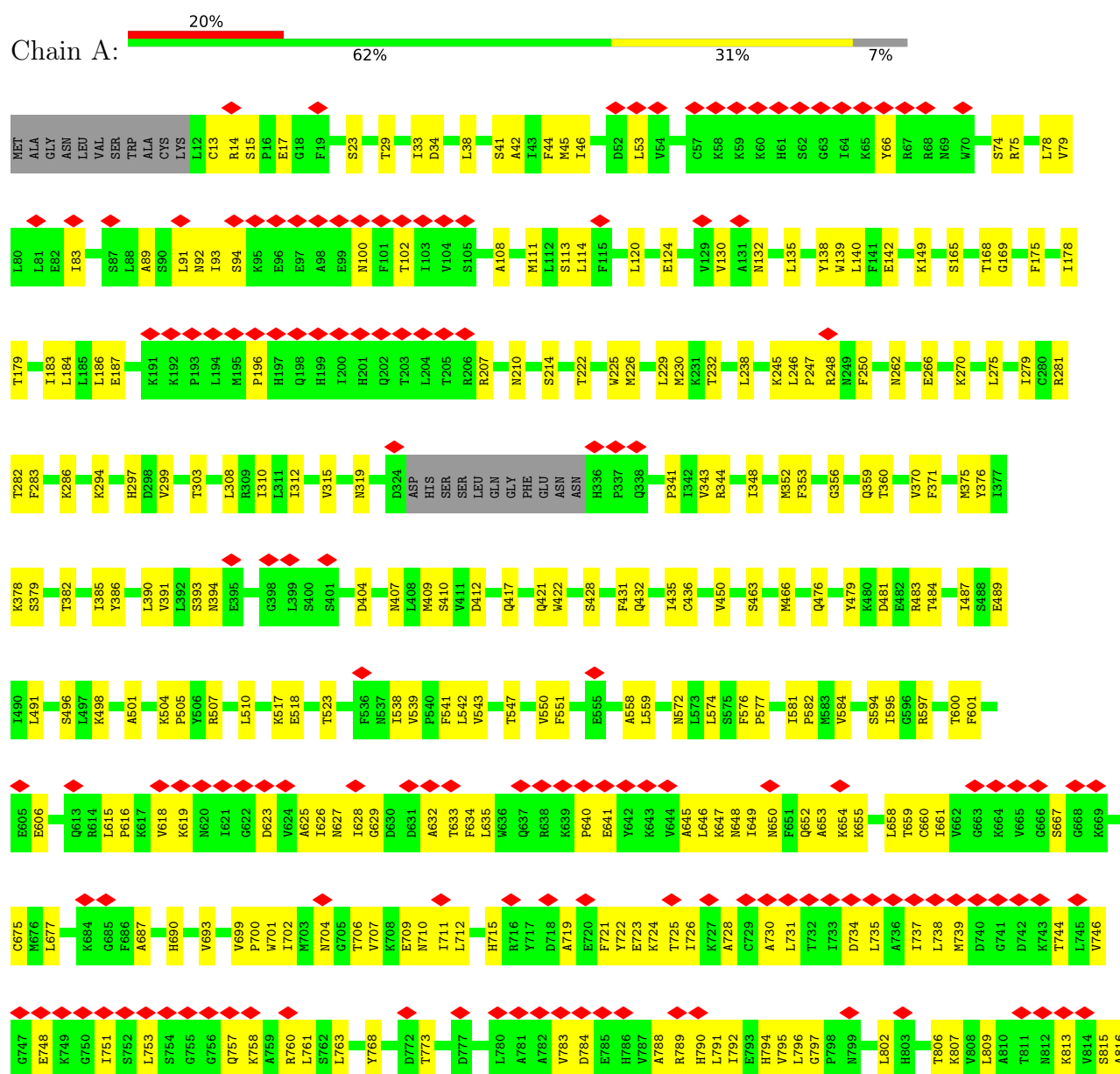
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1516	ASP	-	expression tag	UNP P39109
A	1517	TYR	-	expression tag	UNP P39109
A	1518	LYS	-	expression tag	UNP P39109
A	1519	ASP	-	expression tag	UNP P39109
A	1520	HIS	-	expression tag	UNP P39109
A	1521	ASP	-	expression tag	UNP P39109
A	1522	GLY	-	expression tag	UNP P39109
A	1523	ASP	-	expression tag	UNP P39109
A	1524	TYR	-	expression tag	UNP P39109
A	1525	LYS	-	expression tag	UNP P39109
A	1526	ASP	-	expression tag	UNP P39109
A	1527	HIS	-	expression tag	UNP P39109
A	1528	ASP	-	expression tag	UNP P39109
A	1529	ILE	-	expression tag	UNP P39109
A	1530	ASP	-	expression tag	UNP P39109
A	1531	TYR	-	expression tag	UNP P39109
A	1532	LYS	-	expression tag	UNP P39109
A	1533	ASP	-	expression tag	UNP P39109
A	1534	ASP	-	expression tag	UNP P39109
A	1535	ASP	-	expression tag	UNP P39109
A	1536	ASP	-	expression tag	UNP P39109
A	1537	LYS	-	expression tag	UNP P39109

3 Residue-property plots

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

• Molecule 1: Metal resistance protein YCF1



E1487	C1419	I1351	T1279	G1184	F1068	L940	ILE	E1488	L1420	P1352	R1280	L1189	V1084	Y942	S818	E1489	A1421	Q1354	E1284	R1198	T1088	K944	GLY	E1490	L1425	S1356	D1286	L1199	T1089	C946	S822	E1491	R1422	D1355	L1285	L1204	M1208	V1209	S1100	Y951	L1492	V1426	P1427	V1357	L1287	N1293	L1213	Y1103	F954	ASN	G829	N828	D827	L826	G1493	V1428	S1428	K1429	E1360	H1291	I1290	K1290	M1208	F1359	V1358	W1358	L1288	V1288	L1287	D1286	E1284	R1280	L1189	V1084	Y942	S818	E1487	L1501	F1502	Y1503	S1504	L1505	C1506	D1441	V1442	E1443	T1444	D1445	K1446	V1447	V1448	L1511	V1512	ASN	T1451	I1452	R1453	T1454	A1455	F1456	K1457	D1458	R1459	T1460	T1461	I1461	M1394	S1395	N1396	D1397	G1398	L1399	D1400	A1401	Q1402	L1403	T1404	E1405	G1406	G1407	G1408	N1409	D1410	S1411	V1412	R1415	A1486	N1482	D1481	L1480	V1479	D1475	S1474	D1473	M1472	I1471	N1469	L1468	R1467	H1466	A1465	ASP	ILE	ASP	TYR	LYS	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP	ASP
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	124864	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	45	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	4.508	Depositor
Minimum map value	-2.776	Depositor
Average map value	0.009	Depositor
Map value standard deviation	0.113	Depositor
Recommended contour level	0.879	Depositor
Map size (Å)	263.68, 263.68, 263.68	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.03, 1.03, 1.03	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	A	0.20	0/11443	0.39	0/15530

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	11214	0	11296	336	0
All	All	11214	0	11296	336	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 15.

