Full wwPDB EM Validation Report ⓘ

Oct 14, 2024 – 07:42 AM EDT

PDB ID	:	7L7K
EMDB ID	:	EMD-23215
Title	:	Cryo-EM structure of protein encoded by vaccine candidate BNT162b2
Authors	:	Lees, J.A.; Han, S.
Deposited on	:	2020-12-28
Resolution	:	3.29 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:

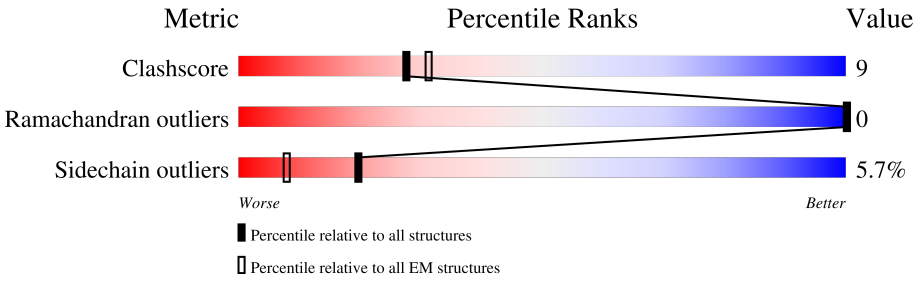
EMDB validation analysis	:	0.0.1.dev113
MolProbity	:	4.02b-467
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

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 3.29 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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

Mol	Chain	Length	Quality of chain
1	A	1273	57% 19% 24%
1	B	1273	58% 18% 24%
1	C	1273	58% 17% 24%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 22124 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 Spike glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	973	Total	C	N	O	S	0	0
			7370	4708	1225	1403	34		
1	C	973	Total	C	N	O	S	0	0
			7384	4720	1225	1405	34		
1	A	973	Total	C	N	O	S	0	0
			7370	4708	1225	1403	34		

There are 6 discrepancies between the modelled and reference sequences:

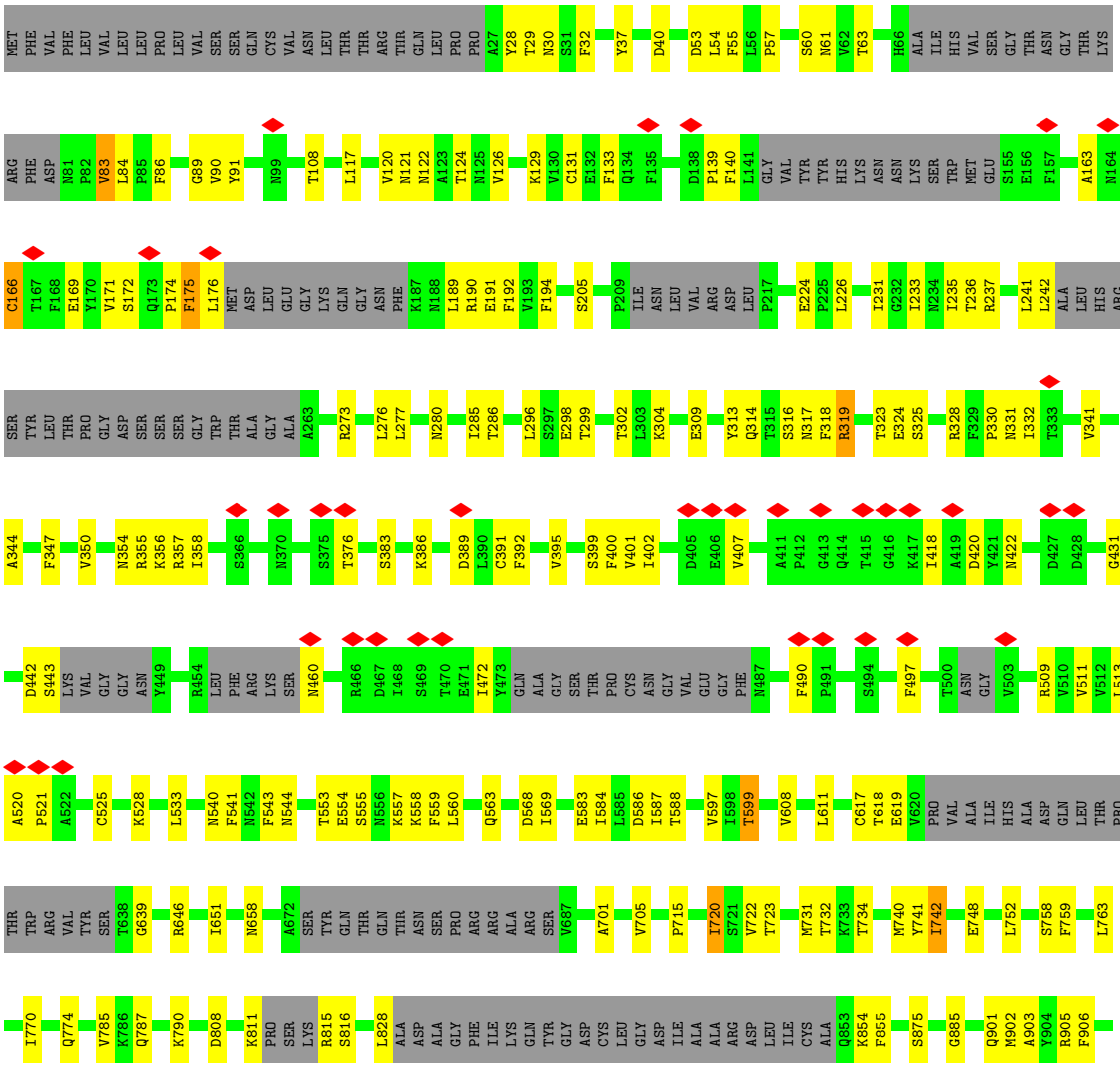
Chain	Residue	Modelled	Actual	Comment	Reference
B	986	PRO	LYS	conflict	UNP P0DTC2
B	987	PRO	VAL	conflict	UNP P0DTC2
C	986	PRO	LYS	conflict	UNP P0DTC2
C	987	PRO	VAL	conflict	UNP P0DTC2
A	986	PRO	LYS	conflict	UNP P0DTC2
A	987	PRO	VAL	conflict	UNP P0DTC2

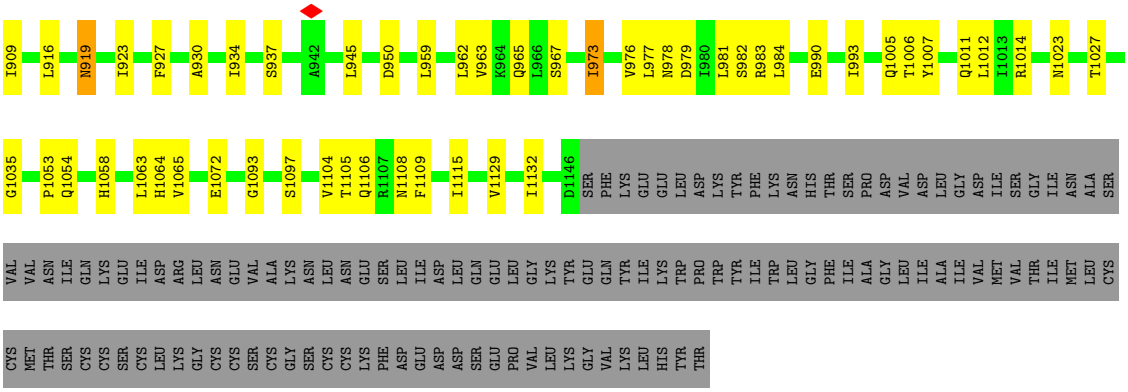
3 Residue-property plots (i)

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: Spike glycoprotein

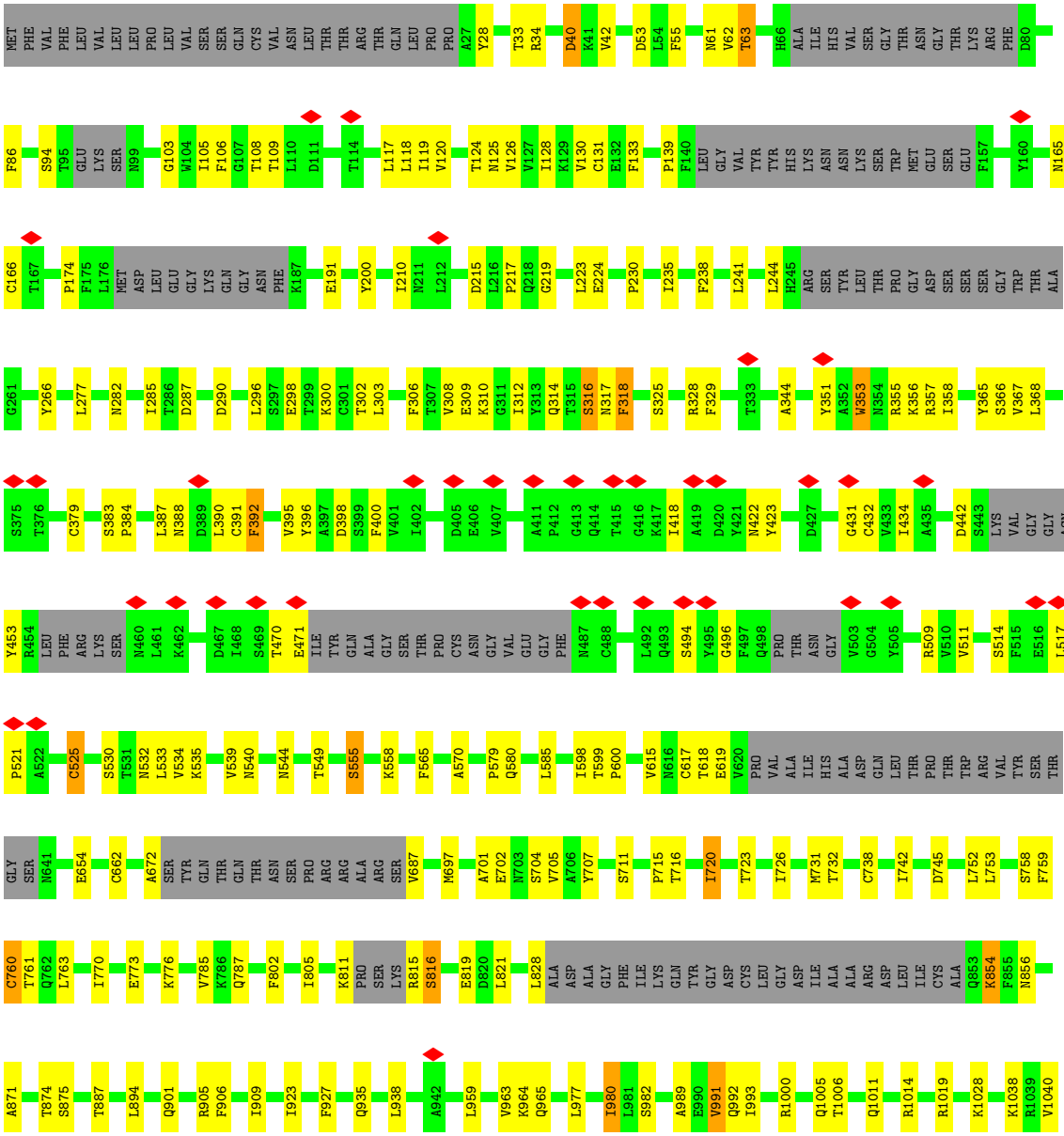
Chain B:

