



Full wwPDB EM Validation Report ⓘ

May 26, 2024 – 12:54 PM EDT

PDB ID : 7S7B
EMDB ID : EMD-24882
Title : Human Nuclear exosome targeting (NEXT) complex homodimer bound to RNA (substrate 1)
Authors : Puno, M.R.; Lima, C.D.
Deposited on : 2021-09-15
Resolution : 4.06 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 ⓘ 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.dev92
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

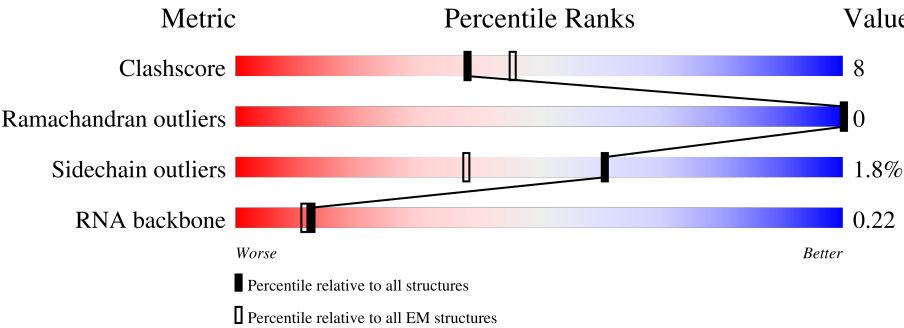
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.06 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.

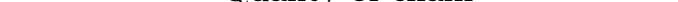


Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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	1045	70% 18% 12%
1	E	1045	70% 18% 12%
2	B	618	10% 48% 9% 43%
2	F	618	18% 46% 10% 43%
3	C	83	5% 66% 27% • 5%
3	G	83	48% 61% 33% • 5%
4	D	46	13% 11% 76%

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Mol	Chain	Length	Quality of chain
4	H	46	 <p>7% 9% 7% 9% 76%</p>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 22217 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 Exosome RNA helicase MTR4.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	924	Total	C	N	O	S	0	0
			7376	4700	1256	1369	51		
1	E	924	Total	C	N	O	S	0	0
			7376	4700	1256	1369	51		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP P42285
A	-1	GLY	-	expression tag	UNP P42285
A	0	ASP	-	expression tag	UNP P42285
E	-2	SER	-	expression tag	UNP P42285
E	-1	GLY	-	expression tag	UNP P42285
E	0	ASP	-	expression tag	UNP P42285

- Molecule 2 is a protein called Zinc finger CCHC domain-containing protein 8,Zinc finger CCHC domain-containing protein 8.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	351	Total	C	N	O	S	0	0
			2847	1802	486	544	15		
2	F	352	Total	C	N	O	S	0	0
			2856	1807	487	547	15		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	SER	-	expression tag	UNP Q6NZY4
B	-1	GLY	-	expression tag	UNP Q6NZY4
B	0	ASP	-	expression tag	UNP Q6NZY4
F	-2	SER	-	expression tag	UNP Q6NZY4
F	-1	GLY	-	expression tag	UNP Q6NZY4
F	0	ASP	-	expression tag	UNP Q6NZY4

- Molecule 3 is a protein called RNA-binding protein 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	79	Total	C	N	O	S	0	0
			644	424	109	110	1		
3	G	79	Total	C	N	O	S	0	0
			644	424	109	110	1		

There are 6 discrepancies between the modelled and reference sequences:

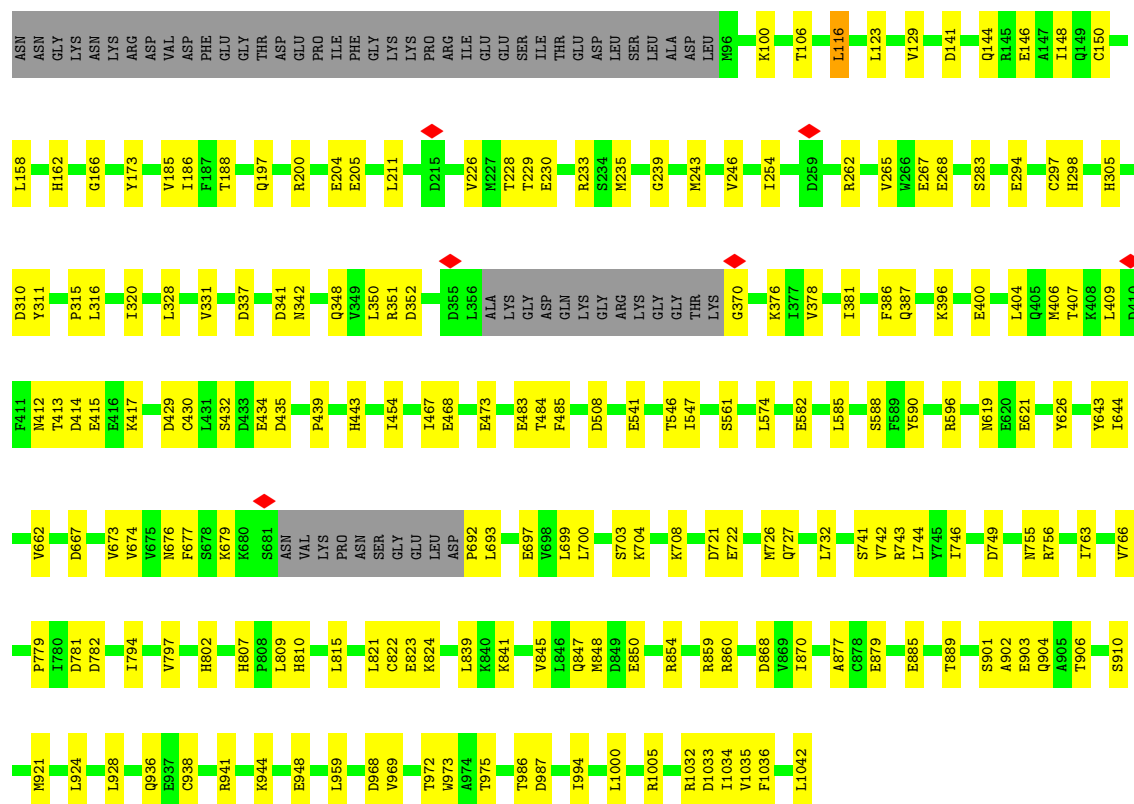
Chain	Residue	Modelled	Actual	Comment	Reference
C	4	SER	-	expression tag	UNP Q9Y580
C	5	GLY	-	expression tag	UNP Q9Y580
C	6	ASP	-	expression tag	UNP Q9Y580
G	4	SER	-	expression tag	UNP Q9Y580
G	5	GLY	-	expression tag	UNP Q9Y580
G	6	ASP	-	expression tag	UNP Q9Y580

- Molecule 4 is a RNA chain called RNA (46-MER).

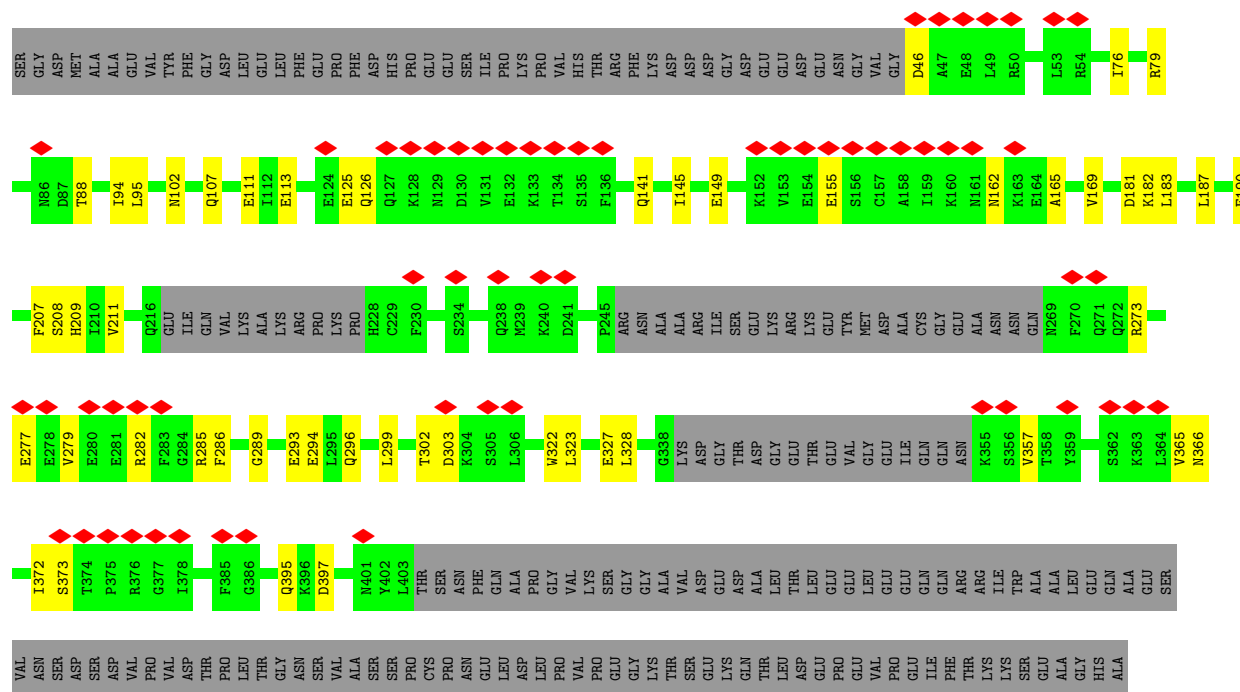
Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	11	Total	C	N	O	P	0	0
			236	107	46	72	11		
4	H	11	Total	C	N	O	P	0	0
			236	107	46	72	11		