All (336) 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:A:629:GLY:O	1:A:650:ASN:HA	1.47	1.11
1:A:391:VAL:HG12	1:A:606:GLU:HB2	1.60	0.84
1:A:78:LEU:HB2	1:A:183:ILE:HD12	1.65	0.79
1:A:1084:VAL:HG11	1:A:1215:TYR:HD2	1.51	0.75
1:A:1367:ILE:HG22	1:A:1425:LEU:HD22	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:LEU:HD13	1:A:120:LEU:HD22	1.66	0.75
1:A:1255:LEU:HD12	1:A:1343:TYR:HB2	1.66	0.75
1:A:632:ALA:HB3	1:A:649:ILE:HB	1.72	0.72
1:A:210:ASN:HD21	1:A:247:PRO:HG3	1.55	0.71
1:A:378:LYS:HB2	1:A:417:GLN:HG3	1.71	0.71
1:A:646:LEU:HD11	1:A:649:ILE:HD11	1.73	0.71
1:A:1379:TRP:HE1	1:A:1399:LEU:HD21	1.56	0.70
1:A:655:LYS:HA	1:A:806:THR:HB	1.73	0.70
1:A:83:ILE:HD11	1:A:111:MET:HB2	1.74	0.68
1:A:792:ILE:HG21	1:A:813:LYS:HZ3	1.59	0.68
1:A:640:PRO:O	1:A:641:GLU:HG2	1.92	0.67
1:A:802:LEU:O	1:A:807:LYS:NZ	2.26	0.67
1:A:635:LEU:HA	1:A:645:ALA:HB3	1.77	0.67
1:A:498:LYS:NZ	1:A:1252:GLU:OE1	2.28	0.66
1:A:1020:HIS:O	1:A:1020:HIS:ND1	2.29	0.66
1:A:409:MET:HE2	1:A:1132:PHE:CD2	2.30	0.66
1:A:1275:ASN:ND2	1:A:1329:ASN:OD1	2.27	0.66
1:A:1331:VAL:HA	1:A:1336:ALA:HA	1.78	0.66
1:A:1356:SER:HG	1:A:1419:CYS:HG	1.36	0.66
1:A:626:ILE:O	1:A:690:HIS:ND1	2.28	0.65
1:A:1479:VAL:HG23	1:A:1486:ALA:HB3	1.77	0.65
1:A:1351:ILE:HG13	1:A:1432:VAL:HG23	1.78	0.65
1:A:1052:ARG:HG3	1:A:1246:TYR:HE1	1.60	0.65
1:A:375:MET:HE1	1:A:1167:TYR:HD1	1.61	0.64
1:A:207:ARG:O	1:A:245:LYS:NZ	2.31	0.64
1:A:1363:VAL:HG23	1:A:1403:LEU:HD11	1.80	0.64
1:A:1420:LEU:HD22	1:A:1448:VAL:HG22	1.80	0.64
1:A:34:ASP:HA	1:A:38:LEU:HB2	1.80	0.63
1:A:294:LYS:NZ	1:A:428:SER:OG	2.31	0.63
1:A:661:ILE:HG22	1:A:824:ALA:HB3	1.81	0.63
1:A:139:TRP:HE1	1:A:183:ILE:HG23	1.64	0.62
1:A:1116:GLU:OE1	1:A:1116:GLU:N	2.33	0.62
1:A:795:VAL:HG23	1:A:796:LEU:HG	1.80	0.62
1:A:1307:ARG:HB3	1:A:1505:LEU:HD21	1.81	0.62
1:A:1504:SER:HA	1:A:1507:MET:HE3	1.82	0.61
1:A:379:SER:HB3	1:A:1163:MET:HE3	1.83	0.61
1:A:1278:SER:HB3	1:A:1291:HIS:H	1.66	0.61
1:A:677:LEU:HD12	1:A:1144:GLY:HA2	1.82	0.61
1:A:1420:LEU:HD21	1:A:1451:THR:HG21	1.82	0.61
1:A:83:ILE:HB	1:A:114:LEU:HD22	1.83	0.60
1:A:412:ASP:O	1:A:594:SER:OG	2.18	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1280:ARG:HG2	1:A:1287:LEU:HA	1.82	0.60
1:A:476:GLN:NE2	1:A:518:GLU:OE2	2.35	0.60
1:A:1167:TYR:HB3	1:A:1168:PRO:HD3	1.83	0.59
1:A:738:LEU:HD12	1:A:744:THR:HG21	1.84	0.59
1:A:1084:VAL:HG11	1:A:1215:TYR:CD2	2.37	0.59
1:A:760:ARG:HH11	1:A:783:VAL:HG23	1.68	0.59
1:A:1052:ARG:HG3	1:A:1246:TYR:CE1	2.38	0.59
1:A:1266:TRP:HB2	1:A:1267:PRO:HD3	1.85	0.58
1:A:491:LEU:HD13	1:A:1049:ILE:HG23	1.84	0.58
1:A:1274:PHE:HB2	1:A:1294:ILE:H	1.69	0.58
1:A:706:THR:HG23	1:A:709:GLU:H	1.69	0.58
1:A:315:VAL:HG22	1:A:1199:LEU:HD22	1.86	0.57
1:A:356:GLY:O	1:A:360:THR:HG23	2.04	0.57
1:A:1088:THR:HG22	1:A:1089:THR:HG23	1.87	0.57
1:A:108:ALA:HA	1:A:111:MET:SD	2.45	0.57
1:A:262:ASN:ND2	1:A:282:THR:OG1	2.36	0.57
1:A:632:ALA:HB1	1:A:634:PHE:CE1	2.40	0.57
1:A:422:TRP:HB3	1:A:584:VAL:HG22	1.87	0.57
1:A:650:ASN:H	1:A:831:ILE:HD12	1.69	0.57
1:A:1256:ILE:HG23	1:A:1260:HIS:CE1	2.40	0.57
1:A:701:TRP:NE1	1:A:1141:THR:OG1	2.38	0.57
1:A:483:ARG:HG3	1:A:510:LEU:HD11	1.87	0.57
1:A:1479:VAL:HG11	1:A:1502:PHE:HB2	1.87	0.57
1:A:937:ASN:O	1:A:941:GLU:HG2	2.05	0.56
1:A:130:VAL:HG22	1:A:214:SER:HA	1.86	0.56
1:A:836:THR:OG1	1:A:838:ASP:OD1	2.17	0.56
1:A:385:ILE:HG23	1:A:601:PHE:HE2	1.71	0.56
1:A:404:ASP:OD1	1:A:597:ARG:NH2	2.39	0.56
1:A:42:ALA:O	1:A:46:ILE:HG12	2.05	0.56
1:A:723:GLU:HA	1:A:726:ILE:HG22	1.88	0.56
1:A:1321:ARG:HH21	1:A:1338:ASN:HA	1.71	0.56
1:A:17:GLU:OE1	1:A:29:THR:HG21	2.06	0.55
1:A:700:PRO:O	1:A:758:LYS:NZ	2.36	0.55
1:A:693:VAL:HG12	1:A:773:THR:HB	1.88	0.55
1:A:1266:TRP:CD1	1:A:1335:ILE:HG21	2.41	0.55
1:A:826:LEU:HD22	1:A:831:ILE:HG12	1.88	0.55
1:A:1198:ARG:NH1	1:A:1208:MET:SD	2.79	0.55
1:A:248:ARG:HH11	1:A:248:ARG:HG2	1.72	0.55
1:A:626:ILE:HG21	1:A:806:THR:HG21	1.87	0.55
1:A:626:ILE:HG13	1:A:627:ASN:N	2.23	0.54
1:A:947:ASN:OD1	1:A:949:LYS:NZ	2.39	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ARG:HH11	1:A:187:GLU:HA	1.73	0.54
1:A:633:THR:HG22	1:A:648:ASN:H	1.72	0.54
1:A:407:ASN:HA	1:A:410:SER:OG	2.08	0.54
1:A:813:LYS:HZ1	1:A:816:ALA:HB2	1.73	0.54
1:A:975:TRP:NE1	1:A:979:ASN:HD21	2.06	0.54
1:A:1006:THR:O	1:A:1009:GLN:HG3	2.07	0.54
1:A:1348:LYS:NZ	1:A:1427:PRO:O	2.32	0.54
1:A:507:ARG:HD3	1:A:1034:LEU:HD12	1.89	0.54
1:A:572:ASN:ND2	1:A:1218:GLN:OE1	2.41	0.54
1:A:1226:ILE:O	1:A:1230:THR:OG1	2.26	0.53
1:A:1272:ILE:HD12	1:A:1274:PHE:HE1	1.71	0.53
1:A:91:LEU:O	1:A:94:SER:OG	2.26	0.53
1:A:421:GLN:N	1:A:421:GLN:OE1	2.38	0.53
1:A:1143:ARG:HG3	1:A:1148:GLN:HE21	1.72	0.53
1:A:1260:HIS:ND1	1:A:1344:ASP:OD2	2.38	0.53
1:A:222:THR:O	1:A:1179:ARG:NH1	2.42	0.53
1:A:739:MET:N	1:A:739:MET:SD	2.75	0.53
1:A:1456:PHE:HB3	1:A:1459:ARG:HB2	1.89	0.53
1:A:266:GLU:OE2	1:A:270:LYS:HD3	2.08	0.53
1:A:315:VAL:O	1:A:319:ASN:ND2	2.42	0.53
1:A:619:LYS:HD2	1:A:655:LYS:HD2	1.90	0.53
1:A:757:GLN:HG3	1:A:760:ARG:NH2	2.24	0.53
1:A:344:ARG:O	1:A:348:ILE:HG13	2.08	0.52
1:A:702:ILE:HG21	1:A:746:VAL:HG11	1.90	0.52
1:A:551:PHE:CD2	1:A:559:LEU:HD12	2.44	0.52