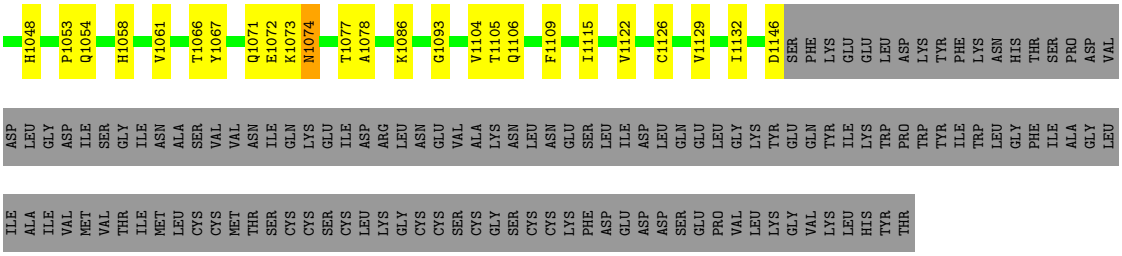




• Molecule 1: Spike glycoprotein

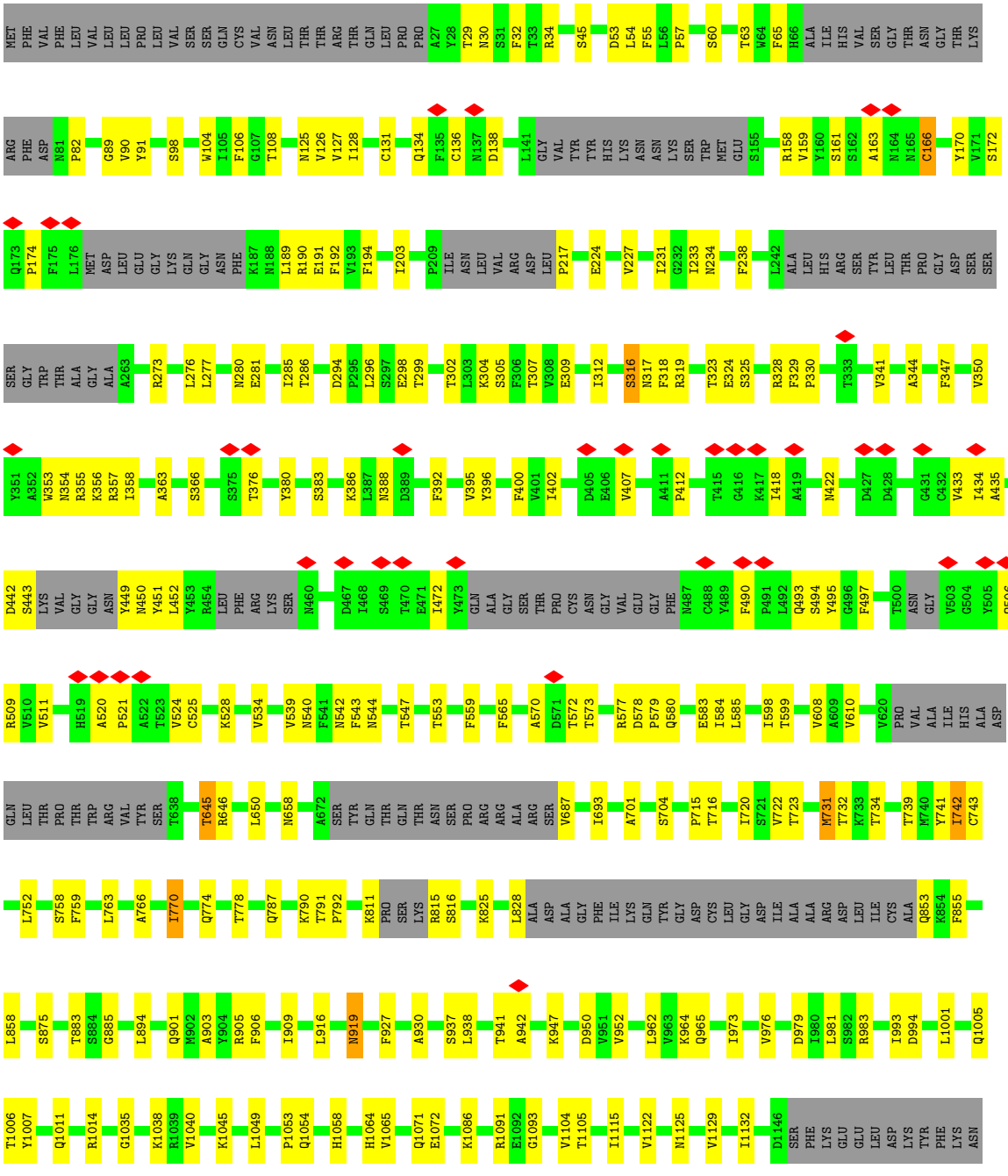
Chain C:





• Molecule 1: Spike glycoprotein

Chain A:



HIS THR SER PRO ASP VAL ASP LEU GLY ASP ILE SER GLY ILE ASN ALA SER VAL VAL ASN ILE GLN LYS GLU ILE ASP ARG LEU ASN GLU VAL ALA LYS ASN LEU ASN GLU SER LEU ILE ASP LEU GLN GLU LEU GLY LYS TYR GLU GLN TYR ILE LYS TRP PRO TRP TYR ILE TRP LEU

GLY PHE ILE ALA GLY LEU ILE ALA ILE VAL MET VAL THR ILE MET LEU CYS CYS MET THR SER CYS CYS SER CYS LEU LYS GLY CYS CYS SER CYS GLY SER CYS CYS LYS PHE ASP GLU ASP ASP SER GLU PRO VAL LEU LYS GLY VAL LYS LEU HIS TYR THR

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	58295	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 ($e^-/Å^2$)	50.22	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	0.002	Depositor
Minimum map value	-0.001	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.000	Depositor
Recommended contour level	0.0003	Depositor
Map size (Å)	378.0, 378.0, 378.0	wwPDB
Map dimensions	450, 450, 450	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.84, 0.84, 0.84	Depositor

5 Model quality

5.1 Standard geometry

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	0.27	0/7527	0.46	0/10258
1	B	0.28	0/7527	0.47	0/10258
1	C	0.27	0/7540	0.46	0/10279
All	All	0.27	0/22594	0.46	0/30795