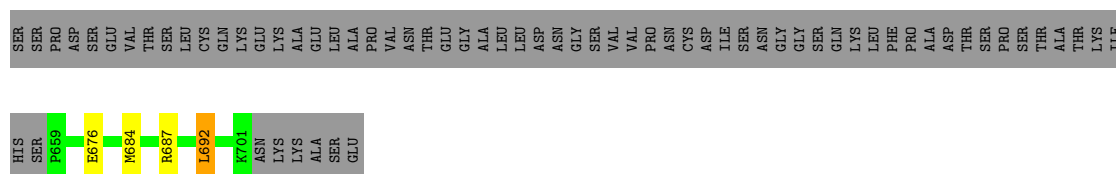
- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
5	B	1	Total	Zn	0
			1	1	
5	F	1	Total	Zn	0
			1	1	

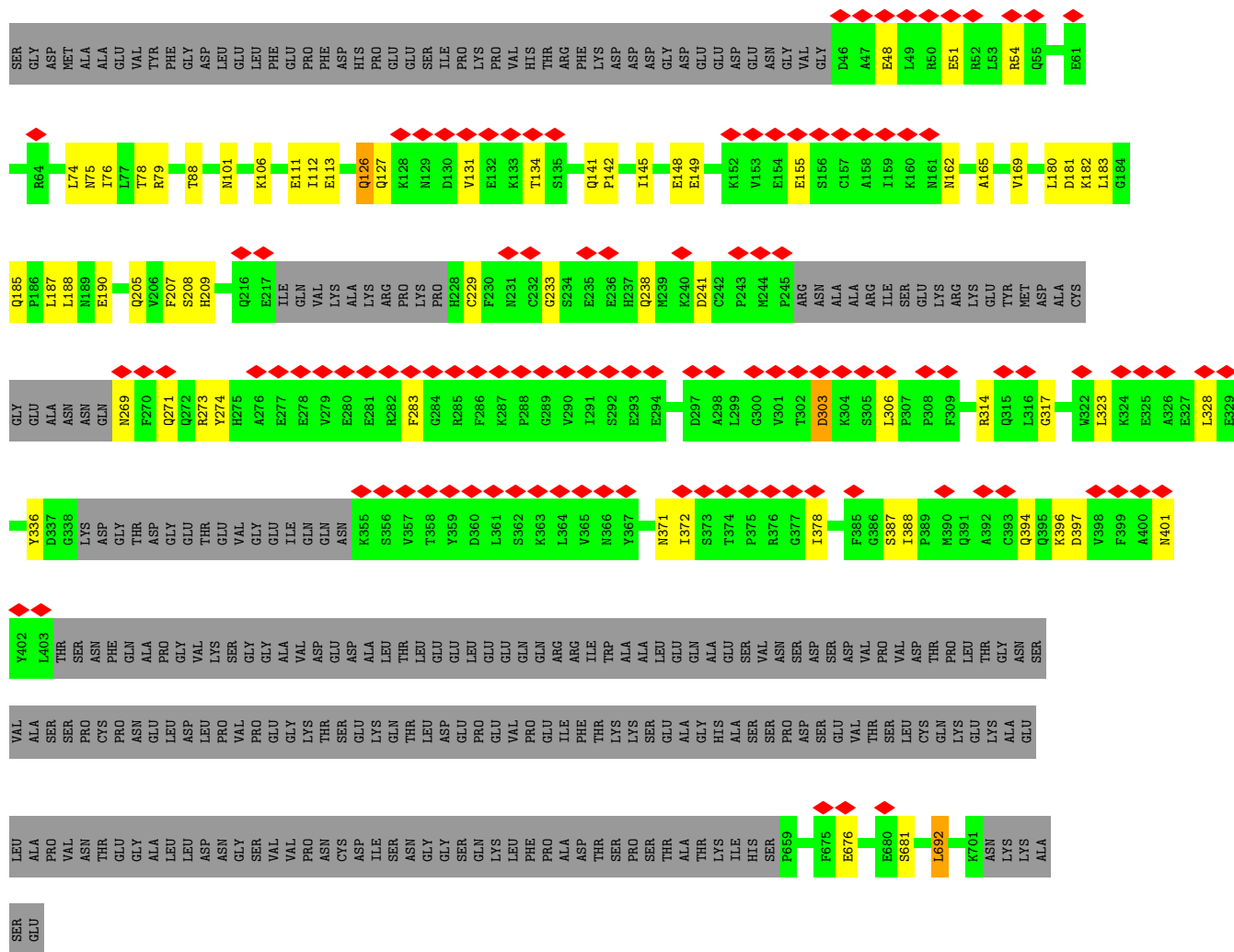


● Molecule 2: Zinc finger CCHC domain-containing protein 8, Zinc finger CCHC domain-containing protein 8





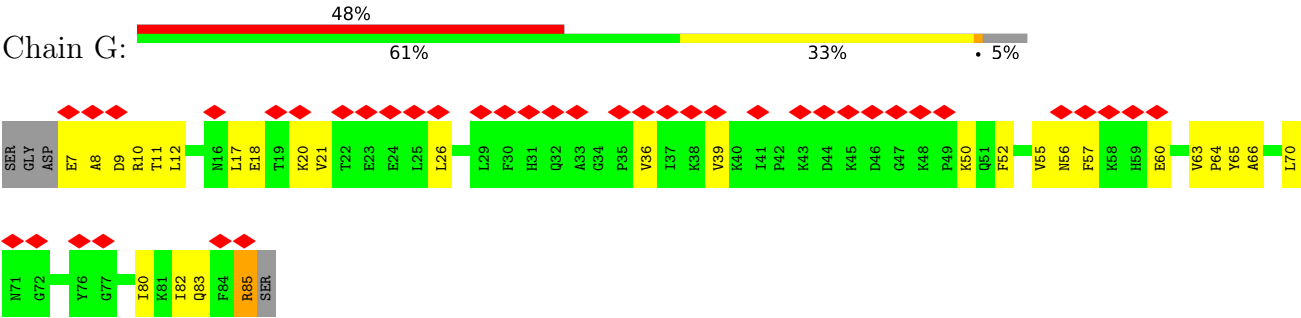
- Molecule 2: Zinc finger CCHC domain-containing protein 8, Zinc finger CCHC domain-containing protein 8



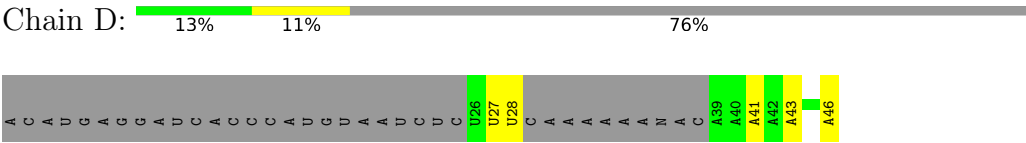
- Molecule 3: RNA-binding protein 7



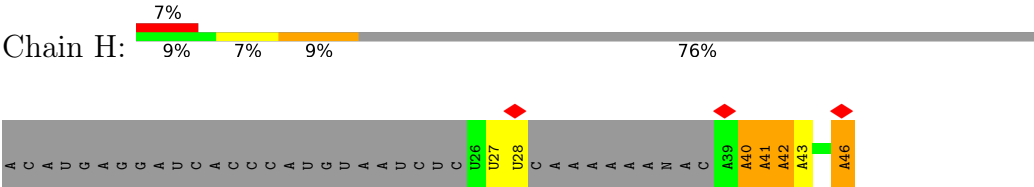
- Molecule 3: RNA-binding protein 7



• Molecule 4: RNA (46-MER)



• Molecule 4: RNA (46-MER)



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	618412	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	67.6	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	117.665	Depositor
Minimum map value	-58.579	Depositor
Average map value	-0.009	Depositor
Map value standard deviation	1.240	Depositor
Recommended contour level	11.5	Depositor
Map size (\AA)	417.79202, 417.79202, 417.79202	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.088, 1.088, 1.088	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: ZN

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	0/7516	0.47	0/10139
1	E	0.29	0/7516	0.47	0/10139
2	B	0.27	0/2907	0.46	0/3919
2	F	0.27	0/2916	0.46	0/3931
3	C	0.26	0/659	0.45	0/886
3	G	0.25	0/659	0.45	0/886
4	D	0.27	0/264	0.65	0/406
4	H	0.27	0/264	0.64	0/406
All	All	0.28	0/22701	0.47	0/30712