1:A:707:VAL:O	1:A:711:ILE:HG12	2.09	0.52
1:A:1390:HIS:CE1	1:A:1394:MET:HE1	2.45	0.52
1:A:667:SER:OG	1:A:828:ASN:N	2.41	0.52
1:A:626:ILE:HD13	1:A:806:THR:HG21	1.91	0.52
1:A:746:VAL:HG13	1:A:753:LEU:HD23	1.90	0.52
1:A:308:LEU:O	1:A:312:ILE:HG13	2.08	0.52
1:A:702:ILE:HD11	1:A:758:LYS:HG3	1.91	0.52
1:A:139:TRP:HZ2	1:A:183:ILE:HD13	1.75	0.51
1:A:1046:ILE:HA	1:A:1049:ILE:HG22	1.91	0.51
1:A:1277:TYR:HE2	1:A:1318:ALA:HB1	1.75	0.51
1:A:731:LEU:O	1:A:735:LEU:N	2.39	0.51
1:A:139:TRP:NE1	1:A:183:ILE:HG23	2.25	0.51
1:A:483:ARG:NH2	1:A:1053:PHE:O	2.43	0.51
1:A:74:SER:HB3	1:A:186:LEU:HD21	1.93	0.51
1:A:14:ARG:NH2	1:A:100:ASN:OD1	2.44	0.51
1:A:1288:VAL:HG13	1:A:1289:LEU:HG	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:SER:OG	1:A:169:GLY:HA3	2.10	0.51
1:A:45:MET:HG3	1:A:138:TYR:CD1	2.46	0.51
1:A:210:ASN:ND2	1:A:247:PRO:HG3	2.23	0.51
1:A:827:ASP:O	1:A:830:GLU:N	2.44	0.50
1:A:1494:LEU:O	1:A:1500:SER:OG	2.24	0.50
1:A:229:LEU:O	1:A:232:THR:OG1	2.29	0.50
1:A:954:PHE:CZ	1:A:958:ILE:HD11	2.46	0.50
1:A:1419:CYS:SG	1:A:1422:ARG:NH1	2.85	0.50
1:A:124:GLU:OE2	1:A:132:ASN:N	2.35	0.49
1:A:538:ILE:O	1:A:542:LEU:N	2.44	0.49
1:A:175:PHE:HA	1:A:178:ILE:HG22	1.94	0.49
1:A:498:LYS:HB3	1:A:1346:ARG:HH12	1.77	0.49
1:A:310:ILE:HG22	1:A:348:ILE:HD13	1.93	0.49
1:A:352:MET:HE1	1:A:1213:LEU:HD21	1.93	0.49
1:A:724:LYS:NZ	1:A:728:ALA:HB2	2.27	0.49
1:A:1273:LYS:HG2	1:A:1295:HIS:CD2	2.47	0.49
1:A:83:ILE:HD13	1:A:114:LEU:HD22	1.95	0.49
1:A:813:LYS:NZ	1:A:815:SER:OG	2.29	0.49
1:A:297:HIS:CD2	1:A:359:GLN:HG3	2.48	0.49
1:A:763:LEU:HD23	1:A:791:LEU:HD11	1.94	0.49
1:A:627:ASN:OD1	1:A:653:ALA:HB3	2.12	0.49
1:A:784:ASP:O	1:A:788:ALA:N	2.42	0.49
1:A:721:PHE:O	1:A:725:THR:HG23	2.12	0.49
1:A:939:TYR:OH	1:A:1234:GLU:OE2	2.31	0.49
1:A:1299:ASN:O	1:A:1301:LYS:NZ	2.39	0.49
1:A:262:ASN:HD21	1:A:281:ARG:HB3	1.77	0.48
1:A:431:PHE:O	1:A:435:ILE:HG12	2.13	0.48
1:A:719:ALA:O	1:A:722:TYR:N	2.45	0.48
1:A:788:ALA:O	1:A:792:ILE:HG23	2.13	0.48
1:A:825:LEU:HD13	1:A:847:PRO:HB2	1.94	0.48
1:A:841:THR:HG23	1:A:849:TRP:HA	1.96	0.48
1:A:711:ILE:HG22	1:A:768:TYR:CD2	2.48	0.48
1:A:1262:PRO:CD	1:A:1263:PRO:HD3	2.44	0.48
1:A:275:LEU:O	1:A:279:ILE:HG13	2.14	0.48
1:A:436:CYS:SG	1:A:574:LEU:HB2	2.54	0.48
1:A:484:THR:HA	1:A:487:ILE:HG22	1.95	0.48
1:A:539:VAL:O	1:A:543:VAL:HG23	2.14	0.48
1:A:1262:PRO:HG2	1:A:1263:PRO:HD3	1.94	0.48
1:A:1039:THR:O	1:A:1043:THR:HG23	2.14	0.48
1:A:1056:ASP:OD1	1:A:1242:ARG:NH1	2.47	0.48
1:A:1377:ALA:HA	1:A:1380:ARG:HG2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1431:LEU:O	1:A:1461:ILE:HG13	2.13	0.48
1:A:1026:HIS:HB2	1:A:1057:ILE:HD12	1.95	0.47
1:A:1364:ARG:NH2	1:A:1371:ASN:OD1	2.47	0.47
1:A:1500:SER:O	1:A:1504:SER:N	2.45	0.47
1:A:623:ASP:HA	1:A:655:LYS:HB3	1.97	0.47
1:A:226:MET:HE2	1:A:230:MET:HE2	1.96	0.47
1:A:1176:LEU:HD21	1:A:1227:VAL:HG11	1.96	0.47
1:A:1441:ASP:O	1:A:1444:THR:OG1	2.31	0.47
1:A:634:PHE:CD2	1:A:675:CYS:HB3	2.48	0.47
1:A:1229:MET:O	1:A:1233:VAL:HG23	2.13	0.47
1:A:13:CYS:SG	1:A:102:THR:HG23	2.55	0.47
1:A:450:VAL:HG21	1:A:550:VAL:HG21	1.96	0.47
1:A:647:LYS:NZ	1:A:830:GLU:OE2	2.35	0.47
1:A:479:TYR:CE1	1:A:517:LYS:HG2	2.50	0.47
1:A:792:ILE:O	1:A:797:GLY:N	2.47	0.47
1:A:1068:PHE:HE1	1:A:1233:VAL:HG13	1.79	0.47
1:A:1040:PHE:O	1:A:1044:THR:HG22	2.14	0.47
1:A:1270:GLY:O	1:A:1460:THR:OG1	2.25	0.47
1:A:283:PHE:CE1	1:A:376:TYR:HB3	2.50	0.46
1:A:1153:HIS:NE2	1:A:1157:CYS:SG	2.89	0.46
1:A:1339:GLU:N	1:A:1339:GLU:OE1	2.49	0.46
1:A:179:THR:O	1:A:183:ILE:HG22	2.15	0.46
1:A:504:LYS:HG2	1:A:507:ARG:HH22	1.80	0.46
1:A:826:LEU:HD22	1:A:831:ILE:HA	1.97	0.46
1:A:1304:ILE:HG22	1:A:1312:LYS:HG3	1.98	0.46
1:A:1448:VAL:O	1:A:1452:ILE:HG12	2.14	0.46
1:A:1064:LEU:HD12	1:A:1236:ASN:HB3	1.98	0.46
1:A:1289:LEU:HD13	1:A:1292:ILE:HD12	1.96	0.46
1:A:504:LYS:HB2	1:A:505:PRO:HD3	1.96	0.46
1:A:712:LEU:HD12	1:A:715:HIS:O	2.16	0.46
1:A:944:LYS:HA	1:A:948:PRO:HG3	1.97	0.46
1:A:1263:PRO:C	1:A:1265:GLU:H	2.24	0.46
1:A:790:HIS:O	1:A:794:HIS:ND1	2.32	0.46
1:A:238:LEU:HD23	1:A:238:LEU:HA	1.85	0.46
1:A:626:ILE:N	1:A:653:ALA:O	2.42	0.46
1:A:1278:SER:O	1:A:1326:SER:OG	2.28	0.45
1:A:1280:ARG:HH11	1:A:1287:LEU:HD11	1.80	0.45
1:A:845:ASP:O	1:A:849:TRP:N	2.47	0.45
1:A:1199:LEU:HB2	1:A:1204:LEU:HD23	1.99	0.45
1:A:1356:SER:OG	1:A:1419:CYS:SG	2.52	0.45
1:A:813:LYS:HD2	1:A:815:SER:H	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1110:TYR:CD2	1:A:1176:LEU:HD22	2.51	0.45
1:A:341:PRO:HB2	1:A:343:VAL:HG12	1.98	0.45
1:A:634:PHE:N	1:A:646:LEU:O	2.50	0.45
1:A:1305:VAL:O	1:A:1479:VAL:HA	2.17	0.45
1:A:730:ALA:HB2	1:A:790:HIS:CE1	2.52	0.45
1:A:1277:TYR:CE2	1:A:1318:ALA:HB1	2.52	0.45
1:A:629:GLY:O	1:A:650:ASN:CA	2.41	0.45
1:A:44:PHE:CD2	1:A:138:TYR:HB2	2.52	0.44
1:A:135:LEU:HG	1:A:135:LEU:O	2.17	0.44
1:A:701:TRP:CD2	1:A:1138:GLY:HA3	2.52	0.44
1:A:1447:VAL:O	1:A:1451:THR:HG22	2.17	0.44
1:A:660:CYS:SG	1:A:809:LEU:HD23	2.57	0.44
1:A:1452:ILE:O	1:A:1456:PHE:HB2	2.17	0.44
1:A:1503:TYR:CD2	1:A:1507:MET:HE2	2.52	0.44
1:A:238:LEU:O	1:A:1120:LEU:HD11	2.16	0.44
1:A:523:THR:HG22	1:A:1019:ILE:HD13	1.99	0.44