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

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	7370	0	7024	155	0
1	B	7370	0	7024	138	0
1	C	7384	0	7061	143	0
All	All	22124	0	21109	409	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (409) 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:C:103:GLY:HA3	1:C:119:ILE:O	1.53	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:391:CYS:HA	1:B:525:CYS:HB3	1.53	0.91
1:C:105:ILE:HG12	1:C:241:LEU:HD11	1.58	0.85
1:C:391:CYS:HA	1:C:525:CYS:HB3	1.59	0.85
1:A:347:PHE:HD2	1:A:509:ARG:HD3	1.48	0.78
1:B:646:ARG:O	1:B:646:ARG:NH1	2.21	0.73
1:A:126:VAL:HG23	1:A:174:PRO:HA	1.70	0.72
1:A:646:ARG:O	1:A:646:ARG:NH1	2.22	0.71
1:A:131:CYS:HA	1:A:166:CYS:HB3	1.76	0.68
1:A:1093:GLY:HA3	1:A:1105:THR:O	1.92	0.67
1:A:277:LEU:HD23	1:A:285:ILE:HG21	1.76	0.67
1:A:442:ASP:OD1	1:A:509:ARG:NH1	2.28	0.67
1:C:384:PRO:HA	1:C:387:LEU:HD13	1.75	0.67
1:B:418:ILE:HA	1:B:422:ASN:HD22	1.60	0.67
1:A:901:GLN:HE21	1:A:905:ARG:HH21	1.43	0.66
1:A:578:ASP:HB2	1:A:583:GLU:H	1.61	0.66
1:B:296:LEU:HD13	1:B:608:VAL:HG21	1.77	0.66
1:C:367:VAL:HG23	1:C:368:LEU:HD12	1.78	0.66
1:C:28:TYR:HA	1:C:62:VAL:O	1.95	0.65
1:C:418:ILE:HA	1:C:422:ASN:HD22	1.59	0.65
1:A:1011:GLN:OE1	1:A:1014:ARG:NH1	2.29	0.65
1:A:53:ASP:OD1	1:A:54:LEU:N	2.30	0.64
1:C:344:ALA:O	1:C:509:ARG:NH1	2.30	0.64
1:B:977:LEU:HD11	1:B:993:ILE:HG12	1.80	0.64
1:C:296:LEU:HG	1:C:300:LYS:HE3	1.80	0.64
1:B:981:LEU:HD21	1:B:993:ILE:HD11	1.79	0.63
1:C:392:PHE:HD2	1:C:517:LEU:HD11	1.63	0.63
1:C:1106:GLN:HE21	1:C:1109:PHE:HB3	1.63	0.63
1:B:139:PRO:HG2	1:B:241:LEU:HD23	1.81	0.63
1:B:901:GLN:HE21	1:B:905:ARG:HH21	1.47	0.63
1:B:401:VAL:HG12	1:B:509:ARG:HG3	1.80	0.62
1:C:1006:THR:OG1	1:A:1005:GLN:OE1	2.17	0.62
1:C:1129:VAL:HG13	1:C:1132:ILE:HD13	1.81	0.62
1:A:134:GLN:H	1:A:161:SER:HB2	1.62	0.62
1:C:126:VAL:HG23	1:C:174:PRO:HA	1.80	0.62
1:C:1011:GLN:OE1	1:C:1014:ARG:NH1	2.33	0.62
1:A:366:SER:HB3	1:A:388:ASN:HD21	1.63	0.62
1:B:1005:GLN:OE1	1:A:1006:THR:OG1	2.18	0.62
1:C:901:GLN:HE21	1:C:905:ARG:HH21	1.46	0.62
1:A:34:ARG:NH1	1:A:217:PRO:O	2.33	0.62
1:B:790:LYS:HE3	1:A:704:SER:HB3	1.82	0.62
1:B:442:ASP:OD1	1:B:509:ARG:NH1	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:347:PHE:HD2	1:B:509:ARG:HD3	1.66	0.61
1:B:358:ILE:HB	1:B:395:VAL:HB	1.82	0.61
1:B:978:ASN:OD1	1:A:547:THR:OG1	2.19	0.61
1:A:316:SER:OG	1:A:317:ASN:N	2.33	0.61
1:B:420:ASP:O	1:B:460:ASN:N	2.34	0.61
1:C:328:ARG:HH11	1:C:530:SER:HB3	1.66	0.61
1:A:280:ASN:ND2	1:A:286:THR:OG1	2.34	0.61
1:C:701:ALA:HB3	1:A:787:GLN:HG3	1.81	0.61
1:A:131:CYS:HB3	1:A:163:ALA:HA	1.82	0.61
1:A:402:ILE:HD11	1:A:407:VAL:HA	1.82	0.61
1:C:119:ILE:HG13	1:C:128:ILE:HD12	1.83	0.60
1:B:57:PRO:HG3	1:B:273:ARG:HD2	1.83	0.60
1:C:434:ILE:HB	1:C:511:VAL:HG23	1.83	0.60
1:B:1006:THR:OG1	1:C:1005:GLN:OE1	2.18	0.60
1:B:108:THR:HG23	1:B:236:THR:HB	1.83	0.60
1:B:1011:GLN:OE1	1:B:1014:ARG:NH1	2.35	0.60
1:C:358:ILE:HB	1:C:395:VAL:HB	1.84	0.60
1:C:131:CYS:HA	1:C:166:CYS:HB3	1.84	0.59
1:A:363:ALA:N	1:A:525:CYS:O	2.34	0.59
1:A:434:ILE:HB	1:A:511:VAL:HG23	1.83	0.59
1:B:280:ASN:ND2	1:B:286:THR:OG1	2.35	0.59
1:A:296:LEU:HD13	1:A:608:VAL:HG21	1.84	0.59
1:C:672:ALA:O	1:C:687:VAL:N	2.36	0.59
1:C:977:LEU:HD11	1:C:993:ILE:HG12	1.85	0.59
1:A:358:ILE:HB	1:A:395:VAL:HB	1.83	0.59
1:B:722:VAL:HG22	1:B:930:ALA:HB1	1.85	0.59
1:A:722:VAL:HG22	1:A:930:ALA:HB1	1.85	0.58
1:C:1053:PRO:O	1:C:1054:GLN:NE2	2.36	0.58
1:B:316:SER:OG	1:B:317:ASN:N	2.35	0.58
1:A:1104:VAL:HG23	1:A:1115:ILE:HG12	1.84	0.58
1:C:540:ASN:HB3	1:C:549:THR:HG22	1.85	0.58
1:C:244:LEU:HD23	1:C:244:LEU:H	1.69	0.58
1:A:1129:VAL:HB	1:A:1132:ILE:HD11	1.84	0.57
1:B:126:VAL:HB	1:B:174:PRO:HB3	1.86	0.57
1:B:787:GLN:HG3	1:A:701:ALA:HB3	1.86	0.57
1:C:442:ASP:OD2	1:C:509:ARG:NE	2.30	0.57
1:B:701:ALA:HB3	1:C:787:GLN:HG2	1.86	0.57
1:A:298:GLU:O	1:A:302:THR:OG1	2.23	0.57
1:B:402:ILE:HD11	1:B:407:VAL:HA	1.86	0.57
1:B:233:ILE:HG22	1:B:235:ILE:HG13	1.85	0.56
1:C:390:LEU:HD21	1:A:983:ARG:HB2	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:VAL:HG13	1:B:356:LYS:HD2	1.86	0.56
1:A:731:MET:HG3	1:A:774:GLN:HG3	1.87	0.56
1:B:319:ARG:NH2	1:C:745:ASP:OD1	2.39	0.56
1:A:347:PHE:CD2	1:A:509:ARG:HD3	2.36	0.56
1:B:328:ARG:HH21	1:B:533:LEU:HB2	1.71	0.56
1:B:560:LEU:H	1:B:563:GLN:HG3	1.70	0.56
1:A:325:SER:HA	1:A:540:ASN:O	2.05	0.56
1:A:543:PHE:CZ	1:A:579:PRO:HD3	2.41	0.56
1:C:760:CYS:SG	1:C:761:THR:N	2.79	0.56
1:B:1104:VAL:HG23	1:B:1115:ILE:HG12	1.88	0.55
1:C:618:THR:OG1	1:C:619:GLU:OE1	2.24	0.55
1:A:57:PRO:HG3	1:A:273:ARG:HD2	1.87	0.55
1:B:741:TYR:HD2	1:B:742:ILE:HG22	1.72	0.55
1:A:449:TYR:O	1:A:494:SER:OG	2.25	0.55
1:A:828:LEU:H	1:A:828:LEU:HD12	1.72	0.55
1:C:1093:GLY:HA3	1:C:1105:THR:O	2.06	0.55
1:B:722:VAL:HG12	1:B:1065:VAL:HG22	1.87	0.55
1:A:439:ASN:HA	1:A:442:ASP:HB2	1.87	0.55
1:B:1093:GLY:HA3	1:B:1105:THR:O	2.07	0.54
1:B:731:MET:HG3	1:B:774:GLN:HG3	1.89	0.54
1:A:534:VAL:HG21	1:A:539:VAL:HG11	1.88	0.54
1:C:329:PHE:O	1:C:580:GLN:NE2	2.32	0.54
1:B:91:TYR:OH	1:B:191:GLU:OE1	2.26	0.54
1:B:108:THR:O	1:B:237:ARG:NH1	2.38	0.54
1:A:108:THR:OG1	1:A:234:ASN:O	2.26	0.54
1:A:323:THR:OG1	1:A:324:GLU:N	2.41	0.54
1:A:722:VAL:HG12	1:A:1065:VAL:HG22	1.90	0.54
1:B:347:PHE:CD2	1:B:509:ARG:HD3	2.42	0.54
1:C:887:THR:HG21	1:C:894:LEU:HG	1.89	0.54
1:B:758:SER:H	1:A:965:GLN:HE22	1.56	0.54
1:C:210:ILE:HG21	1:C:217:PRO:HG3	1.90	0.54
1:B:277:LEU:HD23	1:B:285:ILE:HG21	1.89	0.53
1:B:740:MET:HG3	1:A:319:ARG:HH22	1.73	0.53
1:A:309:GLU:OE2	1:A:309:GLU:N	2.41	0.53
1:B:472:ILE:HA	1:B:490:PHE:HA	1.90	0.53
1:B:553:THR:HG22	1:B:554:GLU:H	1.74	0.53
1:C:34:ARG:NE	1:C:191:GLU:OE1	2.41	0.53
1:C:312:ILE:HD12	1:C:598:ILE:HG12	1.89	0.53
1:C:707:TYR:HB2	1:A:883:THR:HG23	1.90	0.53
1:A:559:PHE:HB2	1:A:577:ARG:HH21	1.73	0.53
1:C:316:SER:OG	1:C:317:ASN:N	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:365:TYR:HD2	1:C:387:LEU:HB3	1.74	0.53
1:A:45:SER:HB3	1:A:281:GLU:HA	1.91	0.53
1:A:125:ASN:OD1	1:A:125:ASN:N	2.42	0.53
1:B:205:SER:HB3	1:B:226:LEU:HD13	1.89	0.52
1:B:89:GLY:HA2	1:B:194:PHE:O	2.10	0.52
1:B:828:LEU:HD12	1:B:828:LEU:H	1.73	0.52
1:A:443:SER:HA	1:A:497:PHE:HB2	1.91	0.52
1:B:174:PRO:HG2	1:B:176:LEU:HG	1.90	0.52
1:A:715:PRO:HA	1:A:1072:GLU:HA	1.91	0.52
1:C:773:GLU:OE2	1:C:1019:ARG:NH1	2.43	0.52