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	7376	0	7472	129	0
1	E	7376	0	7472	121	0
2	B	2847	0	2781	43	0
2	F	2856	0	2787	51	0
3	C	644	0	670	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	644	0	670	18	0
4	D	236	0	120	2	0
4	H	236	0	120	4	0
5	B	1	0	0	0	0
5	F	1	0	0	0	0
All	All	22217	0	22092	355	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 (355) 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:E:320:ILE:HD12	1:E:331:VAL:HG11	1.46	0.92
1:E:230:GLU:OE1	1:E:230:GLU:N	2.10	0.84
1:A:228:THR:OG1	4:D:46:A:OP1	1.96	0.83
1:A:230:GLU:N	1:A:230:GLU:OE1	2.12	0.83
1:E:267:GLU:OE2	1:E:561:SER:OG	1.98	0.82
1:E:859:ARG:NH2	1:E:868:ASP:OD1	2.12	0.82
1:E:676:ASN:OD1	1:E:677:PHE:N	2.10	0.81
1:E:396:LYS:NZ	4:H:43:A:OP2	2.13	0.81
1:A:676:ASN:OD1	1:A:677:PHE:N	2.14	0.80
1:A:903:GLU:N	1:A:903:GLU:OE1	2.14	0.80
1:E:903:GLU:N	1:E:903:GLU:OE1	2.15	0.80
1:E:370:GLY:O	1:E:376:LYS:NZ	2.12	0.79
2:B:95:LEU:HD21	2:F:112:ILE:HD13	1.63	0.79
2:B:111:GLU:OE1	2:B:111:GLU:N	2.16	0.78
1:E:972:THR:O	1:E:975:THR:OG1	2.02	0.77
3:G:9:ASP:O	3:G:85:ARG:NH2	2.18	0.77
1:E:400:GLU:OE2	2:F:273:ARG:NH1	2.19	0.76
2:F:269:ASN:N	2:F:271:GLN:OE1	2.19	0.76
3:C:39:VAL:HG12	3:C:55:VAL:HG12	1.66	0.76
1:A:727:GLN:OE1	1:A:756:ARG:NH1	2.19	0.75
1:A:879:GLU:OE1	1:A:879:GLU:N	2.19	0.75
2:B:285:ARG:NH1	2:B:286:PHE:O	2.19	0.75
1:E:1005:ARG:NH1	2:F:676:GLU:OE1	2.19	0.75
1:E:679:LYS:NZ	1:E:692:PRO:O	2.19	0.74
2:B:279:VAL:O	2:B:282:ARG:NH2	2.21	0.74
2:F:111:GLU:N	2:F:111:GLU:OE1	2.20	0.74
1:E:879:GLU:N	1:E:879:GLU:OE1	2.21	0.73
3:G:8:ALA:O	3:G:85:ARG:NE	2.21	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:60:GLU:OE1	3:C:60:GLU:N	2.20	0.73
1:E:877:ALA:HB2	1:E:889:THR:HG21	1.70	0.73
1:A:200:ARG:NH2	1:A:1042:LEU:O	2.21	0.73
2:F:190:GLU:N	2:F:190:GLU:OE1	2.22	0.73
1:A:1005:ARG:NH1	2:B:676:GLU:OE1	2.22	0.72
1:E:885:GLU:N	1:E:885:GLU:OE1	2.22	0.72
1:A:972:THR:O	1:A:975:THR:OG1	2.07	0.72
1:E:415:GLU:OE1	1:E:415:GLU:N	2.23	0.71
2:B:181:ASP:OD1	2:B:182:LYS:N	2.24	0.71
1:E:473:GLU:HG3	2:F:692:LEU:HD11	1.73	0.70
1:E:901:SER:N	1:E:904:GLN:OE1	2.24	0.70
2:F:317:GLY:O	3:G:65:TYR:OH	2.10	0.70
2:B:107:GLN:NE2	2:F:126:GLN:OE1	2.25	0.69
3:G:7:GLU:O	3:G:11:THR:HG23	1.92	0.69
3:C:9:ASP:O	3:C:85:ARG:NH2	2.26	0.68
1:E:351:ARG:NH2	1:E:708:LYS:O	2.27	0.68
1:A:415:GLU:N	1:A:415:GLU:OE1	2.27	0.68
1:A:370:GLY:O	1:A:376:LYS:NZ	2.15	0.67
1:A:229:THR:OG1	1:A:230:GLU:OE1	2.11	0.67
1:A:434:GLU:N	1:A:434:GLU:OE1	2.27	0.67
3:C:17:LEU:O	3:C:51:GLN:NE2	2.26	0.67
2:F:371:ASN:O	2:F:396:LYS:N	2.26	0.67
1:E:973:TRP:O	1:E:1032:ARG:NH2	2.27	0.67
1:A:859:ARG:NH2	1:A:868:ASP:OD1	2.28	0.66
3:G:39:VAL:HG12	3:G:55:VAL:HG12	1.77	0.66
1:A:877:ALA:HB2	1:A:889:THR:HG21	1.77	0.66
2:F:238:GLN:N	2:F:241:ASP:OD1	2.29	0.66
1:A:351:ARG:NH2	1:A:708:LYS:O	2.29	0.66
1:E:727:GLN:OE1	1:E:756:ARG:NH1	2.29	0.66
1:E:434:GLU:N	1:E:434:GLU:OE1	2.27	0.66
3:C:30:PHE:CE2	3:C:55:VAL:HG11	2.31	0.66
3:C:7:GLU:O	3:C:11:THR:HG23	1.96	0.65
1:A:807:HIS:ND1	1:A:809:LEU:HD22	2.11	0.65
1:A:885:GLU:OE1	1:A:885:GLU:N	2.30	0.65
1:E:294:GLU:OE1	1:E:596:ARG:NH1	2.30	0.65
2:F:169:VAL:HG23	2:F:169:VAL:O	1.97	0.64
2:B:190:GLU:N	2:B:190:GLU:OE1	2.31	0.64
2:F:141:GLN:NE2	2:F:183:LEU:O	2.29	0.64
2:F:314:ARG:NE	2:F:387:SER:OG	2.29	0.64
1:A:473:GLU:HG3	2:B:692:LEU:HD11	1.79	0.64
3:G:18:GLU:OE1	3:G:20:LYS:N	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:742:VAL:HG23	1:E:742:VAL:O	1.99	0.63
1:A:241:GLU:N	1:A:241:GLU:OE1	2.31	0.63
1:A:884:ASP:OD1	1:A:885:GLU:N	2.31	0.63
2:F:74:LEU:HD23	2:F:74:LEU:O	1.99	0.63
1:A:849:ASP:OD1	1:A:850:GLU:N	2.32	0.62
1:E:673:VAL:HG11	1:E:746:ILE:HD13	1.81	0.62
1:A:404:LEU:HD11	2:B:277:GLU:OE2	1.99	0.62
2:F:397:ASP:O	2:F:401:ASN:ND2	2.32	0.62
1:A:396:LYS:NZ	4:D:43:A:OP2	2.26	0.62
2:F:303:ASP:N	2:F:303:ASP:OD1	2.33	0.61
1:A:742:VAL:HG23	1:A:742:VAL:O	2.01	0.61
1:E:412:ASN:O	1:E:417:LYS:NZ	2.32	0.61
1:E:483:GLU:OE2	4:H:42:A:O2'	2.14	0.61
1:E:574:LEU:HB2	1:E:870:ILE:HD11	1.83	0.61
1:A:267:GLU:OE2	1:A:561:SER:OG	2.12	0.61
1:A:617:ILE:HG12	1:A:821:LEU:HD11	1.83	0.60
1:E:200:ARG:NH2	1:E:1042:LEU:O	2.34	0.60
2:F:88:THR:HG23	2:F:165:ALA:HB2	1.82	0.60
1:E:744:LEU:HD21	1:E:766:VAL:HG21	1.84	0.60
1:A:809:LEU:HD23	1:A:810:HIS:N	2.17	0.59
2:F:208:SER:O	2:F:209:HIS:ND1	2.34	0.59
2:F:314:ARG:O	2:F:394:GLN:NE2	2.35	0.59
1:A:626:TYR:OH	1:A:630:ARG:NH2	2.36	0.59
3:C:18:GLU:OE1	3:C:19:THR:N	2.35	0.59
3:C:12:LEU:HD21	3:C:82:ILE:HG21	1.85	0.59
1:A:973:TRP:O	1:A:1032:ARG:NH2	2.35	0.59
1:E:262:ARG:O	1:E:265:VAL:HG22	2.02	0.58
1:E:404:LEU:HD21	2:F:274:TYR:HA	1.84	0.58
2:B:94:ILE:HG22	2:B:95:LEU:H	1.68	0.58
1:E:994:ILE:CD1	1:E:1035:VAL:HG12	2.33	0.58
1:A:578:GLU:N	1:A:578:GLU:OE1	2.37	0.58
3:C:11:THR:HG22	3:C:56:ASN:OD1	2.03	0.58
3:C:72:GLY:N	3:C:80:ILE:O	2.36	0.58
1:E:162:HIS:NE2	1:E:310:ASP:OD1	2.36	0.58
1:E:348:GLN:NE2	1:E:352:ASP:OD1	2.36	0.58
3:G:11:THR:HG22	3:G:56:ASN:OD1	2.04	0.57
1:E:197:GLN:NE2	1:E:468:GLU:OE1	2.38	0.57
1:E:341:ASP:OD1	1:E:342:ASN:N	2.38	0.57
1:A:938:CYS:SG	1:A:941:ARG:NH2	2.78	0.57
1:E:141:ASP:OD1	1:E:141:ASP:N	2.36	0.57
2:F:78:THR:HG23	2:F:148:GLU:O	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:621:GLU:OE1	1:E:621:GLU:N	2.35	0.57
1:E:229:THR:OG1	1:E:230:GLU:OE1	2.16	0.56
3:G:57:PHE:CD1	3:G:63:VAL:HG22	2.40	0.56
2:B:76:ILE:HD12	2:F:145:ILE:HD11	1.88	0.56
2:F:79:ARG:NH2	2:F:149:GLU:OE1	2.38	0.56
1:A:644:ILE:HD11	1:A:779:PRO:HG3	1.88	0.56
1:A:621:GLU:OE1	1:A:621:GLU:N	2.37	0.55
1:A:266:TRP:O	1:A:269:THR:OG1	2.19	0.55
1:A:341:ASP:OD1	1:A:342:ASN:N	2.37	0.55
1:E:936:GLN:OE1	1:E:959:LEU:HD12	2.06	0.55
1:A:233:ARG:NH2	1:A:568:ASN:OD1	2.39	0.55
2:B:88:THR:HG23	2:B:165:ALA:HB2	1.88	0.55
2:B:299:LEU:HA	3:C:75:LEU:HD13	1.88	0.55
1:E:644:ILE:HD11	1:E:779:PRO:HG3	1.87	0.55
1:E:823:GLU:OE1	1:E:824:LYS:N	2.39	0.55
1:A:747:PRO:HD2	1:A:750:LEU:HD21	1.88	0.55
1:E:582:GLU:OE1	1:E:582:GLU:N	2.35	0.55
1:E:821:LEU:HD12	1:E:822:CYS:N	2.21	0.55
1:A:679:LYS:HZ2	1:A:692:PRO:HB2	1.72	0.54
1:E:400:GLU:O	1:E:404:LEU:HD23	2.06	0.54
1:A:602:VAL:HG22	1:A:839:LEU:HD11	1.90	0.54
1:E:741:SER:O	2:F:180:LEU:HD23	2.07	0.54
1:A:662:VAL:HG11	1:A:700:LEU:HD11	1.90	0.54
2:B:145:ILE:HD11	2:F:76:ILE:HD12	1.90	0.54
1:A:782:ASP:OD1	2:F:75:ASN:ND2	2.38	0.54
1:E:228:THR:OG1	4:H:46:A:OP1	2.19	0.54
2:F:127:GLN:HA	2:F:131:VAL:HG22	1.90	0.53
1:E:877:ALA:CB	1:E:889:THR:HG21	2.38	0.53
1:E:994:ILE:HD12	1:E:1035:VAL:HA	1.91	0.53
2:F:207:PHE:O	2:F:208:SER:OG	2.20	0.53
2:B:208:SER:O	2:B:209:HIS:ND1	2.41	0.53
3:G:17:LEU:HA	3:G:80:ILE:HG22	1.89	0.53
1:E:699:LEU:O	1:E:699:LEU:HD23	2.09	0.53
1:E:1033:ASP:OD1	1:E:1034:ILE:N	2.40	0.53
2:F:371:ASN:OD1	2:F:372:ILE:N	2.41	0.53
2:F:229:CYS:O	2:F:233:GLY:N	2.36	0.53
1:A:150:CYS:HG	1:A:305:HIS:CG	2.27	0.53
2:F:314:ARG:NE	2:F:388:ILE:O	2.41	0.53
2:B:169:VAL:HG23	2:B:169:VAL:O	2.09	0.53
1:A:699:LEU:O	1:A:699:LEU:HD23	2.09	0.52
1:E:699:LEU:HD21	1:E:726:MET:HG3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:235:MET:O	1:E:239:GLY:N	2.42	0.52
1:E:233:ARG:NE	1:E:268:GLU:OE1	2.38	0.52
1:A:294:GLU:OE1	1:A:596:ARG:NH1	2.39	0.52
1:A:994:ILE:CD1	1:A:1035:VAL:HG12	2.39	0.52
1:A:141:ASP:OD1	1:A:141:ASP:N	2.38	0.52
1:E:320:ILE:CD1	1:E:331:VAL:HG11	2.29	0.52
3:C:9:ASP:OD1	3:C:10:ARG:N	2.43	0.52
1:A:877:ALA:CB	1:A:889:THR:HG21	2.40	0.51
1:A:626:TYR:CD1	1:A:809:LEU:HD21	2.45	0.51
1:A:1033:ASP:OD1	1:A:1034:ILE:N	2.43	0.51
1:A:925:THR:OG1	1:A:926:GLU:N	2.42	0.51
1:E:439:PRO:O	1:E:443:HIS:ND1	2.42	0.51
1:A:230:GLU:OE2	1:A:262:ARG:NH2	2.43	0.51
1:A:259:ASP:O	1:A:261:GLU:N	2.43	0.50
1:A:465:GLU:HB3	1:A:1040:LEU:HD11	1.93	0.50
1:E:404:LEU:O	1:E:407:THR:HG22	2.12	0.50
1:A:918:SER:O	1:A:965:HIS:NE2	2.43	0.50
1:A:162:HIS:NE2	1:A:310:ASP:OD1	2.37	0.50
1:A:594:HIS:NE2	1:A:846:LEU:HD21	2.26	0.50
1:A:753:VAL:O	1:A:753:VAL:HG22	2.12	0.50
1:E:673:VAL:HG23	1:E:763:ILE:HD11	1.94	0.50
1:A:279:TYR:OH	1:A:300:HIS:NE2	2.42	0.49
1:A:721:ASP:OD1	1:A:722:GLU:N	2.46	0.49
2:B:125:GLU:OE1	2:B:126:GLN:N	2.46	0.49
1:E:749:ASP:O	1:E:755:ASN:ND2	2.41	0.49
1:E:815:LEU:HD12	1:E:815:LEU:C	2.33	0.49
2:B:322:TRP:NE1	3:C:31:HIS:O	2.45	0.49
2:F:181:ASP:OD1	2:F:182:LYS:N	2.46	0.49
1:E:994:ILE:HD13	1:E:1035:VAL:HG12	1.95	0.49
1:E:243:MET:O	1:E:246:VAL:HG22	2.13	0.48
1:A:412:ASN:O	1:A:417:LYS:NZ	2.46	0.48