1:A:633:THR:HA	1:A:647:LYS:HA	1.99	0.44
1:A:375:MET:HE1	1:A:1167:TYR:CD1	2.48	0.44
1:A:1321:ARG:NE	1:A:1337:ILE:O	2.49	0.44
1:A:1379:TRP:NE1	1:A:1399:LEU:HD21	2.28	0.44
1:A:576:PHE:HB3	1:A:577:PRO:HD3	1.99	0.44
1:A:541:PHE:CZ	1:A:999:GLY:HA3	2.52	0.44
1:A:813:LYS:O	1:A:817:LEU:HD23	2.17	0.44
1:A:1103:TYR:OH	1:A:1226:ILE:HG23	2.18	0.44
1:A:1454:THR:HA	1:A:1457:LYS:HZ3	1.82	0.44
1:A:370:VAL:HG13	1:A:371:PHE:HD1	1.82	0.43
1:A:1097:ILE:HB	1:A:1098:PRO:HD3	2.00	0.43
1:A:1257:VAL:O	1:A:1260:HIS:NE2	2.51	0.43
1:A:41:SER:OG	1:A:142:GLU:OE2	2.22	0.43
1:A:789:ARG:HA	1:A:792:ILE:HG12	1.98	0.43
1:A:948:PRO:HA	1:A:951:VAL:HG12	2.00	0.43
1:A:712:LEU:HD13	1:A:768:TYR:CE2	2.52	0.43
1:A:615:LEU:HG	1:A:616:PRO:HD2	2.00	0.43
1:A:1319:LEU:HD23	1:A:1432:VAL:HG11	1.99	0.43
1:A:1360:GLU:HA	1:A:1404:THR:HA	2.01	0.43
1:A:824:ALA:HB2	1:A:834:GLN:HG3	2.00	0.43
1:A:66:TYR:HB3	1:A:196:PRO:N	2.34	0.43
1:A:1489:ASP:OD1	1:A:1489:ASP:N	2.51	0.43
1:A:250:PHE:HE1	1:A:375:MET:HE2	1.83	0.43
1:A:428:SER:O	1:A:432:GLN:HG3	2.18	0.43
1:A:659:THR:OG1	1:A:822:SER:O	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1126:SER:OG	1:A:1127:PRO:HD3	2.17	0.43
1:A:1158:ARG:HD3	1:A:1158:ARG:HA	1.88	0.43
1:A:1262:PRO:CG	1:A:1263:PRO:HD3	2.48	0.43
1:A:15:SER:OG	1:A:17:GLU:O	2.33	0.43
1:A:23:SER:HB3	1:A:29:THR:HG22	2.01	0.43
1:A:1259:GLY:HA3	1:A:1340:ILE:HD12	2.01	0.43
1:A:595:ILE:HD13	1:A:595:ILE:HA	1.93	0.43
1:A:1273:LYS:HE2	1:A:1295:HIS:NE2	2.34	0.43
1:A:1429:LYS:HD2	1:A:1429:LYS:H	1.84	0.43
1:A:45:MET:HE2	1:A:45:MET:HB3	1.67	0.42
1:A:699:VAL:HG13	1:A:699:VAL:O	2.19	0.42
1:A:1052:ARG:H	1:A:1052:ARG:HG2	1.71	0.42
1:A:1059:LYS:HA	1:A:1063:LEU:HB2	2.00	0.42
1:A:1358:VAL:HA	1:A:1422:ARG:HH22	1.83	0.42
1:A:353:PHE:HB2	1:A:1189:LEU:HB2	2.00	0.42
1:A:634:PHE:HD2	1:A:675:CYS:HB3	1.84	0.42
1:A:1035:ARG:HD2	1:A:1255:LEU:HD21	2.01	0.42
1:A:489:GLU:OE2	1:A:1359:PHE:HA	2.19	0.42
1:A:1342:LEU:HD13	1:A:1346:ARG:NH2	2.34	0.42
1:A:1495:LEU:HD12	1:A:1502:PHE:HE2	1.83	0.42
1:A:382:THR:HG23	1:A:409:MET:SD	2.59	0.42
1:A:731:LEU:O	1:A:731:LEU:HD23	2.19	0.42
1:A:558:ALA:HB2	1:A:983:GLY:O	2.20	0.42
1:A:632:ALA:O	1:A:648:ASN:N	2.53	0.42
1:A:33:ILE:HG21	1:A:149:LYS:HB2	2.02	0.42
1:A:930:GLU:HG2	1:A:931:GLN:N	2.35	0.42
1:A:935:LYS:HE3	1:A:935:LYS:HB3	1.92	0.42
1:A:1180:LEU:O	1:A:1184:GLY:N	2.46	0.42
1:A:1307:ARG:NH1	1:A:1310:ALA:HB2	2.35	0.42
1:A:390:LEU:HD23	1:A:1148:GLN:HB3	2.01	0.42
1:A:625:ALA:HA	1:A:654:LYS:HA	2.02	0.42
1:A:79:VAL:HG22	1:A:139:TRP:CH2	2.55	0.42
1:A:1429:LYS:HD2	1:A:1429:LYS:N	2.34	0.42
1:A:1456:PHE:HA	1:A:1459:ARG:HG3	2.02	0.42
1:A:710:ASN:HB3	1:A:761:LEU:HD11	2.02	0.42
1:A:1412:VAL:HA	1:A:1415:ARG:NH1	2.35	0.42
1:A:1490:SER:O	1:A:1494:LEU:HG	2.20	0.42
1:A:1503:TYR:O	1:A:1507:MET:HG3	2.20	0.42
1:A:139:TRP:HZ2	1:A:183:ILE:HG21	1.85	0.41
1:A:543:VAL:O	1:A:547:THR:HG22	2.20	0.41
1:A:832:THR:OG1	1:A:833:GLN:N	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1271:ASP:OD1	1:A:1271:ASP:N	2.53	0.41
1:A:315:VAL:HG21	1:A:1209:VAL:HG11	2.02	0.41
1:A:501:ALA:HB2	1:A:1346:ARG:HB2	2.02	0.41
1:A:597:ARG:O	1:A:600:THR:HG22	2.20	0.41
1:A:1420:LEU:HD21	1:A:1451:THR:CG2	2.47	0.41
1:A:92:ASN:OD1	1:A:168:THR:OG1	2.34	0.41
1:A:140:LEU:HD12	1:A:184:LEU:HD21	2.03	0.41
1:A:1352:ILE:HD12	1:A:1433:LEU:HD13	2.03	0.41
1:A:1362:THR:HB	1:A:1400:ASP:HA	2.01	0.41
1:A:74:SER:O	1:A:78:LEU:HG	2.19	0.41
1:A:1237:ILE:O	1:A:1237:ILE:HG13	2.20	0.41
1:A:481:ASP:HA	1:A:484:THR:HG22	2.01	0.41
1:A:1428:SER:C	1:A:1459:ARG:HH21	2.29	0.41
1:A:46:ILE:HD11	1:A:113:SER:HA	2.02	0.41
1:A:393:SER:OG	1:A:394:ASN:N	2.52	0.41
1:A:634:PHE:HZ	1:A:687:ALA:HB2	1.86	0.41
1:A:1404:THR:N	1:A:1408:GLY:O	2.51	0.41
1:A:246:LEU:HD13	1:A:250:PHE:HB3	2.03	0.41
1:A:386:TYR:OH	1:A:1131:HIS:NE2	2.45	0.41
1:A:658:LEU:HD11	1:A:809:LEU:HD22	2.02	0.41
1:A:1100:SER:O	1:A:1104:ILE:HG23	2.21	0.41
1:A:1107:GLN:HG3	1:A:1227:VAL:HG22	2.03	0.41
1:A:1321:ARG:NH2	1:A:1338:ASN:O	2.54	0.41
1:A:299:VAL:O	1:A:303:THR:HG23	2.21	0.41
1:A:1443:GLU:HA	1:A:1446:LYS:NZ	2.36	0.41
1:A:225:TRP:HH2	1:A:1171:ASN:HD22	1.67	0.40
1:A:286:LYS:HA	1:A:286:LYS:HD3	1.72	0.40
1:A:704:ASN:ND2	1:A:748:GLU:OE1	2.54	0.40
1:A:734:ASP:HA	1:A:737:ILE:HD12	2.02	0.40
1:A:463:SER:HA	1:A:466:MET:HG2	2.03	0.40
1:A:942:TYR:O	1:A:946:CYS:HB2	2.21	0.40
1:A:1358:VAL:HA	1:A:1422:ARG:NH2	2.37	0.40
1:A:83:ILE:HD13	1:A:114:LEU:HB3	2.03	0.40
1:A:581:ILE:N	1:A:582:PRO:HD2	2.36	0.40
1:A:628:ILE:HG13	1:A:652:GLN:OE1	2.21	0.40
1:A:1284:GLU:OE2	1:A:1284:GLU:N	2.49	0.40
1:A:89:ALA:O	1:A:93:ILE:HG12	2.21	0.40
1:A:496:SER:HB2	1:A:1357:GLN:OE1	2.21	0.40
1:A:498:LYS:HB3	1:A:1346:ARG:NH1	2.36	0.40
1:A:618:VAL:O	1:A:618:VAL:HG12	2.21	0.40
1:A:633:THR:HG22	1:A:647:LYS:HG3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:748:GLU:O	1:A:751:ILE:HG23	2.21	0.40
1:A:931:GLN:NE2	1:A:1116:GLU:OE2	2.52	0.40
1:A:1147:GLN:O	1:A:1147:GLN:HG3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1420/1537 (92%)	1322 (93%)	98 (7%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	1215/1353 (90%)	1215 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	HIS