1:B:175:PHE:O	1:B:190:ARG:NH2	2.43	0.51
1:C:1104:VAL:HG23	1:C:1115:ILE:HG12	1.91	0.51
1:A:578:ASP:N	1:A:583:GLU:O	2.41	0.51
1:C:366:SER:OG	1:C:388:ASN:ND2	2.41	0.51
1:B:298:GLU:O	1:B:302:THR:OG1	2.26	0.51
1:B:965:GLN:HE22	1:C:758:SER:H	1.56	0.51
1:A:231:ILE:HG22	1:A:233:ILE:HG12	1.91	0.51
1:A:741:TYR:HD2	1:A:742:ILE:HG22	1.76	0.51
1:C:351:TYR:HB3	1:C:453:TYR:HA	1.91	0.51
1:B:53:ASP:OD2	1:B:54:LEU:N	2.44	0.51
1:B:618:THR:OG1	1:B:619:GLU:OE2	2.20	0.51
1:C:139:PRO:HB3	1:C:241:LEU:HD23	1.91	0.51
1:B:29:THR:HG22	1:B:30:ASN:H	1.75	0.51
1:B:331:ASN:OD1	1:B:332:ILE:N	2.44	0.51
1:C:1048:HIS:HA	1:C:1066:THR:HG22	1.93	0.51
1:C:742:ILE:HG21	1:C:753:LEU:HD13	1.94	0.50
1:A:190:ARG:HB3	1:A:192:PHE:HE2	1.76	0.50
1:C:131:CYS:HB2	1:C:133:PHE:CE1	2.47	0.50
1:C:353:TRP:HB3	1:C:400:PHE:HB3	1.94	0.50
1:B:599:THR:HB	1:B:608:VAL:HG22	1.94	0.50
1:C:715:PRO:HA	1:C:1072:GLU:HA	1.94	0.49
1:B:885:GLY:HA2	1:B:901:GLN:NE2	2.26	0.49
1:B:53:ASP:HB3	1:B:55:PHE:CE2	2.47	0.49
1:C:517:LEU:HB3	1:C:518:LEU:HD23	1.93	0.49
1:C:418:ILE:HG23	1:C:422:ASN:HB2	1.94	0.49
1:B:325:SER:HA	1:B:540:ASN:O	2.13	0.49
1:C:828:LEU:HD12	1:C:828:LEU:H	1.77	0.49
1:A:53:ASP:HB3	1:A:55:PHE:CE2	2.48	0.49
1:B:299:THR:HG22	1:B:597:VAL:HG21	1.94	0.49
1:B:770:ILE:HD11	1:B:1012:LEU:HD23	1.94	0.49
1:C:353:TRP:NE1	1:C:422:ASN:O	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:VAL:HG22	1:A:356:LYS:HZ2	1.78	0.49
1:A:741:TYR:CD2	1:A:742:ILE:HG22	2.48	0.49
1:C:532:ASN:OD1	1:C:532:ASN:N	2.46	0.49
1:A:29:THR:HG22	1:A:30:ASN:H	1.77	0.49
1:A:312:ILE:HD12	1:A:598:ILE:HG12	1.95	0.49
1:C:555:SER:O	1:C:555:SER:OG	2.30	0.49
1:C:40:ASP:HB3	1:C:42:VAL:HG13	1.95	0.48
1:B:439:ASN:HA	1:B:442:ASP:HB2	1.95	0.48
1:C:33:THR:OG1	1:C:219:GLY:O	2.17	0.48
1:C:398:ASP:O	1:C:511:VAL:HA	2.13	0.48
1:A:1053:PRO:O	1:A:1054:GLN:NE2	2.46	0.48
1:B:555:SER:HB2	1:B:557:LYS:HG2	1.94	0.48
1:A:350:VAL:HG22	1:A:402:ILE:HG22	1.94	0.48
1:B:166:CYS:HB2	1:B:169:GLU:OE2	2.14	0.48
1:A:128:ILE:HG23	1:A:170:TYR:HB3	1.96	0.48
1:A:328:ARG:NH2	1:A:578:ASP:OD2	2.46	0.48
1:C:310:LYS:HG3	1:C:600:PRO:HA	1.95	0.48
1:C:452:LEU:HA	1:C:494:SER:HA	1.95	0.48
1:C:821:LEU:HD23	1:C:938:LEU:HB3	1.96	0.48
1:A:203:ILE:HB	1:A:227:VAL:HG22	1.95	0.48
1:A:366:SER:HB3	1:A:388:ASN:ND2	2.29	0.48
1:B:959:LEU:O	1:B:963:VAL:HG23	2.14	0.48
1:C:325:SER:HB3	1:C:540:ASN:ND2	2.28	0.48
1:C:871:ALA:HA	1:C:874:THR:HG22	1.95	0.48
1:A:981:LEU:HD21	1:A:993:ILE:HD11	1.96	0.48
1:A:277:LEU:HD23	1:A:285:ILE:HD13	1.96	0.47
1:A:687:VAL:N	1:A:693:ILE:H	2.12	0.47
1:A:716:THR:OG1	1:A:1071:GLN:O	2.27	0.47
1:A:1093:GLY:CA	1:A:1105:THR:O	2.59	0.47
1:B:309:GLU:N	1:B:309:GLU:OE1	2.46	0.47
1:C:738:CYS:HB3	1:C:760:CYS:HB2	1.95	0.47
1:A:330:PRO:HD3	1:A:544:ASN:ND2	2.29	0.47
1:C:742:ILE:O	1:C:1000:ARG:NH1	2.42	0.47
1:A:170:TYR:CE1	1:A:172:SER:HB2	2.50	0.47
1:B:354:ASN:OD1	1:B:355:ARG:N	2.47	0.47
1:B:741:TYR:CD2	1:B:742:ILE:HG22	2.49	0.47
1:C:224:GLU:N	1:C:224:GLU:OE1	2.44	0.47
1:C:619:GLU:OE1	1:C:619:GLU:N	2.48	0.47
1:C:707:TYR:HB3	1:A:792:PRO:HG3	1.96	0.47
1:A:276:LEU:HD11	1:A:304:LYS:HA	1.97	0.47
1:C:379:CYS:HA	1:C:432:CYS:HA	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:534:VAL:HG21	1:C:539:VAL:HG21	1.96	0.47
1:B:1035:GLY:HA3	1:A:1040:VAL:HG21	1.96	0.47
1:A:1086:LYS:HD2	1:A:1122:VAL:HG21	1.95	0.47
1:A:885:GLY:HA2	1:A:901:GLN:NE2	2.29	0.47
1:C:53:ASP:HB2	1:C:55:PHE:CE2	2.50	0.47
1:A:354:ASN:OD1	1:A:355:ARG:N	2.48	0.47
1:B:299:THR:CG2	1:B:597:VAL:HG21	2.45	0.46
1:C:418:ILE:HD13	1:C:422:ASN:ND2	2.30	0.46
1:A:720:ILE:HD11	1:A:927:PHE:HB2	1.96	0.46
1:B:559:PHE:HB3	1:B:563:GLN:HB2	1.97	0.46
1:B:934:ILE:HD13	1:B:1063:LEU:HD22	1.97	0.46
1:C:705:VAL:HB	1:A:883:THR:HG21	1.96	0.46
1:B:715:PRO:HA	1:B:1072:GLU:HA	1.97	0.46
1:B:740:MET:HG3	1:A:319:ARG:NH2	2.31	0.46
1:C:533:LEU:HD13	1:C:535:LYS:HZ3	1.79	0.46
1:B:83:VAL:O	1:B:84:LEU:HD13	2.15	0.46
1:A:418:ILE:HA	1:A:422:ASN:HD22	1.79	0.46
1:C:716:THR:OG1	1:C:1071:GLN:O	2.25	0.46
1:B:330:PRO:HD3	1:B:544:ASN:ND2	2.31	0.46
1:C:277:LEU:HD22	1:C:285:ILE:HD13	1.97	0.46
1:C:811:LYS:O	1:C:815:ARG:N	2.49	0.46
1:A:383:SER:HB3	1:A:386:LYS:HD2	1.98	0.46
1:B:224:GLU:OE1	1:B:224:GLU:N	2.46	0.46
1:B:984:LEU:O	1:A:386:LYS:NZ	2.37	0.46
1:C:309:GLU:OE1	1:C:309:GLU:N	2.47	0.46
1:A:811:LYS:HD2	1:A:815:ARG:HB2	1.98	0.46
1:B:323:THR:C	1:B:324:GLU:HG3	2.36	0.45
1:C:379:CYS:HB3	1:C:431:GLY:O	2.16	0.45
1:C:470:THR:OG1	1:C:471:GLU:N	2.50	0.45
1:C:959:LEU:O	1:C:963:VAL:HG23	2.17	0.45
1:A:542:ASN:HA	1:A:547:THR:HG22	1.99	0.45
1:C:449:TYR:HA	1:C:496:GLY:HA2	1.98	0.45
1:A:91:TYR:OH	1:A:191:GLU:OE1	2.33	0.45
1:A:329:PHE:O	1:A:580:GLN:NE2	2.31	0.45
1:C:106:PHE:HD1	1:C:238:PHE:HB2	1.81	0.45
1:C:704:SER:HB3	1:A:790:LYS:HE3	1.98	0.45
1:A:739:THR:O	1:A:743:CYS:HB2	2.17	0.45
1:A:451:TYR:HB2	1:A:495:TYR:HB2	1.98	0.45
1:A:472:ILE:HA	1:A:490:PHE:HA	1.98	0.45
1:B:131:CYS:HB2	1:B:133:PHE:CZ	2.52	0.45
1:C:365:TYR:CD2	1:C:387:LEU:HB3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:356:LYS:HE3	1:C:356:LYS:HB2	1.85	0.45
1:C:991:VAL:HG22	1:C:992:GLN:HE21	1.82	0.45
1:B:583:GLU:OE1	1:B:584:ILE:N	2.47	0.44
1:C:200:TYR:HA	1:C:230:PRO:HA	1.99	0.44
1:A:376:THR:OG1	1:A:435:ALA:HB3	2.17	0.44
1:A:400:PHE:HD1	1:A:402:ILE:HG23	1.82	0.44
1:C:418:ILE:HD13	1:C:422:ASN:HD22	1.82	0.44
1:A:350:VAL:HG12	1:A:422:ASN:HB3	2.00	0.44
1:C:965:GLN:HE22	1:A:758:SER:H	1.64	0.44
1:C:1073:LYS:HB3	1:C:1073:LYS:HE2	1.72	0.44
1:C:1093:GLY:CA	1:C:1105:THR:O	2.66	0.44
1:B:126:VAL:O	1:B:171:VAL:HA	2.18	0.44
1:B:383:SER:HB3	1:B:386:LYS:HD2	1.99	0.44
1:B:431:GLY:N	1:B:513:LEU:O	2.50	0.44
1:C:117:LEU:HD12	1:C:130:VAL:HG12	2.00	0.44
1:A:722:VAL:HA	1:A:1064:HIS:O	2.18	0.44
1:B:586:ASP:OD1	1:B:587:ILE:N	2.50	0.44
1:C:759:PHE:O	1:C:763:LEU:HG	2.18	0.44
1:A:825:LYS:NZ	1:A:942:ALA:HA	2.33	0.44
1:A:1005:GLN:HE21	1:A:1005:GLN:HB2	1.56	0.44
1:B:276:LEU:HD11	1:B:304:LYS:HA	2.00	0.44
1:B:350:VAL:HG12	1:B:422:ASN:HB3	2.00	0.44
1:B:386:LYS:HD3	1:C:982:SER:O	2.18	0.44
1:A:766:ALA:O	1:A:770:ILE:HG23	2.17	0.44
1:B:1005:GLN:HE21	1:B:1005:GLN:HB2	1.56	0.43
1:C:108:THR:HG23	1:C:109:THR:H	1.82	0.43
1:C:570:ALA:HB2	1:A:964:LYS:HA	1.99	0.43
1:A:57:PRO:O	1:A:60:SER:HB2	2.18	0.43
1:B:344:ALA:HB3	1:B:347:PHE:CZ	2.53	0.43
1:C:287:ASP:HB3	1:C:306:PHE:CE2	2.53	0.43