1:E:148:ILE:HD12	1:E:173:TYR:CG	2.48	0.48
3:G:66:ALA:O	3:G:70:LEU:N	2.39	0.48
1:A:125:LEU:HD21	1:A:305:HIS:HE1	1.79	0.48
1:A:188:THR:HG21	1:A:229:THR:HG22	1.95	0.48
1:E:845:VAL:HG23	1:E:848:MET:HB2	1.95	0.48
4:H:40:A:O2'	4:H:41:A:OP2	2.22	0.48
1:A:673:VAL:HG11	1:A:746:ILE:HD13	1.95	0.48
1:A:508:ASP:OD1	1:A:508:ASP:N	2.47	0.48
1:A:673:VAL:HG23	1:A:763:ILE:HD11	1.95	0.48
1:A:378:VAL:HG11	1:A:406:MET:SD	2.54	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:876:VAL:HG21	1:A:1014:ILE:CG2	2.44	0.48
1:A:699:LEU:HD21	1:A:726:MET:HG3	1.96	0.47
1:E:673:VAL:HG12	1:E:674:VAL:O	2.14	0.47
2:B:296:GLN:NE2	2:B:302:THR:O	2.47	0.47
2:B:395:GLN:NE2	2:B:397:ASP:OD2	2.47	0.47
1:A:400:GLU:OE2	2:B:273:ARG:NH1	2.47	0.47
3:C:30:PHE:CD2	3:C:55:VAL:HG11	2.49	0.47
2:B:141:GLN:NE2	2:B:183:LEU:O	2.40	0.47
1:E:1042:LEU:O	2:F:681:SER:OG	2.31	0.47
1:E:337:ASP:OD1	1:E:337:ASP:N	2.48	0.47
1:E:429:ASP:OD1	1:E:430:CYS:N	2.48	0.47
1:E:741:SER:O	1:E:741:SER:OG	2.32	0.47
1:E:673:VAL:HG11	1:E:746:ILE:CD1	2.45	0.47
1:A:588:SER:O	1:A:590:TYR:N	2.49	0.46
1:A:1035:VAL:HG23	1:A:1036:PHE:CD1	2.50	0.46
1:E:200:ARG:NH1	1:E:204:GLU:OE2	2.46	0.46
1:A:743:ARG:NH2	2:F:113:GLU:OE2	2.42	0.46
3:G:12:LEU:HD12	3:G:83:GLN:O	2.15	0.46
1:A:901:SER:N	1:A:904:GLN:OE1	2.49	0.46
3:C:11:THR:O	3:C:85:ARG:NH2	2.43	0.46
3:G:60:GLU:OE1	3:G:60:GLU:N	2.41	0.46
1:A:839:LEU:C	1:A:839:LEU:HD23	2.36	0.46
1:E:315:PRO:C	1:E:316:LEU:HD22	2.35	0.46
1:A:815:LEU:C	1:A:815:LEU:HD12	2.35	0.46
1:E:144:GLN:NE2	1:E:166:GLY:O	2.49	0.46
1:A:910:SER:HB3	1:A:1000:LEU:HD22	1.98	0.46
2:B:293:GLU:OE1	2:B:294:GLU:N	2.49	0.46
1:E:588:SER:O	1:E:590:TYR:N	2.48	0.46
1:E:508:ASP:OD1	1:E:508:ASP:N	2.47	0.46
1:E:619:ASN:N	1:E:619:ASN:OD1	2.48	0.46
1:A:680:LYS:O	1:A:681:SER:OG	2.21	0.46
1:E:781:ASP:OD1	1:E:782:ASP:N	2.49	0.46
1:E:188:THR:HG21	1:E:229:THR:HG22	1.98	0.46
1:E:968:ASP:OD1	1:E:969:VAL:N	2.47	0.46
1:A:821:LEU:C	1:A:821:LEU:HD12	2.36	0.45
1:A:928:LEU:N	1:A:928:LEU:HD12	2.31	0.45
3:G:36:VAL:HG22	3:G:57:PHE:CD2	2.51	0.45
2:B:113:GLU:OE2	1:E:743:ARG:NH2	2.49	0.45
2:F:306:LEU:HD13	2:F:378:ILE:HG21	1.98	0.45
2:F:155:GLU:OE2	2:F:162:ASN:ND2	2.49	0.45
1:A:703:SER:O	1:A:704:LYS:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:46:ASP:OD1	2:B:46:ASP:N	2.49	0.45
1:A:320:ILE:HD12	1:A:331:VAL:HG11	1.98	0.45
2:B:323:LEU:O	2:B:323:LEU:HD23	2.16	0.45
1:E:378:VAL:HG11	1:E:406:MET:SD	2.57	0.45
1:E:860:ARG:NH1	1:E:948:GLU:OE2	2.50	0.45
1:A:409:LEU:N	1:A:409:LEU:HD12	2.31	0.45
1:E:910:SER:HB3	1:E:1000:LEU:HD22	1.99	0.45
2:F:141:GLN:HB2	2:F:142:PRO:HD2	1.99	0.45
1:A:211:LEU:HD12	1:A:226:VAL:O	2.17	0.44
1:A:262:ARG:O	1:A:265:VAL:HG22	2.17	0.44
1:A:968:ASP:OD1	1:A:969:VAL:N	2.49	0.44
1:E:254:ILE:HG22	1:E:283:SER:HB3	1.98	0.44
2:B:145:ILE:CD1	2:F:76:ILE:HD12	2.47	0.44
2:B:303:ASP:OD1	2:B:303:ASP:N	2.50	0.44
1:E:986:THR:OG1	1:E:987:ASP:N	2.48	0.44
3:G:9:ASP:OD1	3:G:10:ARG:N	2.50	0.44
1:A:469:ILE:HD11	1:A:1040:LEU:HD13	1.98	0.44
1:A:673:VAL:HG11	1:A:746:ILE:CD1	2.47	0.44
1:E:839:LEU:C	1:E:839:LEU:HD23	2.38	0.44
1:E:938:CYS:SG	1:E:941:ARG:NH2	2.91	0.44
1:A:927:GLN:C	1:A:928:LEU:HD12	2.37	0.44
1:E:350:LEU:HD11	1:E:541:GLU:HA	1.99	0.44
1:E:407:THR:HG23	1:E:407:THR:O	2.18	0.44
1:E:454:ILE:O	1:E:467:ILE:HD13	2.18	0.44
3:G:18:GLU:HB3	3:G:21:VAL:HG23	1.99	0.44
1:E:1035:VAL:HG23	1:E:1036:PHE:CD2	2.52	0.44
2:B:372:ILE:HG22	2:B:373:SER:N	2.32	0.44
1:A:231:ILE:O	1:A:235:MET:HG3	2.17	0.43
1:E:123:LEU:HD23	1:E:123:LEU:H	1.83	0.43
1:E:106:THR:OG1	1:E:146:GLU:OE2	2.35	0.43
1:E:703:SER:O	1:E:704:LYS:HB3	2.17	0.43
2:F:131:VAL:HG21	2:F:134:THR:HG21	1.99	0.43
1:A:698:VAL:HG23	1:A:700:LEU:HG	2.00	0.43
1:A:407:THR:O	1:A:407:THR:HG23	2.18	0.43
2:B:327:GLU:OE2	2:B:357:VAL:HG11	2.18	0.43
1:E:662:VAL:HG11	1:E:700:LEU:HD11	2.00	0.43
1:E:903:GLU:O	1:E:906:THR:OG1	2.33	0.43
1:E:721:ASP:OD1	1:E:722:GLU:N	2.49	0.43
1:A:994:ILE:HD12	1:A:1035:VAL:HG12	1.99	0.43
2:B:76:ILE:CD1	2:F:145:ILE:HD11	2.48	0.43
2:B:79:ARG:NH2	2:B:149:GLU:OE1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:158:LEU:HD23	1:E:158:LEU:O	2.19	0.43
1:E:409:LEU:HD12	1:E:409:LEU:N	2.34	0.43
1:E:211:LEU:HD12	1:E:226:VAL:O	2.19	0.43
1:A:243:MET:O	1:A:246:VAL:HG22	2.19	0.43
1:E:413:THR:HG22	1:E:414:ASP:N	2.34	0.43
1:A:252:ASP:OD1	1:A:253:GLU:N	2.51	0.43
1:A:318:HIS:O	1:A:331:VAL:HG22	2.19	0.43
1:A:428:ILE:O	1:A:428:ILE:HG22	2.19	0.43
1:E:697:GLU:OE1	2:F:205:GLN:NE2	2.52	0.43
1:E:902:ALA:O	1:E:906:THR:HG23	2.18	0.43
2:F:323:LEU:HD23	2:F:323:LEU:O	2.19	0.43
1:A:523:ARG:HE	1:A:523:ARG:HA	1.84	0.42
1:A:254:ILE:HG22	1:A:283:SER:HB3	2.02	0.42
1:A:565:LEU:HD21	1:A:585:LEU:HD21	2.01	0.42
1:E:100:LYS:O	1:E:116:LEU:HD23	2.20	0.42
1:A:167:LYS:NZ	1:A:283:SER:O	2.32	0.42
2:B:155:GLU:OE2	2:B:162:ASN:ND2	2.53	0.42
2:B:207:PHE:O	2:B:208:SER:OG	2.25	0.42
2:B:366:ASN:HB3	2:B:372:ILE:HD12	2.02	0.42
1:E:585:LEU:HD21	1:E:847:GLN:HB2	2.02	0.42
1:A:484:THR:HG23	1:A:485:PHE:N	2.35	0.42
2:B:95:LEU:HD21	2:F:112:ILE:CD1	2.43	0.42
3:C:39:VAL:CG1	3:C:55:VAL:HG12	2.45	0.42
1:E:150:CYS:HG	1:E:305:HIS:CG	2.33	0.42
1:A:607:ASN:O	1:A:610:GLU:HG3	2.19	0.42
1:A:163:THR:O	1:A:164:SER:OG	2.16	0.42
1:A:978:THR:HG23	1:A:981:HIS:H	1.85	0.42
1:E:129:VAL:HG23	1:E:129:VAL:O	2.20	0.41
1:A:565:LEU:N	1:A:565:LEU:HD22	2.35	0.41
1:A:744:LEU:HD21	1:A:766:VAL:HG21	2.02	0.41
1:A:821:LEU:HA	1:A:824:LYS:HG2	2.03	0.41
2:B:94:ILE:O	2:B:95:LEU:C	2.57	0.41
2:B:328:LEU:N	2:B:328:LEU:HD12	2.35	0.41
1:E:188:THR:CG2	1:E:229:THR:HG22	2.50	0.41
1:E:928:LEU:HD12	1:E:928:LEU:N	2.35	0.41
1:A:505:ASP:OD2	1:A:510:ARG:NE	2.45	0.41
1:A:741:SER:O	1:A:741:SER:OG	2.30	0.41
1:A:845:VAL:HG23	1:A:848:MET:HB2	2.03	0.41
1:E:484:THR:HG23	1:E:485:PHE:N	2.35	0.41
1:E:626:TYR:CD1	1:E:809:LEU:HD21	2.56	0.41
2:F:169:VAL:O	2:F:169:VAL:CG2	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:328:LEU:HD12	2:F:328:LEU:N	2.35	0.41
2:F:78:THR:HA	2:F:148:GLU:O	2.20	0.41
2:F:101:ASN:OD1	2:F:106:LYS:NZ	2.39	0.41
1:A:285:THR:HG23	1:A:285:THR:O	2.20	0.41
1:A:404:LEU:O	1:A:407:THR:HG22	2.20	0.41
1:A:574:LEU:HB2	1:A:870:ILE:HD11	2.03	0.41
1:A:1035:VAL:HG23	1:A:1036:PHE:HD1	1.84	0.41
1:E:432:SER:N	1:E:435:ASP:OD2	2.54	0.41
1:E:850:GLU:OE2	1:E:854:ARG:NH2	2.54	0.41
1:A:413:THR:HG22	1:A:414:ASP:N	2.36	0.41
1:A:794:ILE:HA	1:A:797:VAL:HG12	2.03	0.41
1:A:936:GLN:OE1	1:A:959:LEU:HD12	2.21	0.41
1:A:994:ILE:HD13	1:A:1035:VAL:HG12	2.02	0.41
2:B:289:GLY:N	3:C:32:GLN:OE1	2.41	0.41
1:E:185:VAL:HG22	1:E:186:ILE:N	2.35	0.41
1:E:328:LEU:O	2:F:336:TYR:N	2.46	0.41
1:E:546:THR:HG23	1:E:547:ILE:N	2.36	0.41
1:A:731:VAL:HG12	2:B:211:VAL:HG21	2.04	0.41
3:C:37:ILE:N	3:C:56:ASN:O	2.49	0.41
2:F:48:GLU:O	2:F:51:GLU:HG3	2.20	0.41
1:A:662:VAL:HG21	1:A:731:VAL:HG11	2.03	0.40
2:F:387:SER:OG	2:F:388:ILE:N	2.54	0.40
1:A:203:TYR:O	1:A:207:GLN:NE2	2.54	0.40
1:A:429:ASP:OD1	1:A:430:CYS:N	2.54	0.40
2:B:684:MET:O	2:B:687:ARG:HG2	2.21	0.40
1:E:315:PRO:O	1:E:316:LEU:HD22	2.20	0.40
1:A:483:GLU:HB3	1:A:520:MET:HE1	2.03	0.40
1:A:902:ALA:O	1:A:906:THR:HG23	2.22	0.40
3:C:46:ASP:OD1	3:C:46:ASP:N	2.55	0.40
1:E:667:ASP:OD1	1:E:667:ASP:N	2.54	0.40
1:A:206:PHE:O	1:A:207:GLN:HB2	2.21	0.40
1:A:590:TYR:CE2	1:A:846:LEU:HD22	2.56	0.40
1:A:924:LEU:N	1:A:924:LEU:HD12	2.37	0.40
2:B:365:VAL:HG12	2:B:366:ASN:N	2.36	0.40
1:E:381:ILE:HG23	1:E:386:PHE:HB2	2.03	0.40
1:E:794:ILE:HA	1:E:797:VAL:HG12	2.04	0.40
3:G:12:LEU:HD21	3:G:82:ILE:HG21	2.04	0.40
3:G:21:VAL:HG11	3:G:26:LEU:HD21	2.02	0.40
3:G:63:VAL:HB	3:G:64:PRO:HD3	2.03	0.40
1:A:914:PHE:CD1	1:A:966:LEU:HD12	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	918/1045 (88%)	849 (92%)	69 (8%)	0	100	100
1	E	918/1045 (88%)	855 (93%)	63 (7%)	0	100	100
2	B	341/618 (55%)	304 (89%)	37 (11%)	0	100	100
2	F	342/618 (55%)	314 (92%)	28 (8%)	0	100	100
3	C	77/83 (93%)	69 (90%)	8 (10%)	0	100	100
3	G	77/83 (93%)	71 (92%)	6 (8%)	0	100	100
All	All	2673/3492 (76%)	2462 (92%)	211 (8%)	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	815/911 (90%)	805 (99%)	10 (1%)	71	83
1	E	815/911 (90%)	799 (98%)	16 (2%)	55	73
2	B	319/545 (58%)	316 (99%)	3 (1%)	78	88
2	F	320/545 (59%)	312 (98%)	8 (2%)	47	68
3	C	70/73 (96%)	67 (96%)	3 (4%)	29	56
3	G	70/73 (96%)	67 (96%)	3 (4%)	29	56
All	All	2409/3058 (79%)	2366 (98%)	43 (2%)	61	77