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Mol	Chain	Res	Type
1	A	262	ASN
1	A	297	HIS
1	A	306	GLN
1	A	319	ASN
1	A	572	ASN
1	A	648	ASN
1	A	834	GLN
1	A	1074	ASN
1	A	1148	GLN
1	A	1299	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](#)

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.

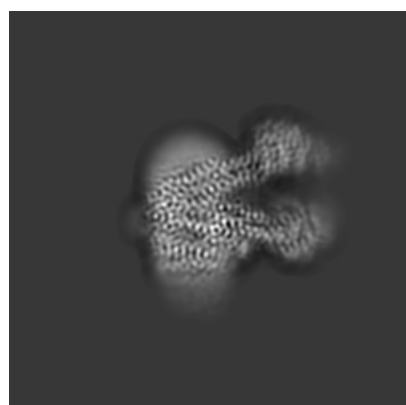
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23932. These allow visual inspection of the internal detail of the map and identification of artifacts.

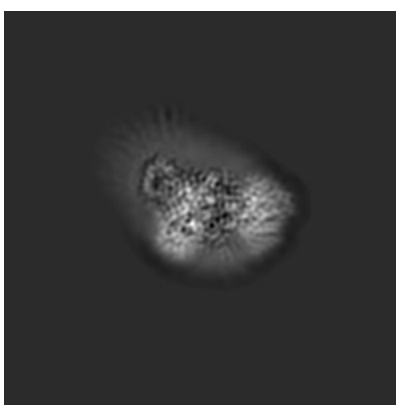
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

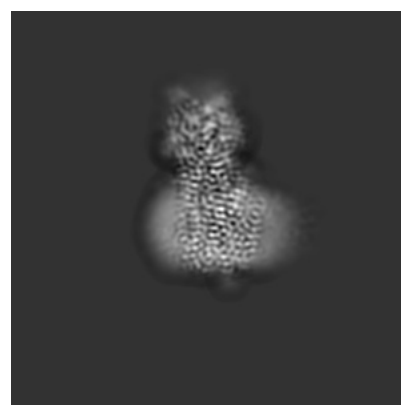
6.1.1 Primary map



X



Y

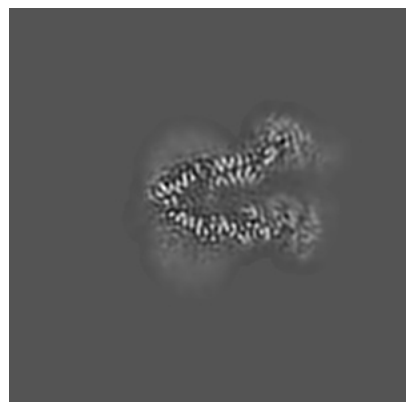


Z

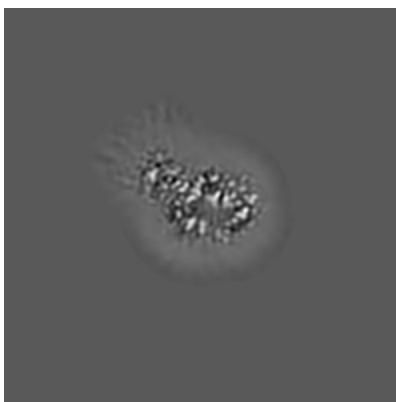
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 128



Y Index: 128

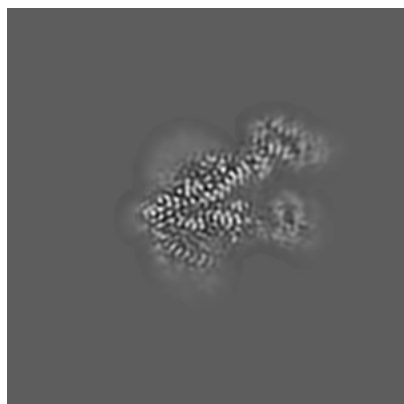


Z Index: 128

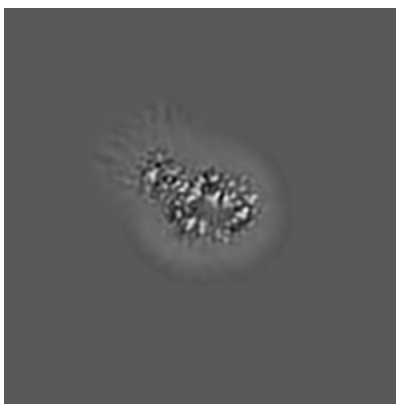
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