1:C:815:ARG:HG2	1:C:819:GLU:HB2	2.00	0.43
1:A:543:PHE:CE1	1:A:579:PRO:HD3	2.52	0.43
1:C:357:ARG:HG3	1:C:396:TYR:HE2	1.84	0.43
1:A:344:ALA:HB3	1:A:347:PHE:CZ	2.53	0.43
1:A:901:GLN:NE2	1:A:905:ARG:HH21	2.14	0.43
1:A:1045:LYS:HA	1:A:1045:LYS:HD3	1.86	0.43
1:B:122:ASN:C	1:B:124:THR:H	2.22	0.43
1:B:277:LEU:HD22	1:B:285:ILE:HD13	1.99	0.43
1:B:976:VAL:HG13	1:B:979:ASP:HB3	1.99	0.43
1:B:982:SER:O	1:A:386:LYS:HD3	2.18	0.43
1:A:328:ARG:HH21	1:A:579:PRO:HD2	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1040:VAL:HG21	1:A:1035:GLY:HA3	1.99	0.43
1:C:1086:LYS:HD2	1:C:1122:VAL:HG21	1.99	0.43
1:A:964:LYS:HE3	1:A:964:LYS:HB2	1.89	0.43
1:C:802:PHE:HD1	1:C:805:ILE:HD11	1.83	0.43
1:A:136:CYS:HB3	1:A:159:VAL:HA	2.01	0.43
1:B:37:TYR:OH	1:B:53:ASP:OD2	2.36	0.43
1:B:568:ASP:O	1:B:569:ILE:HG13	2.19	0.43
1:A:570:ALA:O	1:A:572:THR:HG23	2.19	0.43
1:B:117:LEU:HA	1:B:129:LYS:O	2.19	0.43
1:A:53:ASP:HB3	1:A:55:PHE:HE2	1.84	0.43
1:B:1129:VAL:HB	1:B:1132:ILE:HG13	2.00	0.43
1:A:323:THR:O	1:A:324:GLU:HG2	2.18	0.43
1:A:65:PHE:CD2	1:A:82:PRO:HB3	2.54	0.42
1:B:30:ASN:HB3	1:B:32:PHE:CE1	2.54	0.42
1:B:57:PRO:O	1:B:60:SER:HB2	2.19	0.42
1:B:376:THR:OG1	1:B:435:ALA:HB3	2.19	0.42
1:B:973:ILE:HD11	1:B:983:ARG:NH2	2.34	0.42
1:C:726:ILE:HG12	1:C:1061:VAL:HG22	2.01	0.42
1:A:962:LEU:HD13	1:A:1007:TYR:HB2	2.01	0.42
1:B:309:GLU:O	1:B:313:TYR:OH	2.31	0.42
1:B:231:ILE:HG22	1:B:233:ILE:HG13	2.01	0.42
1:B:399:SER:HB3	1:B:511:VAL:HG12	2.01	0.42
1:B:720:ILE:HD11	1:B:927:PHE:HB2	2.02	0.42
1:A:903:ALA:HB2	1:A:916:LEU:HD23	2.01	0.42
1:B:722:VAL:HA	1:B:1064:HIS:O	2.19	0.42
1:B:1093:GLY:CA	1:B:1105:THR:O	2.68	0.42
1:C:964:LYS:HE2	1:C:964:LYS:HB2	1.76	0.42
1:C:1077:THR:HG22	1:C:1078:ALA:H	1.84	0.42
1:A:439:ASN:ND2	1:A:506:GLN:HE21	2.16	0.42
1:A:906:PHE:O	1:A:909:ILE:HG12	2.20	0.42
1:C:303:LEU:HD12	1:C:308:VAL:HG12	2.02	0.42
1:C:318:PHE:CE1	1:C:615:VAL:HG21	2.55	0.42
1:C:662:CYS:HB2	1:C:697:MET:SD	2.60	0.42
1:B:962:LEU:HD13	1:B:1007:TYR:HB2	2.02	0.42
1:B:945:LEU:HD23	1:B:945:LEU:HA	1.85	0.42
1:A:353:TRP:HB3	1:A:400:PHE:HB3	2.01	0.42
1:A:938:LEU:HD23	1:A:938:LEU:HA	1.82	0.42
1:B:759:PHE:O	1:B:763:LEU:HG	2.19	0.42
1:A:89:GLY:HA2	1:A:194:PHE:O	2.20	0.42
1:A:106:PHE:HA	1:A:238:PHE:HA	2.01	0.42
1:A:1125:ASN:OD1	1:A:1125:ASN:N	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:VAL:HG22	1:B:237:ARG:HB3	2.00	0.42
1:B:553:THR:HG22	1:B:554:GLU:N	2.35	0.42
1:C:854:LYS:HE3	1:C:854:LYS:HB3	1.94	0.42
1:B:906:PHE:O	1:B:909:ILE:HG12	2.20	0.41
1:C:353:TRP:HZ3	1:C:355:ARG:HD2	1.85	0.41
1:C:520:ALA:HB1	1:C:521:PRO:HD2	2.02	0.41
1:C:906:PHE:O	1:C:909:ILE:HG12	2.20	0.41
1:A:610:VAL:O	1:A:650:LEU:HD12	2.20	0.41
1:C:901:GLN:NE2	1:C:905:ARG:HH21	2.16	0.41
1:A:380:TYR:HE2	1:A:412:PRO:HD2	1.85	0.41
1:A:947:LYS:HB3	1:A:947:LYS:HE2	1.85	0.41
1:A:976:VAL:HG13	1:A:979:ASP:HB3	2.02	0.41
1:A:981:LEU:HD23	1:A:981:LEU:HA	1.84	0.41
1:B:357:ARG:HB3	1:B:357:ARG:HH11	1.85	0.41
1:B:811:LYS:HD2	1:B:815:ARG:HB2	2.02	0.41
1:C:418:ILE:HG22	1:C:423:TYR:O	2.20	0.41
1:A:138:ASP:N	1:A:138:ASP:OD1	2.51	0.41
1:A:190:ARG:HB3	1:A:192:PHE:CE2	2.54	0.41
1:B:140:PHE:HA	1:B:242:LEU:C	2.41	0.41
1:C:106:PHE:HB3	1:C:235:ILE:HD13	2.02	0.41
1:B:28:TYR:HD2	1:B:61:ASN:HD22	1.69	0.41
1:B:443:SER:HA	1:B:497:PHE:HB2	2.02	0.41
1:B:520:ALA:HB1	1:B:521:PRO:HD2	2.03	0.41
1:C:53:ASP:HB2	1:C:55:PHE:HE2	1.84	0.41
1:C:124:THR:O	1:C:125:ASN:ND2	2.53	0.41
1:C:816:SER:OG	1:C:819:GLU:HG3	2.21	0.41
1:C:1126:CYS:O	1:C:1132:ILE:HD11	2.20	0.41
1:A:34:ARG:HH12	1:A:217:PRO:HG2	1.85	0.41
1:A:905:ARG:HD3	1:A:1049:LEU:O	2.21	0.41
1:B:1053:PRO:O	1:B:1054:GLN:NE2	2.54	0.41
1:C:215:ASP:OD1	1:C:215:ASP:N	2.54	0.41
1:C:357:ARG:HG3	1:C:396:TYR:CE2	2.56	0.41
1:C:715:PRO:HD3	1:A:894:LEU:HD13	2.03	0.41
1:C:856:ASN:OD1	1:C:856:ASN:N	2.53	0.41
1:C:938:LEU:HD23	1:C:938:LEU:HA	1.82	0.41
1:C:980:ILE:HD11	1:C:989:ALA:HA	2.03	0.41
1:A:233:ILE:HD13	1:A:233:ILE:HA	1.83	0.41
1:A:357:ARG:HG3	1:A:396:TYR:HE1	1.86	0.41
1:A:759:PHE:O	1:A:763:LEU:HG	2.21	0.41
1:B:1106:GLN:HE21	1:B:1109:PHE:HB3	1.85	0.41
1:C:702:GLU:OE2	1:A:790:LYS:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:720:ILE:HD11	1:C:927:PHE:HB2	2.02	0.41
1:A:645:THR:OG1	1:A:646:ARG:N	2.54	0.41
1:B:400:PHE:HD1	1:B:402:ILE:HG23	1.85	0.41
1:C:720:ILE:HG21	1:C:923:ILE:HG23	2.02	0.41
1:A:520:ALA:HB1	1:A:521:PRO:HD2	2.03	0.41
1:B:120:VAL:HG12	1:B:121:ASN:N	2.35	0.41
1:B:131:CYS:HB3	1:B:163:ALA:HA	2.02	0.41
1:B:639:GLY:H	1:B:651:ILE:HG21	1.86	0.41
1:B:901:GLN:NE2	1:B:905:ARG:HH21	2.18	0.41
1:B:1023:ASN:O	1:B:1027:THR:HG23	2.21	0.41
1:C:28:TYR:HD2	1:C:61:ASN:HD22	1.68	0.41
1:A:452:LEU:HD13	1:A:493:GLN:O	2.21	0.41
1:B:190:ARG:HB3	1:B:192:PHE:HE2	1.86	0.41
1:B:528:LYS:HA	1:B:528:LYS:HD2	1.74	0.41
1:A:30:ASN:HB3	1:A:32:PHE:CE1	2.56	0.41
1:A:224:GLU:OE1	1:A:224:GLU:N	2.50	0.41
1:B:541:PHE:HD1	1:B:543:PHE:HB2	1.86	0.40
1:B:916:LEU:HD12	1:B:923:ILE:HD12	2.03	0.40
1:B:959:LEU:HD12	1:B:959:LEU:HA	1.91	0.40
1:C:63:THR:O	1:C:63:THR:OG1	2.39	0.40
1:C:118:LEU:HD11	1:C:120:VAL:HG23	2.02	0.40
1:C:318:PHE:HE1	1:C:615:VAL:HG21	1.86	0.40
1:A:1038:LYS:HE2	1:A:1038:LYS:HB3	1.95	0.40
1:A:341:VAL:HG22	1:A:356:LYS:HD2	2.04	0.40
1:A:528:LYS:HA	1:A:528:LYS:HD2	1.78	0.40
1:A:825:LYS:NZ	1:A:941:THR:O	2.33	0.40
1:A:919:ASN:OD1	1:A:919:ASN:N	2.54	0.40
1:A:1001:LEU:HD12	1:A:1001:LEU:HA	1.92	0.40
1:C:776:LYS:HE2	1:C:776:LYS:HB2	1.73	0.40
1:C:1074:ASN:N	1:C:1074:ASN:OD1	2.55	0.40
1:B:323:THR:OG1	1:B:324:GLU:N	2.54	0.40
1:B:903:ALA:HB2	1:B:916:LEU:HD23	2.04	0.40
1:C:544:ASN:HD21	1:C:579:PRO:HB3	1.86	0.40
1:C:1028:LYS:HE3	1:C:1028:LYS:HB2	1.94	0.40
1:A:296:LEU:O	1:A:299:THR:OG1	2.31	0.40
1:A:577:ARG:HA	1:A:584:ILE:HA	2.04	0.40
1:B:328:ARG:HA	1:B:328:ARG:HD3	1.75	0.40
1:B:919:ASN:OD1	1:B:919:ASN:N	2.54	0.40
1:C:1038:LYS:HE2	1:C:1038:LYS:HB3	1.93	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	945/1273 (74%)	897 (95%)	48 (5%)	0	100	100
1	B	945/1273 (74%)	890 (94%)	55 (6%)	0	100	100
1	C	945/1273 (74%)	895 (95%)	50 (5%)	0	100	100
All	All	2835/3819 (74%)	2682 (95%)	153 (5%)	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	792/1112 (71%)	746 (94%)	46 (6%)	17	44
1	B	792/1112 (71%)	748 (94%)	44 (6%)	17	45
1	C	796/1112 (72%)	751 (94%)	45 (6%)	17	44
All	All	2380/3336 (71%)	2245 (94%)	135 (6%)	20	44