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	116	LEU
1	A	143	PHE
1	A	266	TRP
1	A	387	GLN
1	A	643	TYR
1	A	722	GLU
1	A	809	LEU
1	A	821	LEU
1	A	841	LYS
1	A	1040	LEU
2	B	102	ASN
2	B	187	LEU
2	B	692	LEU
3	C	18	GLU
3	C	81	LYS
3	C	85	ARG
1	E	116	LEU
1	E	205	GLU
1	E	297	CYS
1	E	298	HIS
1	E	311	TYR
1	E	387	GLN
1	E	643	TYR
1	E	693	LEU
1	E	732	LEU
1	E	802	HIS
1	E	807	HIS
1	E	810	HIS
1	E	841	LYS
1	E	921	MET
1	E	924	LEU
1	E	944	LYS
2	F	54	ARG
2	F	126	GLN
2	F	185	GLN
2	F	187	LEU
2	F	188	LEU
2	F	283	PHE
2	F	303	ASP
2	F	692	LEU
3	G	50	LYS
3	G	52	PHE

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Mol	Chain	Res	Type
3	G	85	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	F	271	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	D	9/46 (19%)	3 (33%)	0
4	H	9/46 (19%)	6 (66%)	0
All	All	18/92 (19%)	9 (50%)	0

All (9) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	D	27	U
4	D	28	U
4	D	41	A
4	H	27	U
4	H	28	U
4	H	40	A
4	H	41	A
4	H	42	A
4	H	46	A

There are no RNA pucker outliers to report.