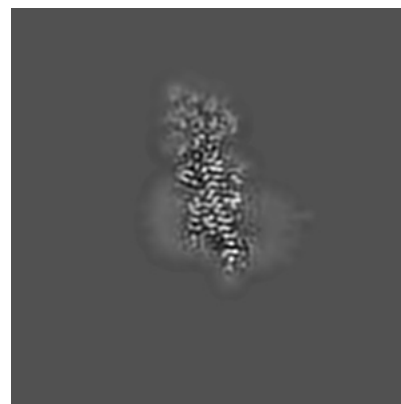
6.3.1 Primary map



X Index: 136



Y Index: 128

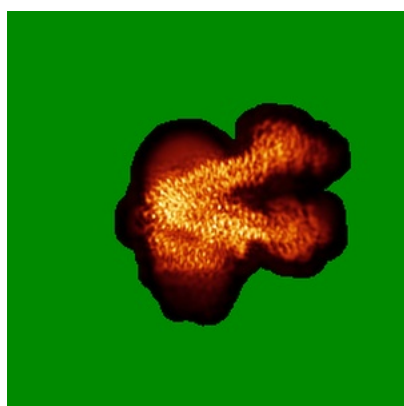


Z Index: 117

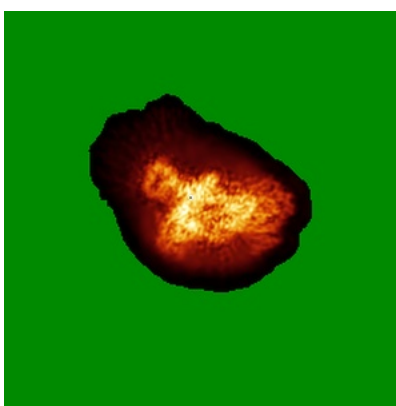
The images above show the largest variance slices of the map in three orthogonal directions.

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

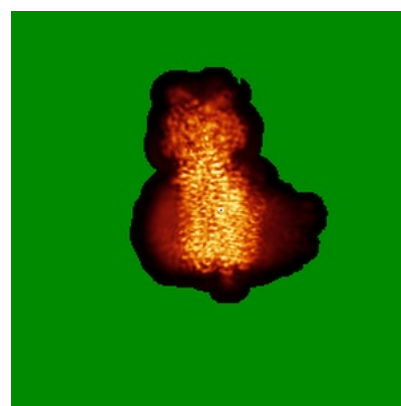
6.4.1 Primary map



X



Y

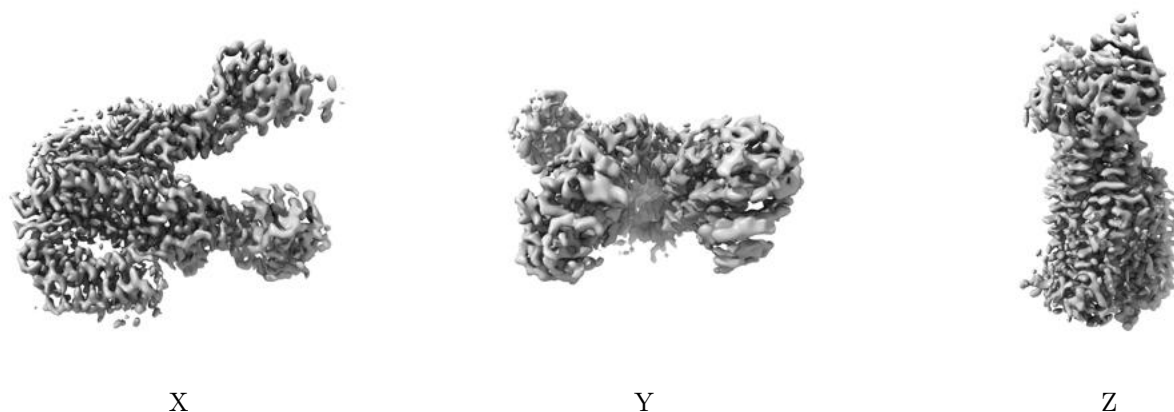


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.879. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

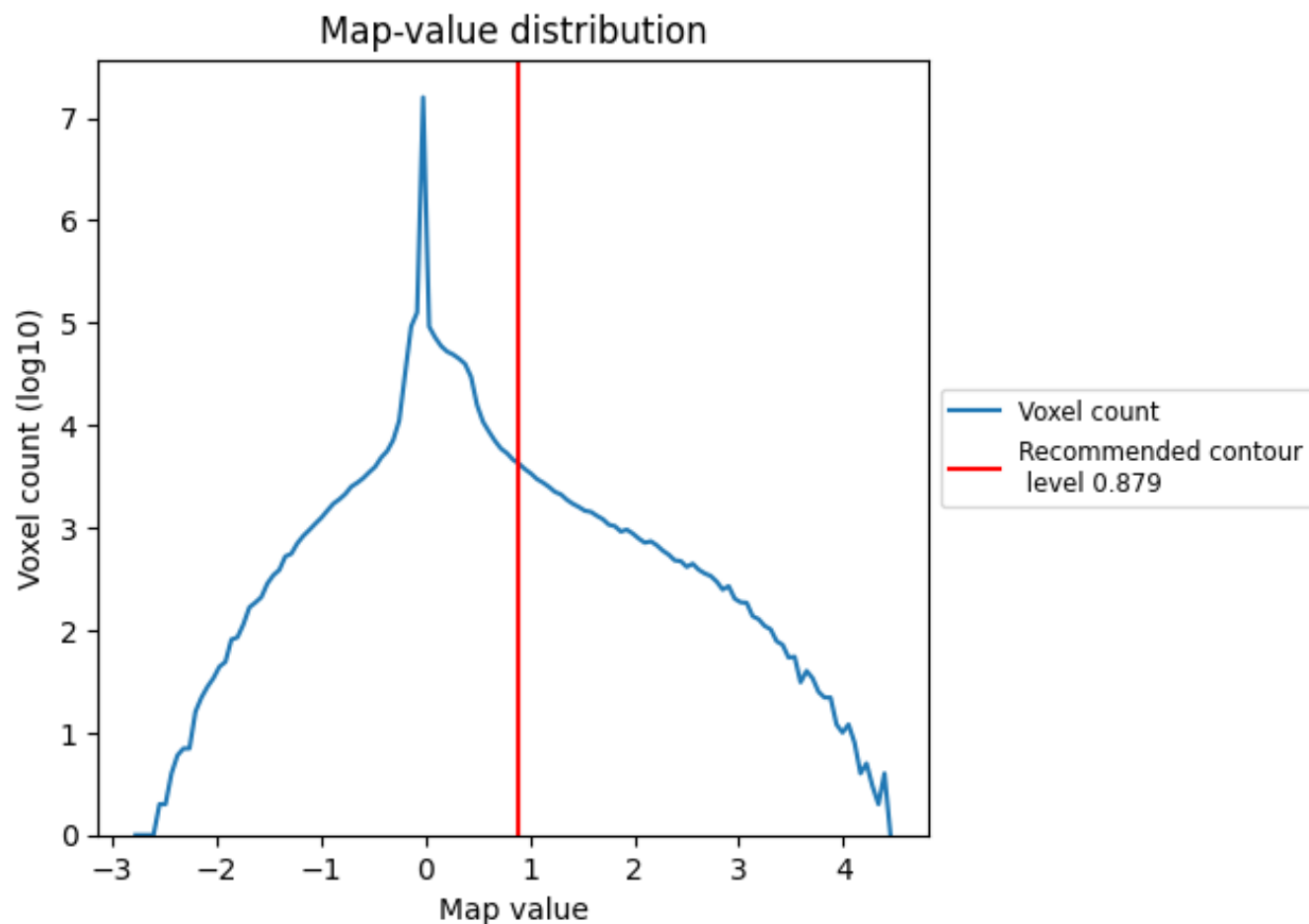
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