All (135) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	40	ASP
1	B	63	THR
1	B	83	VAL
1	B	86	PHE
1	B	90	VAL

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Mol	Chain	Res	Type
1	B	166	CYS
1	B	172	SER
1	B	175	PHE
1	B	189	LEU
1	B	314	GLN
1	B	318	PHE
1	B	319	ARG
1	B	389	ASP
1	B	392	PHE
1	B	558	LYS
1	B	588	THR
1	B	599	THR
1	B	611	LEU
1	B	617	CYS
1	B	658	ASN
1	B	705	VAL
1	B	720	ILE
1	B	723	THR
1	B	732	THR
1	B	734	THR
1	B	742	ILE
1	B	748	GLU
1	B	752	LEU
1	B	785	VAL
1	B	808	ASP
1	B	816	SER
1	B	854	LYS
1	B	855	PHE
1	B	875	SER
1	B	902	MET
1	B	919	ASN
1	B	937	SER
1	B	950	ASP
1	B	967	SER
1	B	973	ILE
1	B	990	GLU
1	B	1058	HIS
1	B	1097	SER
1	B	1108	ASN
1	C	40	ASP
1	C	63	THR
1	C	86	PHE

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Mol	Chain	Res	Type
1	C	94	SER
1	C	165	ASN
1	C	223	LEU
1	C	266	TYR
1	C	282	ASN
1	C	290	ASP
1	C	298	GLU
1	C	302	THR
1	C	314	GLN
1	C	316	SER
1	C	318	PHE
1	C	353	TRP
1	C	383	SER
1	C	392	PHE
1	C	514	SER
1	C	525	CYS
1	C	555	SER
1	C	558	LYS
1	C	565	PHE
1	C	585	LEU
1	C	599	THR
1	C	617	CYS
1	C	654	GLU
1	C	711	SER
1	C	720	ILE
1	C	723	THR
1	C	731	MET
1	C	732	THR
1	C	752	LEU
1	C	760	CYS
1	C	770	ILE
1	C	785	VAL
1	C	816	SER
1	C	854	LYS
1	C	875	SER
1	C	935	GLN
1	C	980	ILE
1	C	991	VAL
1	C	1058	HIS
1	C	1067	TYR
1	C	1074	ASN
1	C	1146	ASP

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Mol	Chain	Res	Type
1	A	63	THR
1	A	90	VAL
1	A	98	SER
1	A	104	TRP
1	A	127	VAL
1	A	158	ARG
1	A	166	CYS
1	A	189	LEU
1	A	294	ASP
1	A	305	SER
1	A	307	THR
1	A	316	SER
1	A	318	PHE
1	A	392	PHE
1	A	433	VAL
1	A	450	ASN
1	A	524	VAL
1	A	553	THR
1	A	565	PHE
1	A	573	THR
1	A	585	LEU
1	A	599	THR
1	A	645	THR
1	A	658	ASN
1	A	723	THR
1	A	731	MET
1	A	732	THR
1	A	734	THR
1	A	742	ILE
1	A	752	LEU
1	A	770	ILE
1	A	778	THR
1	A	791	THR
1	A	816	SER
1	A	853	GLN
1	A	855	PHE
1	A	858	LEU
1	A	875	SER
1	A	919	ASN
1	A	937	SER
1	A	950	ASP
1	A	952	VAL

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Mol	Chain	Res	Type
1	A	973	ILE
1	A	994	ASP
1	A	1058	HIS
1	A	1091	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	81	ASN
1	B	122	ASN
1	B	280	ASN
1	B	317	ASN
1	B	536	ASN
1	B	544	ASN
1	B	644	GLN
1	B	755	GLN
1	B	901	GLN
1	B	954	GLN
1	B	965	GLN
1	B	969	ASN
1	B	1002	GLN
1	B	1010	GLN
1	B	1101	HIS
1	C	66	HIS
1	C	125	ASN
1	C	388	ASN
1	C	422	ASN
1	C	437	ASN
1	C	536	ASN
1	C	644	GLN
1	C	751	ASN
1	C	755	GLN
1	C	901	GLN
1	C	914	ASN
1	C	919	ASN
1	C	965	GLN
1	C	992	GLN
1	C	1002	GLN
1	C	1101	HIS
1	C	1106	GLN
1	A	280	ASN
1	A	317	ASN

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Mol	Chain	Res	Type
1	A	388	ASN
1	A	439	ASN
1	A	506	GLN
1	A	536	ASN
1	A	544	ASN
1	A	644	GLN
1	A	755	GLN
1	A	901	GLN
1	A	954	GLN
1	A	965	GLN
1	A	1054	GLN
1	A	1101	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no 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.

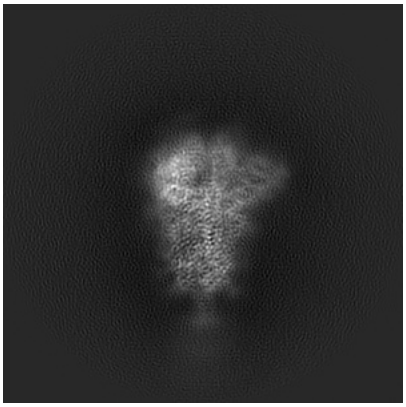
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-23215. These allow visual inspection of the internal detail of the map and identification of artifacts.

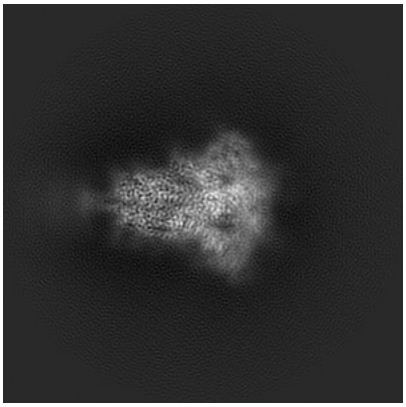
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

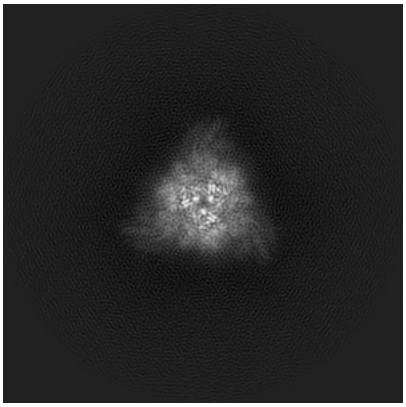
6.1.1 Primary map



X



Y

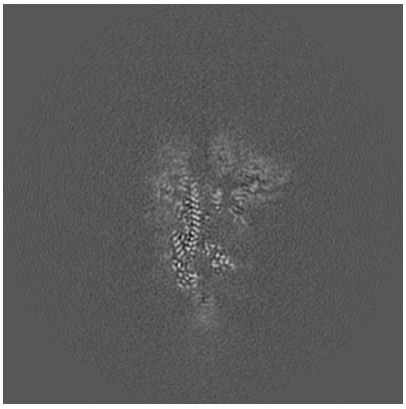


Z

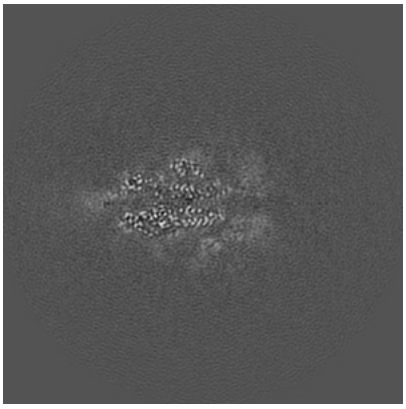
The images above show the map projected in three orthogonal directions.