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

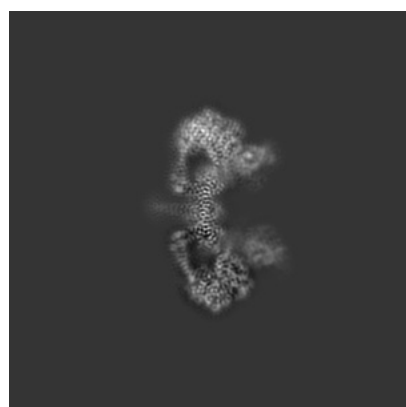
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24882. These allow visual inspection of the internal detail of the map and identification of artifacts.

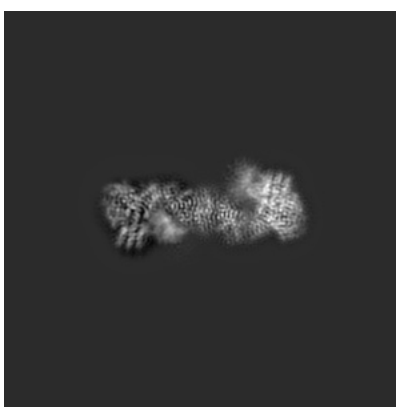
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

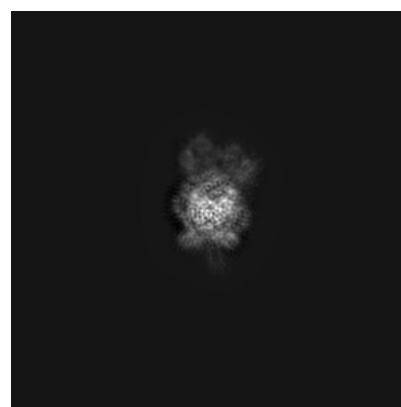
6.1.1 Primary map



X



Y

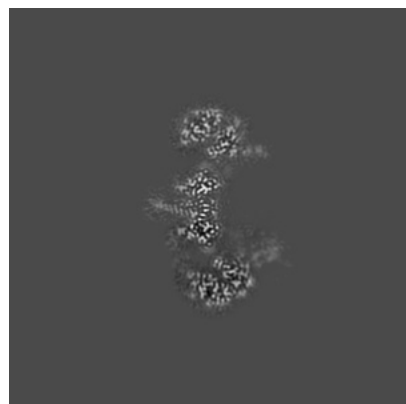


Z

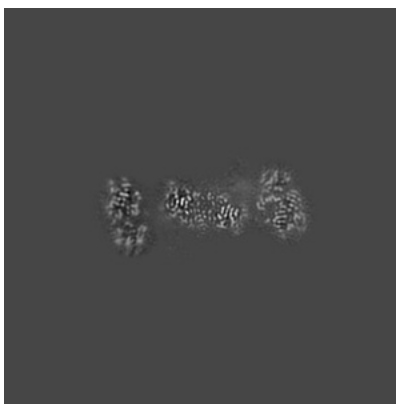
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 192



Y Index: 192

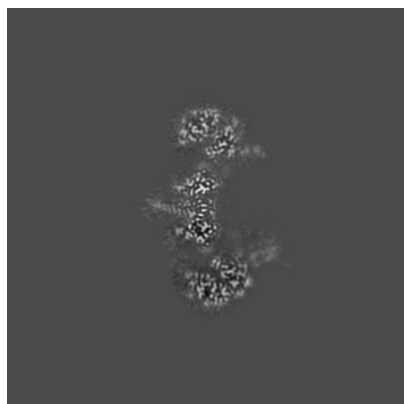


Z Index: 192

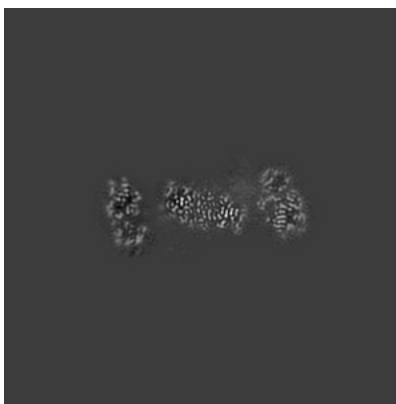
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

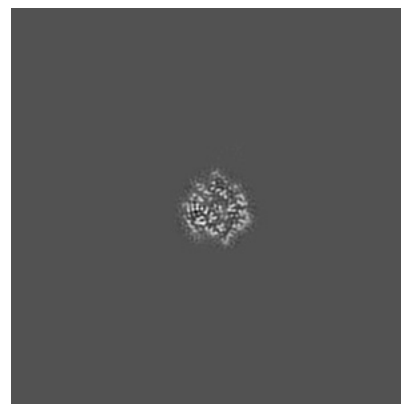
6.3.1 Primary map



X Index: 192



Y Index: 191

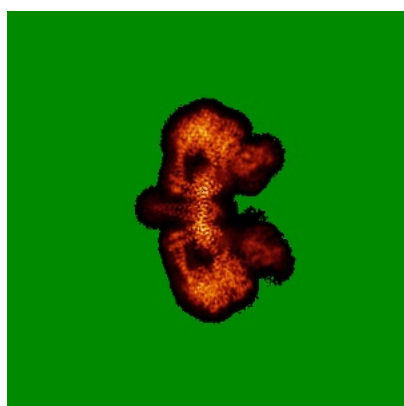


Z Index: 264

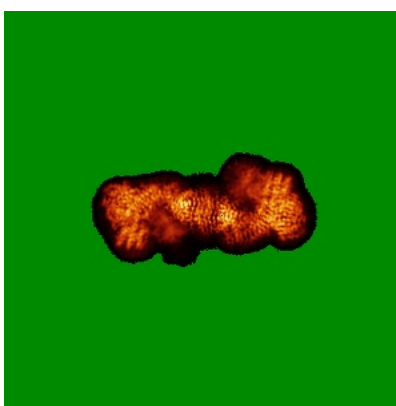
The images above show the largest variance slices of the map in three orthogonal directions.