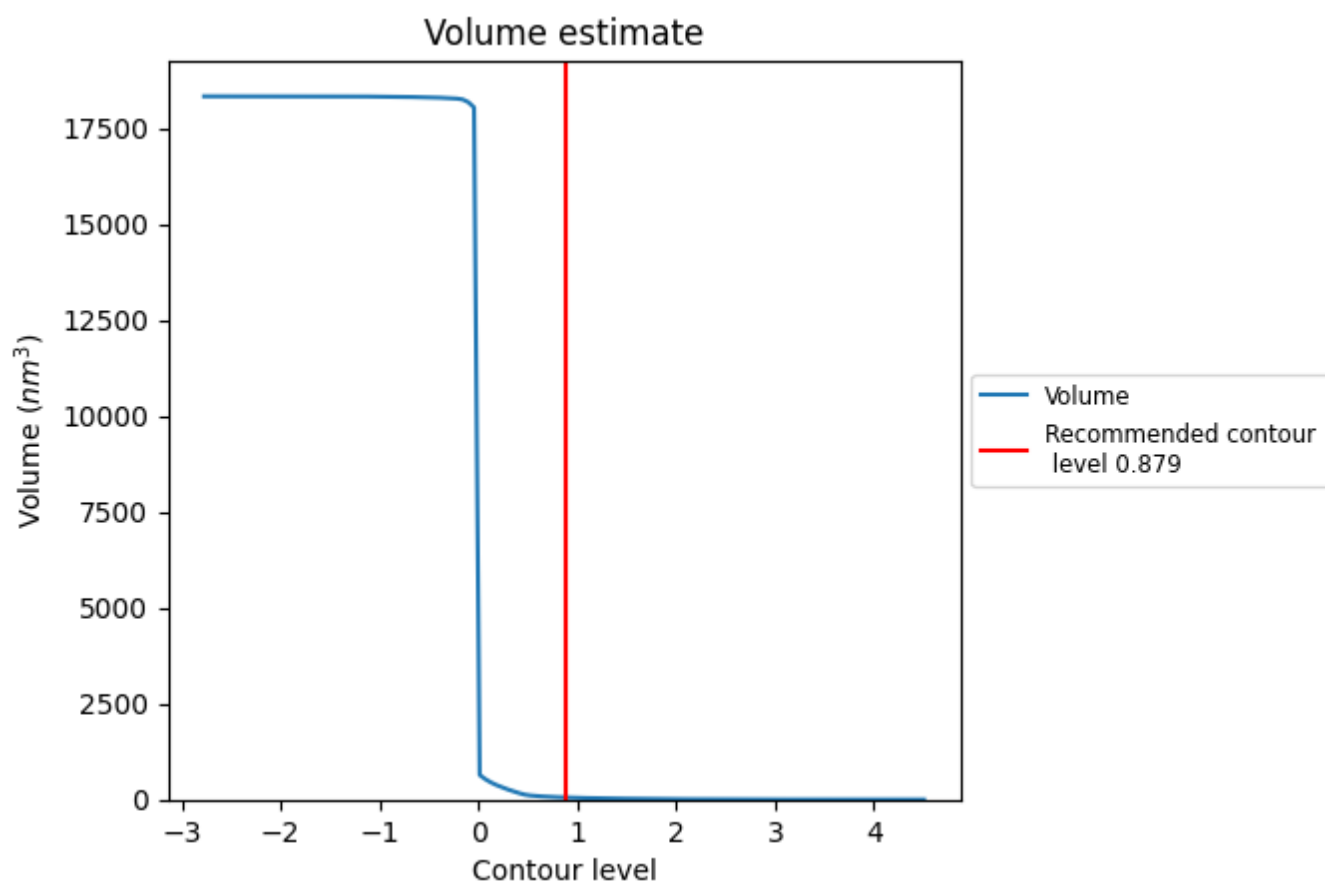
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

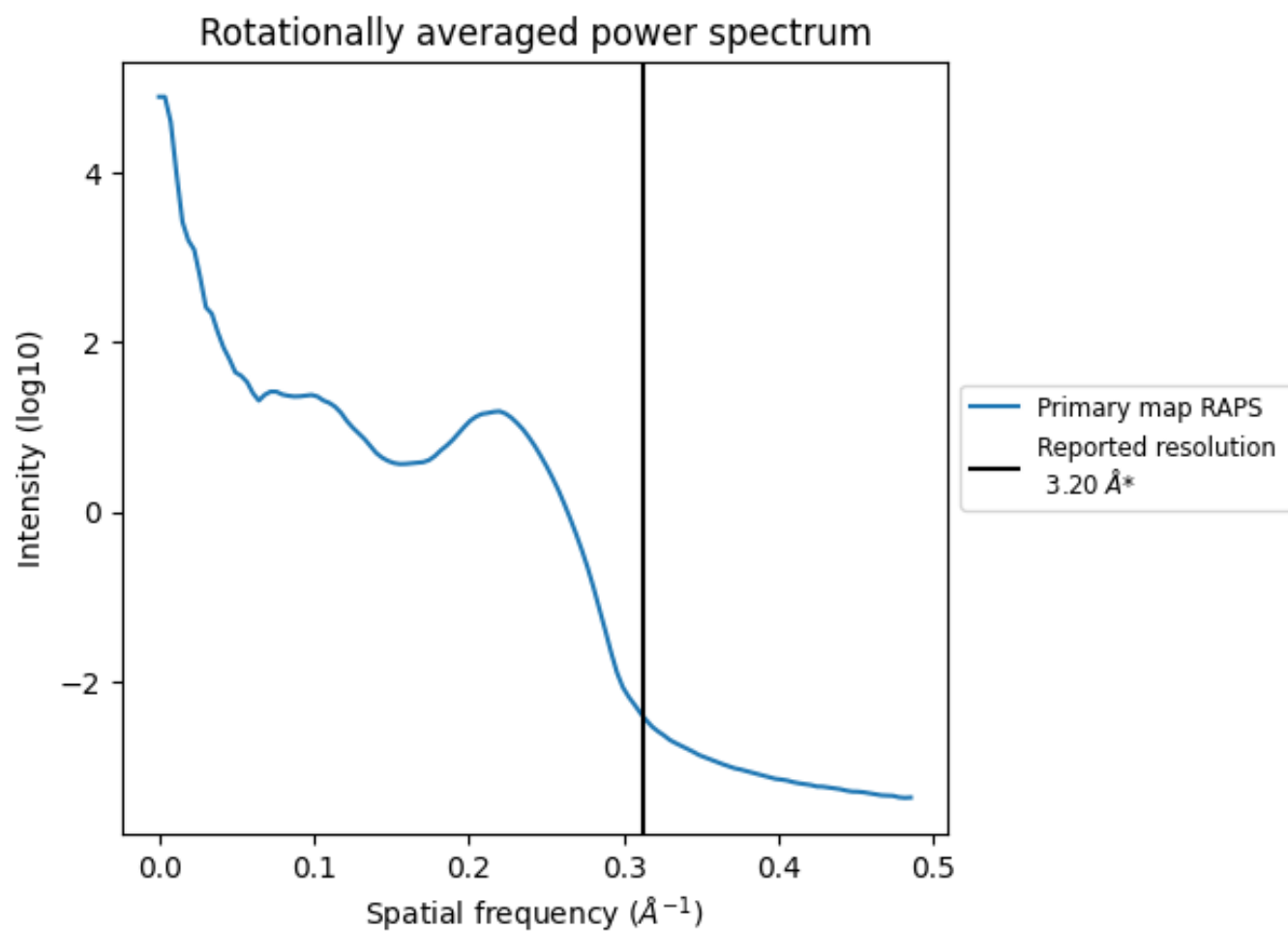
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 55 nm^3 ; this corresponds to an approximate mass of 49 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.312 Å⁻¹

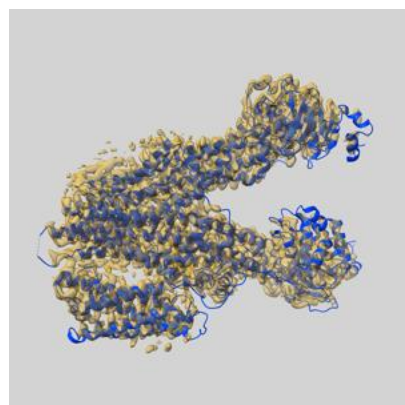
8 Fourier-Shell correlation ⓘ

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

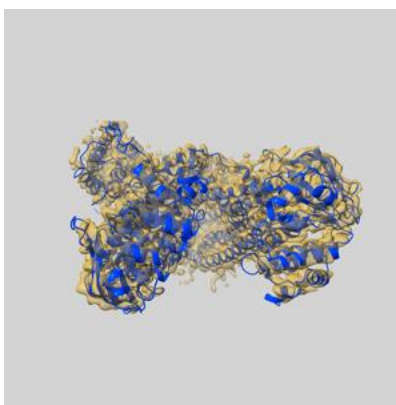
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-23932 and PDB model 7MPE. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