6.2 Central slices

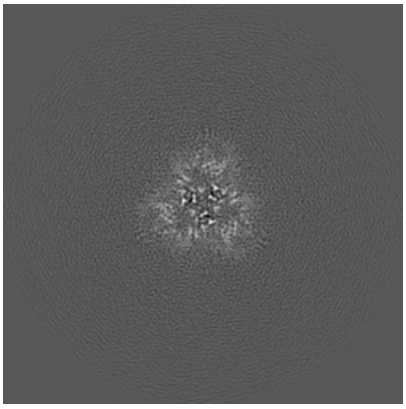
6.2.1 Primary map



X Index: 225



Y Index: 225

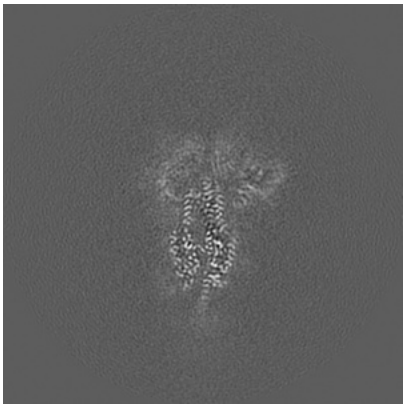


Z Index: 225

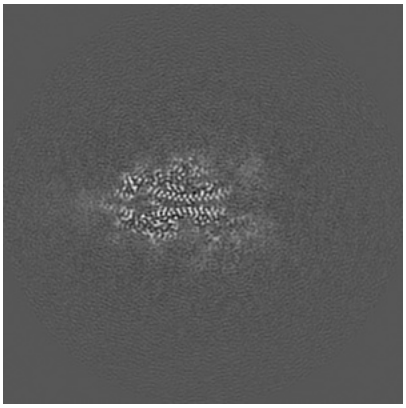
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

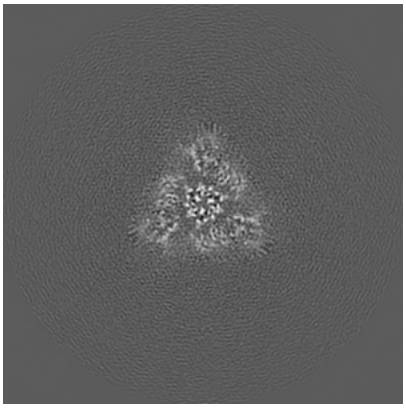
6.3.1 Primary map



X Index: 236



Y Index: 230

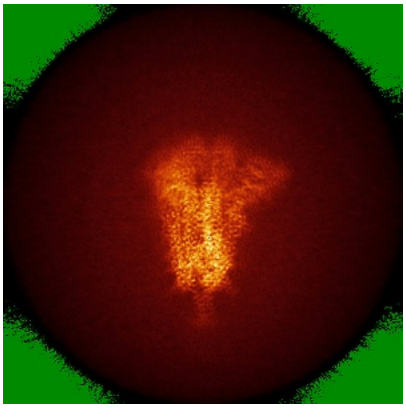


Z Index: 235

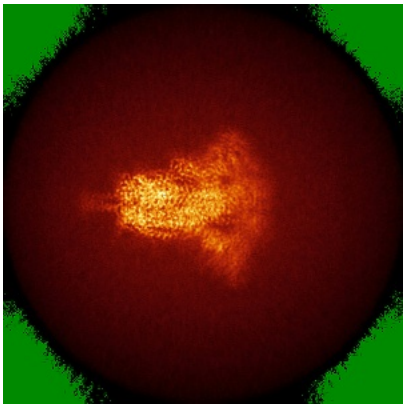
The images above show the largest variance slices of the map in three orthogonal directions.

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

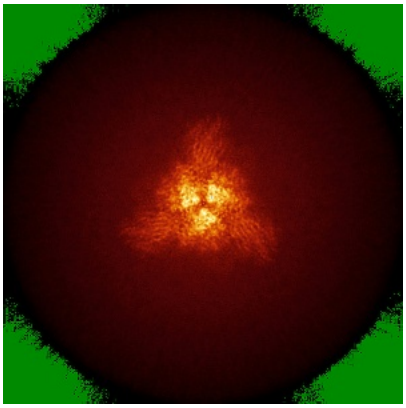
6.4.1 Primary map



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

X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.0003. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

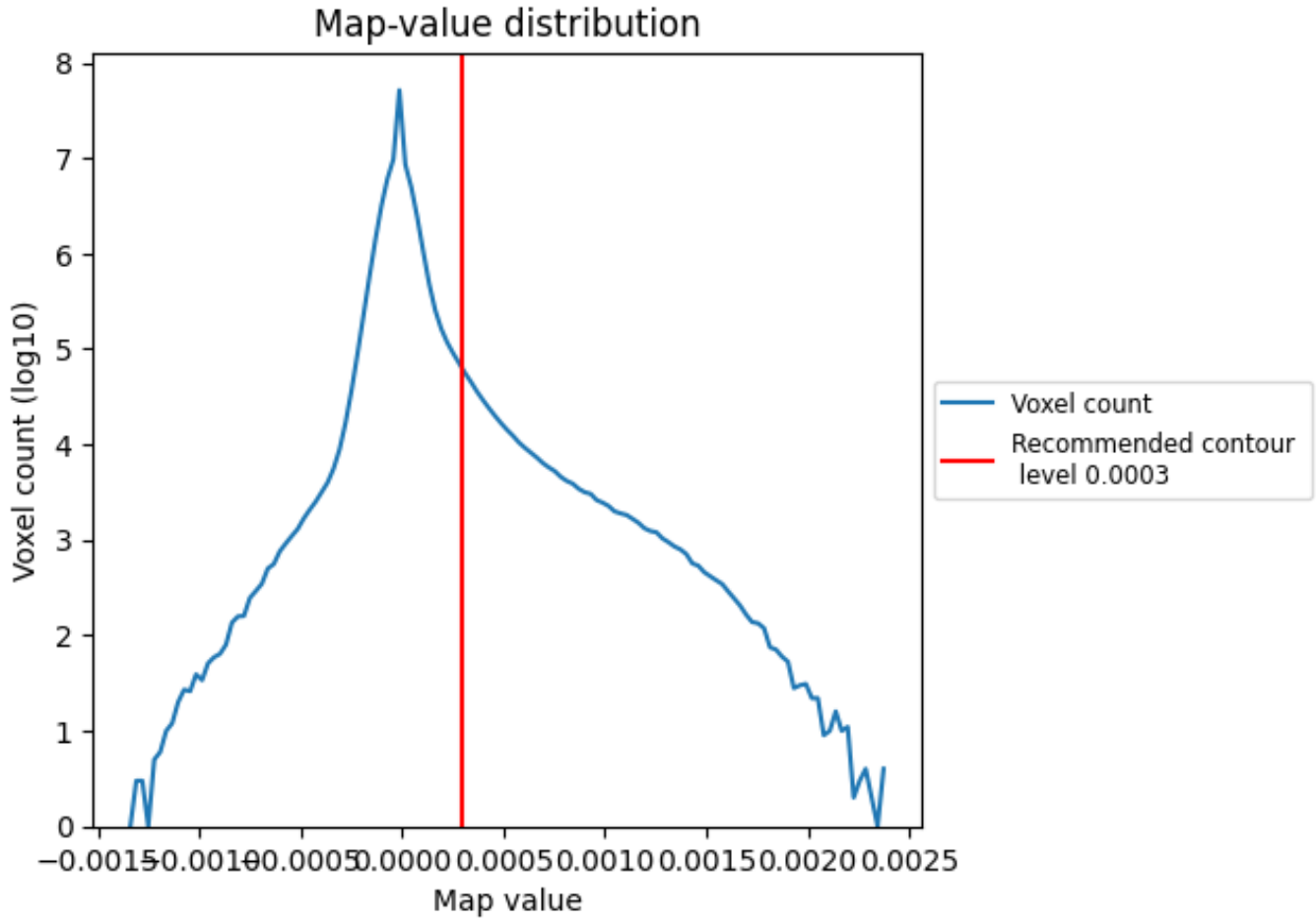
6.6 Mask visualisation (i)

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

7 Map analysis

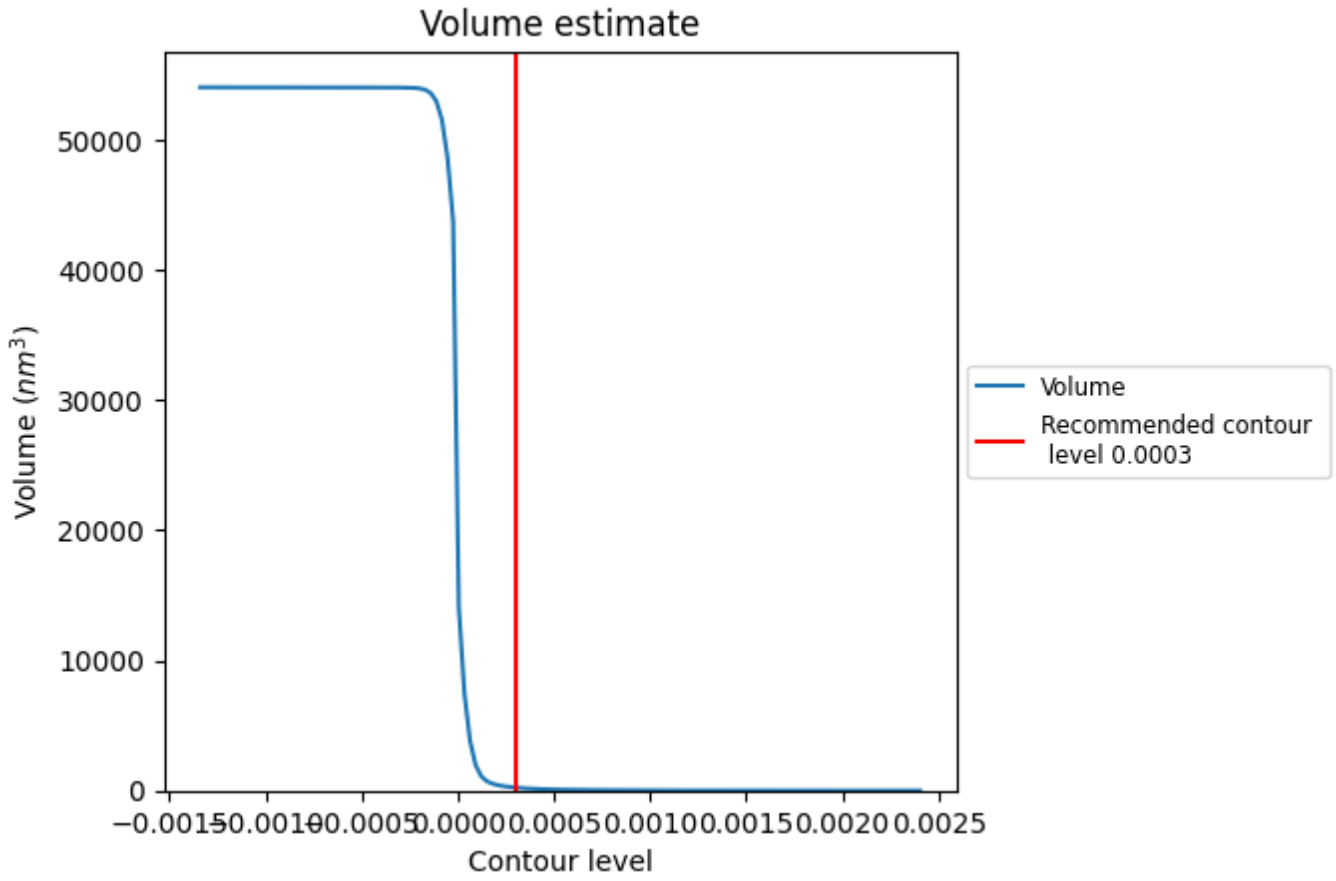
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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