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

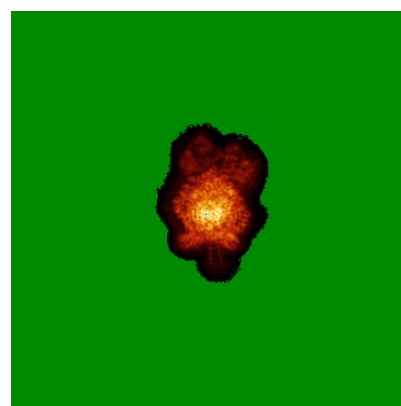
6.4.1 Primary map



X



Y

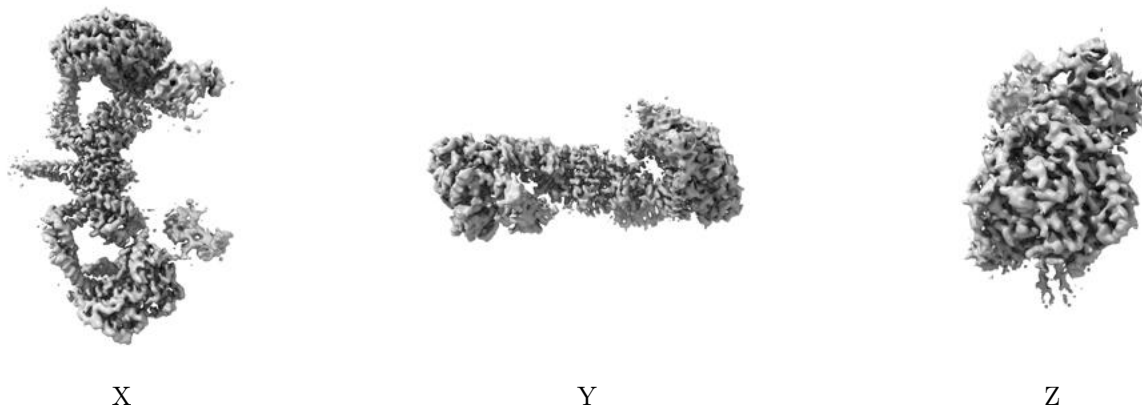


Z

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

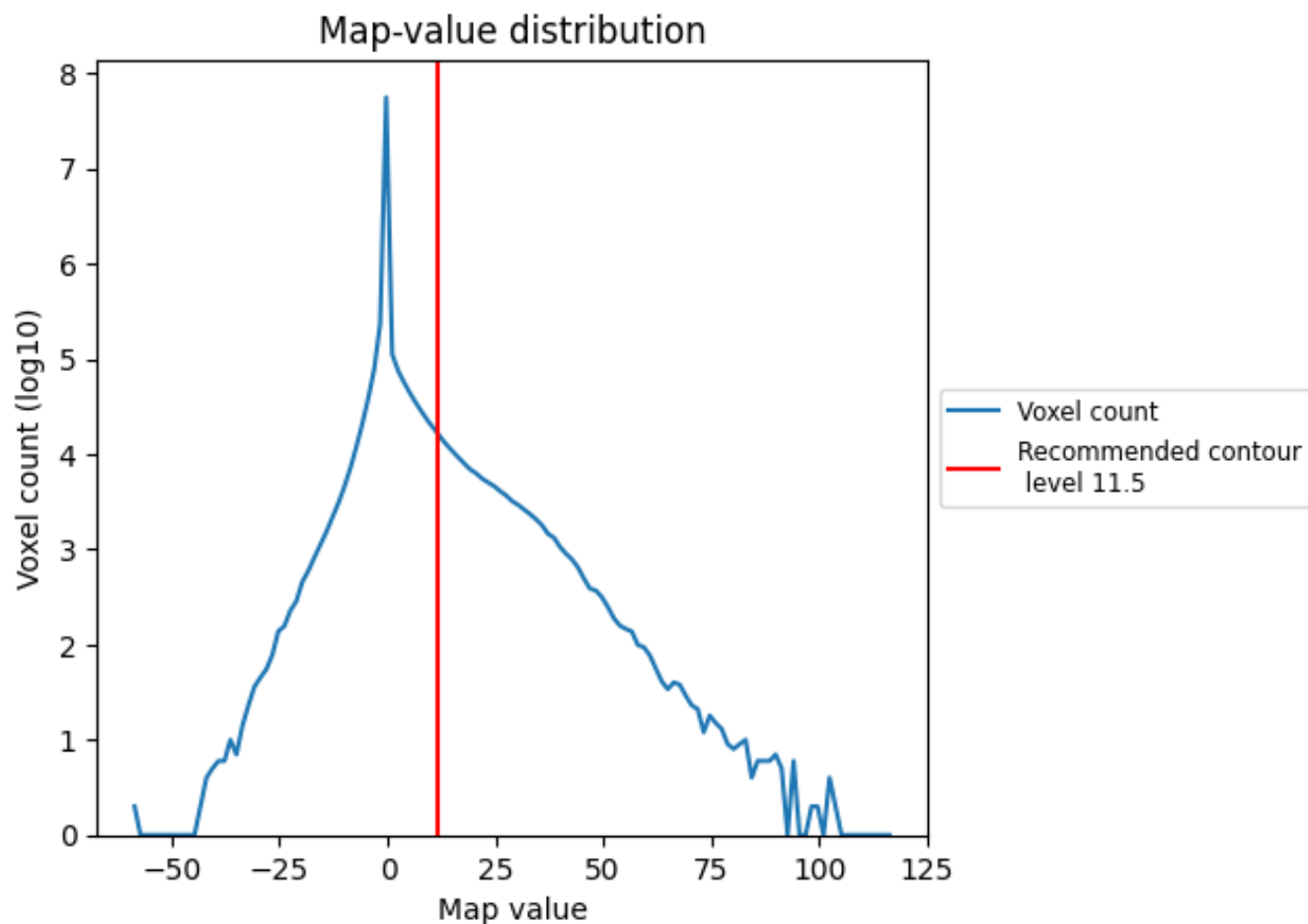
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

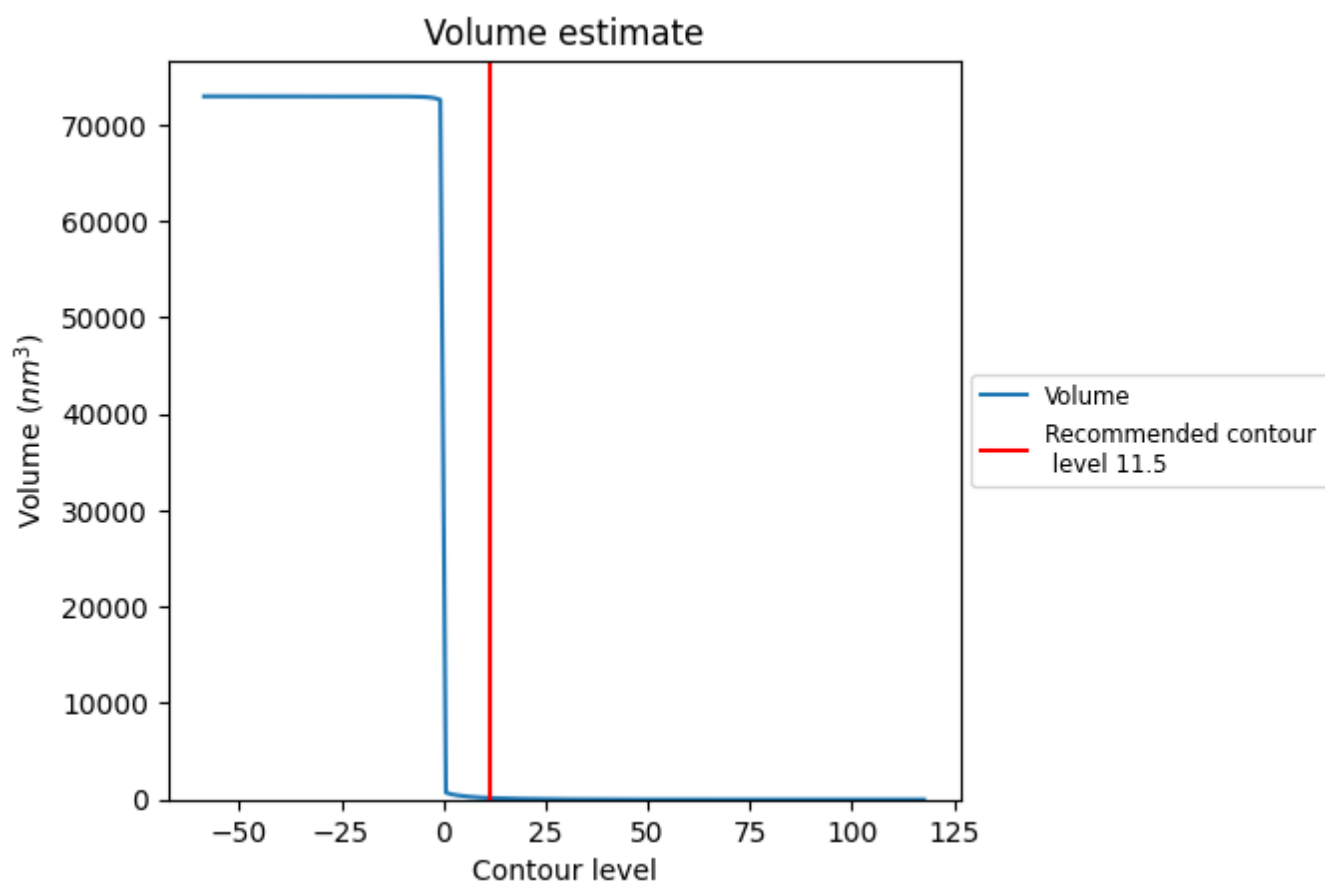
This section contains the results of statistical analysis of the map.

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.

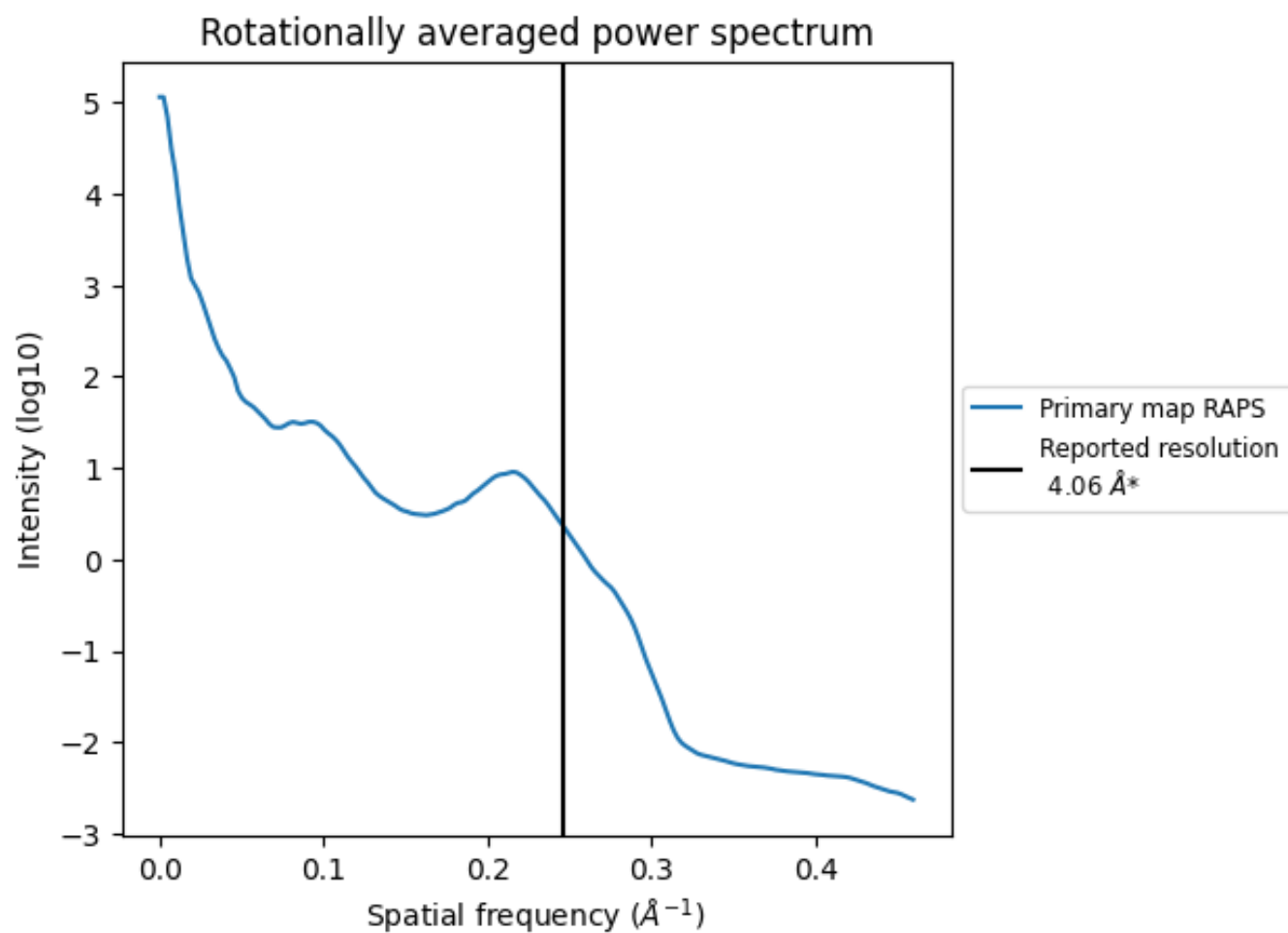
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 161 nm³; this corresponds to an approximate mass of 146 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 0.246 Å⁻¹