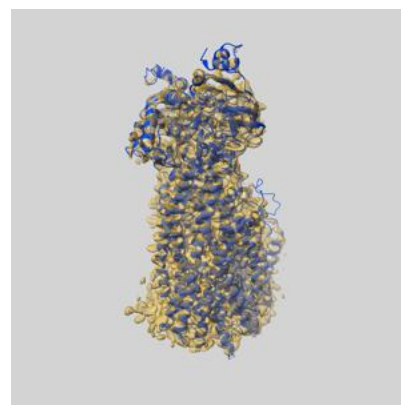
9.1 Map-model overlay [i](#)



X



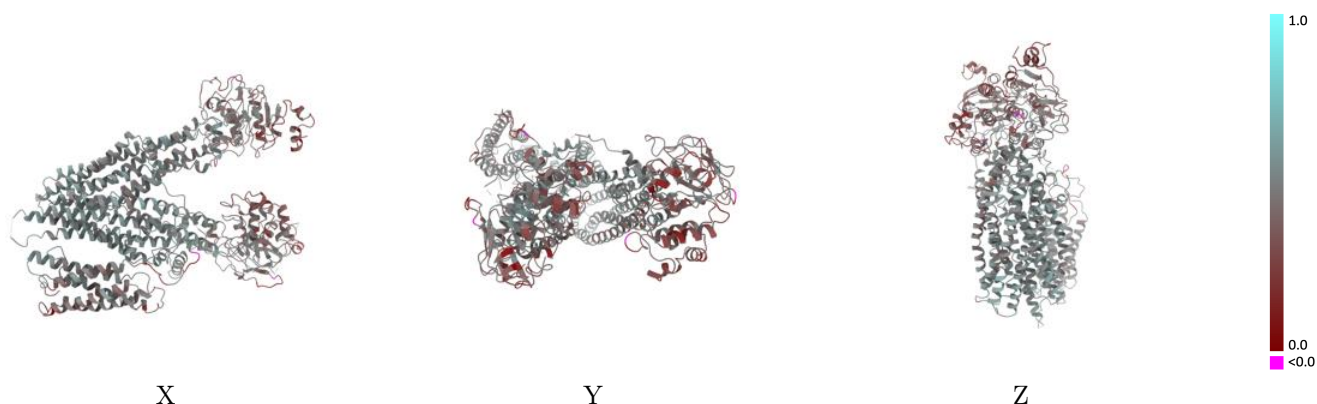
Y



Z

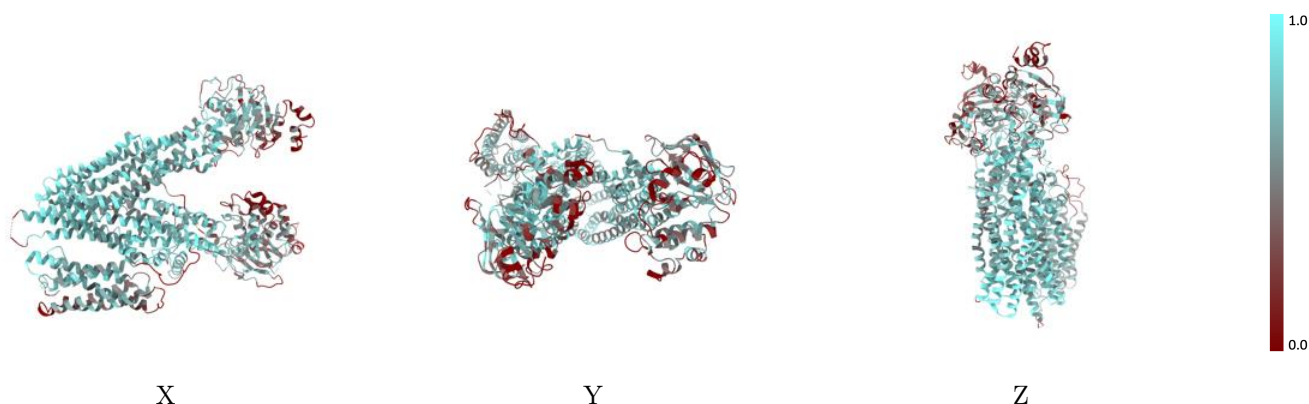
The images above show the 3D surface view of the map at the recommended contour level 0.879 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



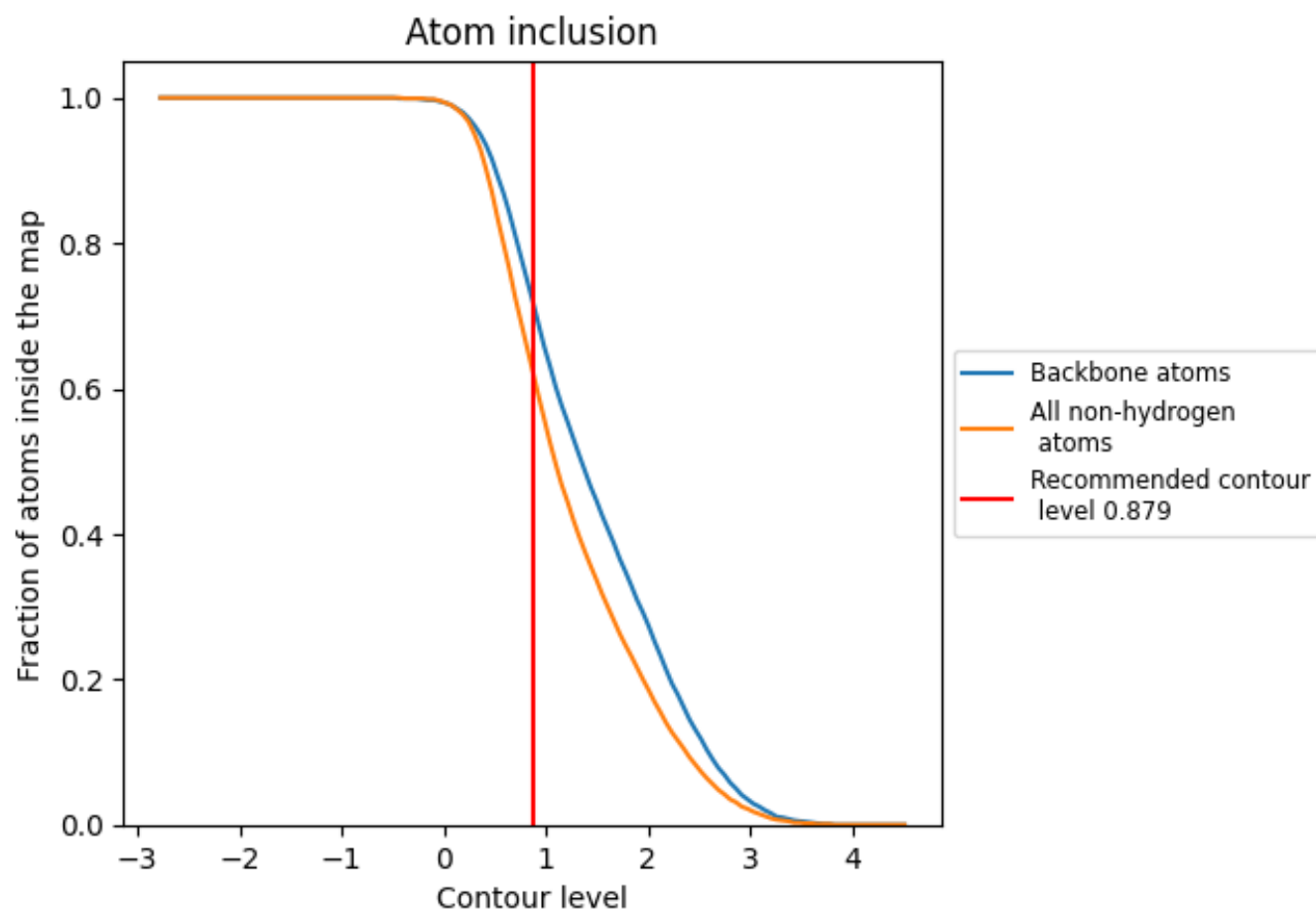
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.879).

9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 62% 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 (0.879) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6170	0.4560
A	0.6170	0.4560