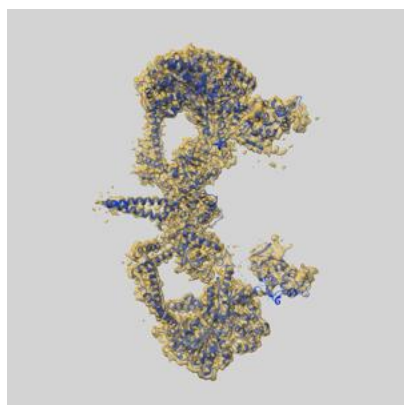
8 Fourier-Shell correlation

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

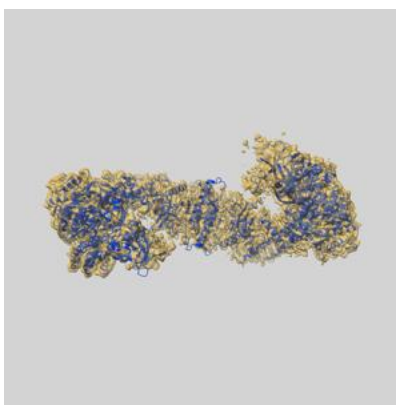
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-24882 and PDB model 7S7B. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

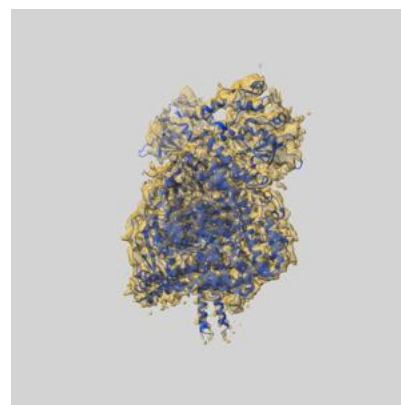
9.1 Map-model overlay [i](#)



X



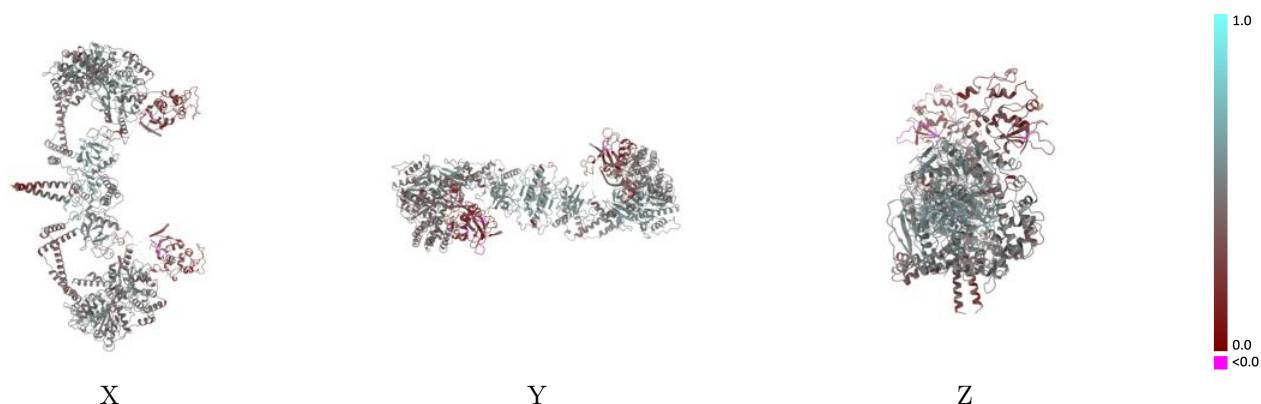
Y



Z

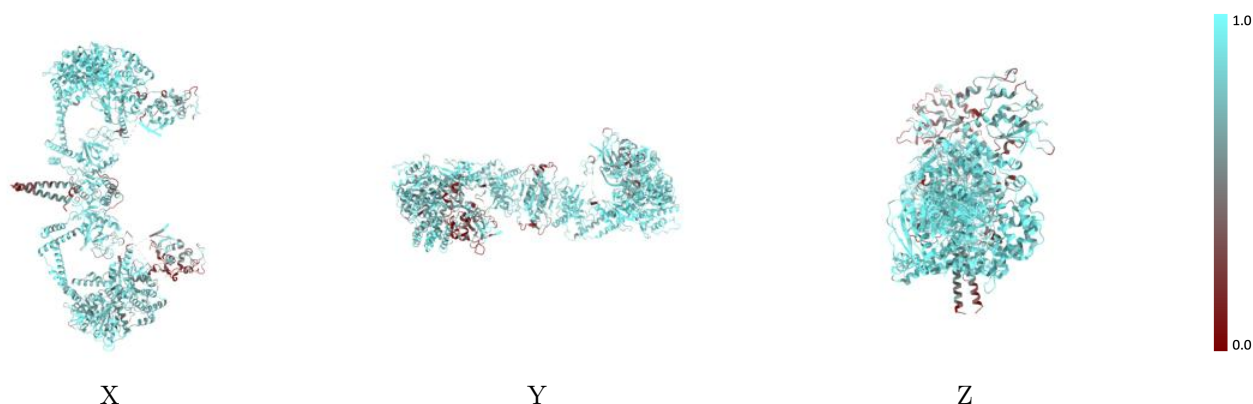
The images above show the 3D surface view of the map at the recommended contour level 11.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



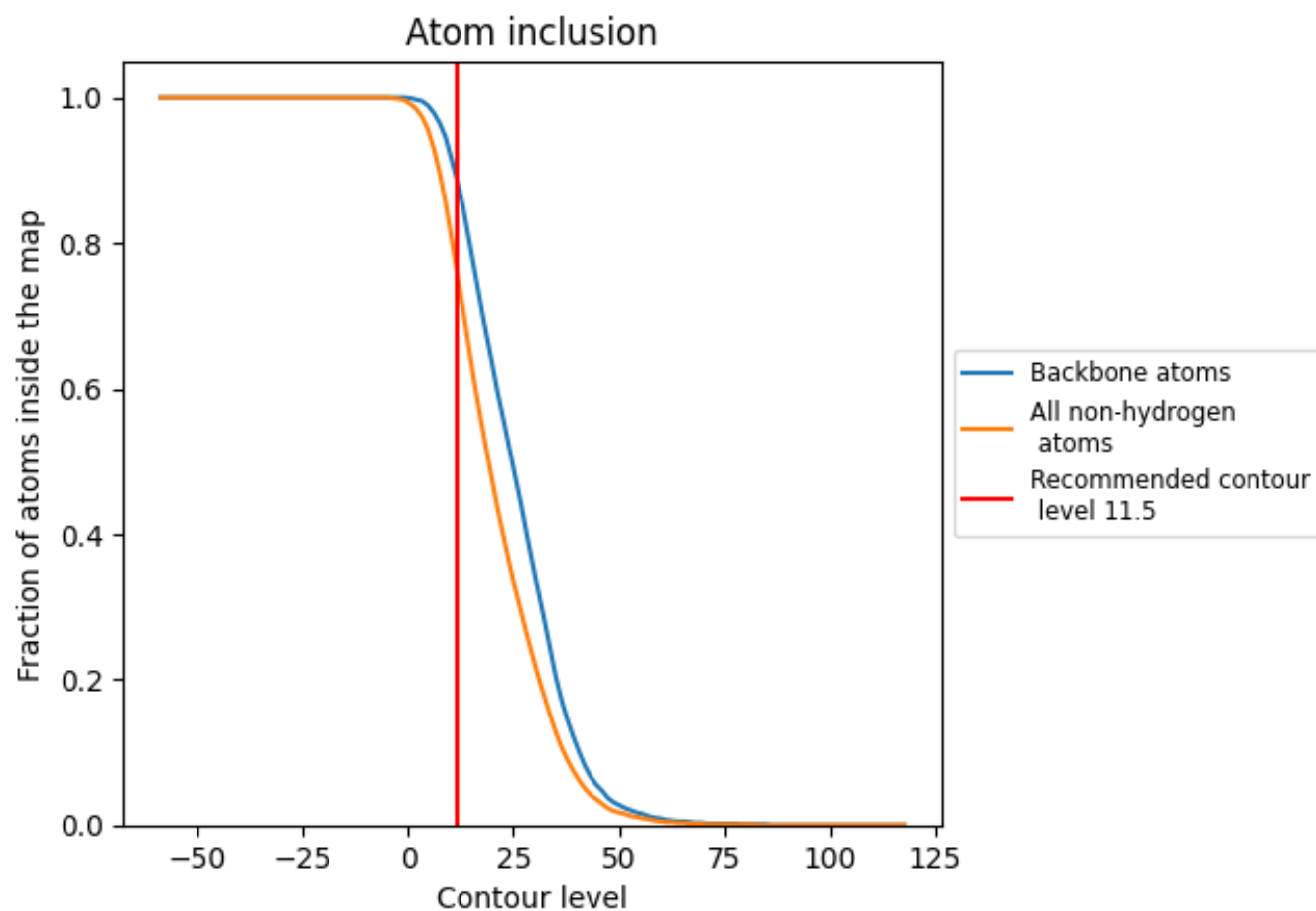
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (11.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 89% of all backbone atoms, 77% 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 (11.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.7660	<div></div> 0.4440
A	<div></div> 0.8920	<div></div> 0.4940
B	<div></div> 0.6730	<div></div> 0.4260
C	<div></div> 0.7410	<div></div> 0.2710
D	<div></div> 0.9150	<div></div> 0.4600
E	<div></div> 0.7910	<div></div> 0.4620
F	<div></div> 0.5500	<div></div> 0.3880
G	<div></div> 0.4320	<div></div> 0.1800
H	<div></div> 0.6610	<div></div> 0.3910

1.0

0.0

<0.0