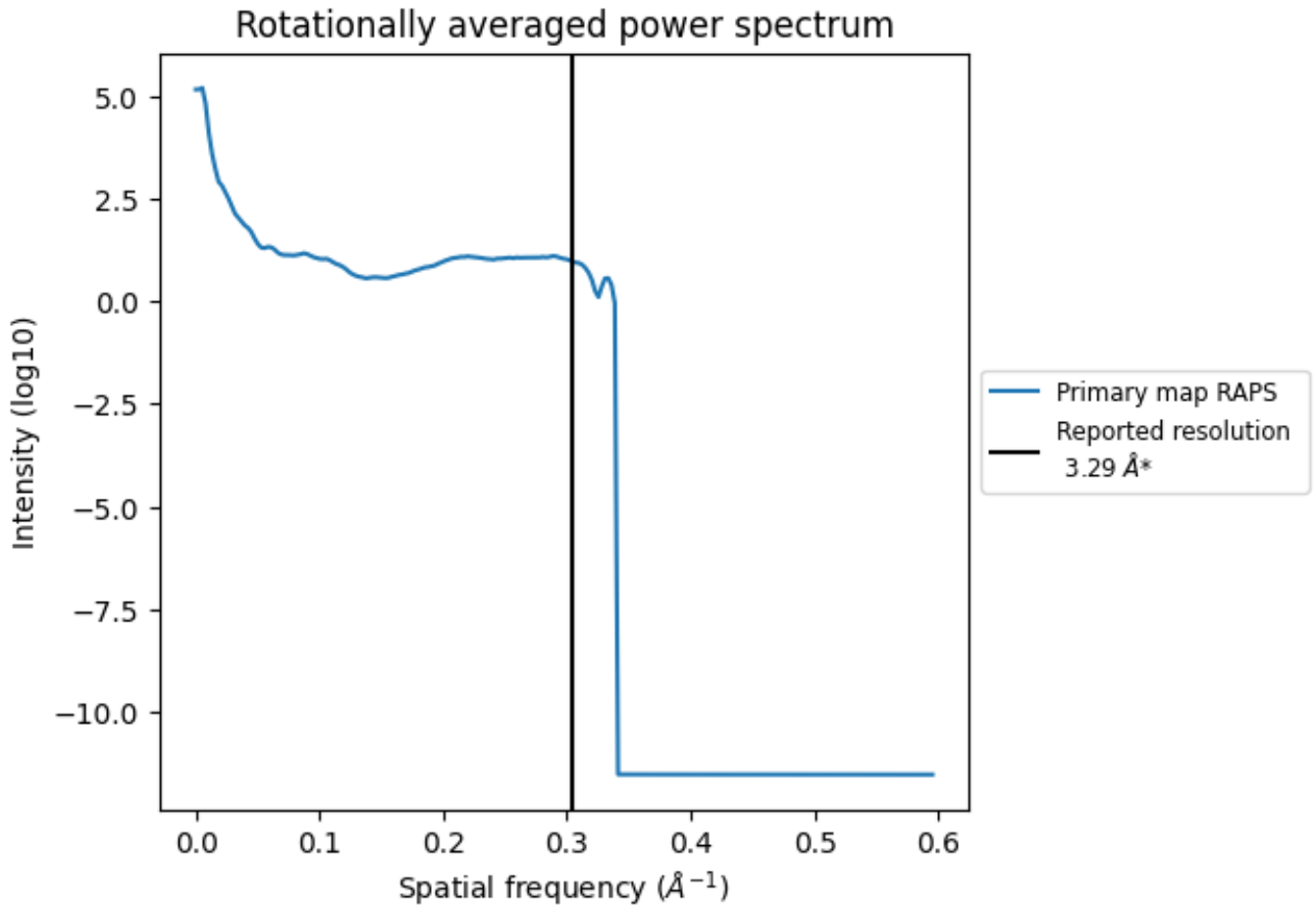
7.2 Volume estimate



The volume at the recommended contour level is 226 nm^3; this corresponds to an approximate mass of 204 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 0.304 $Å^{-1}$

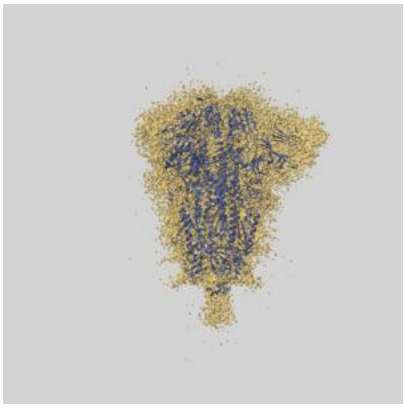
8 Fourier-Shell correlation

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

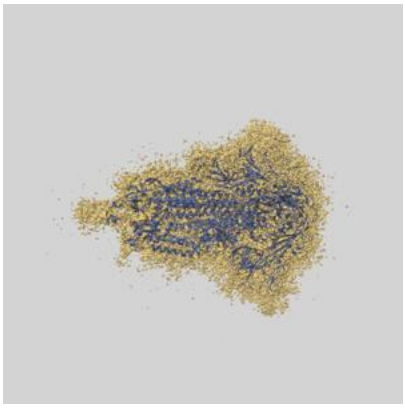
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-23215 and PDB model 7L7K. Per-residue inclusion information can be found in section 3 on page 4.

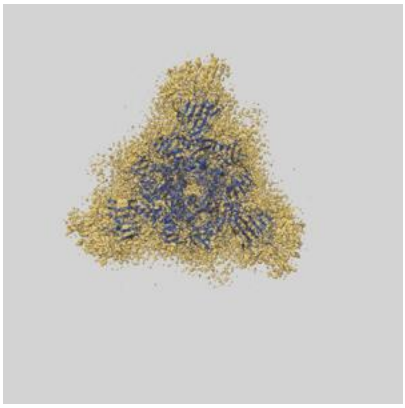
9.1 Map-model overlay



X



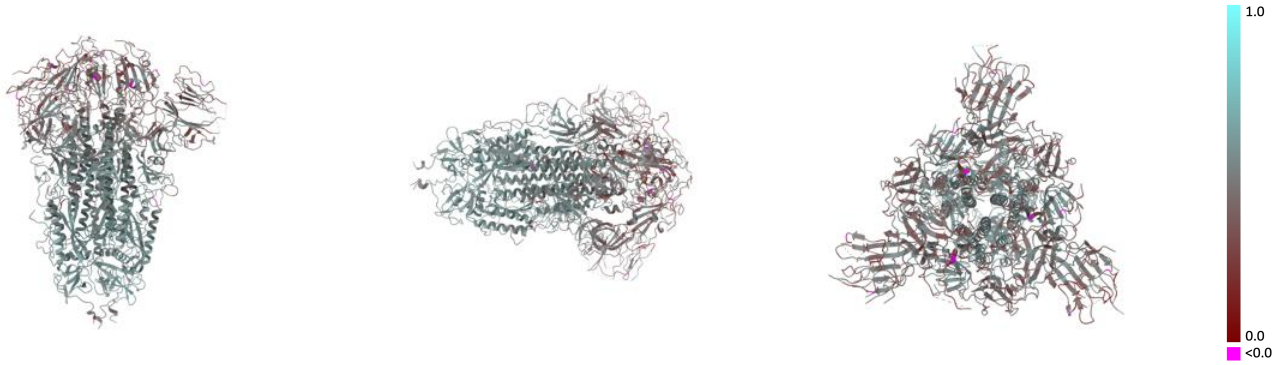
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0003 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

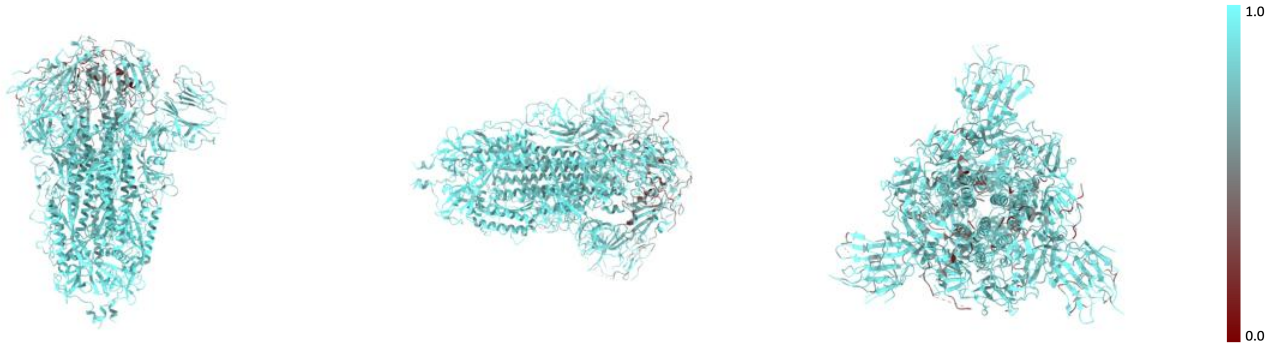
9.2 Q-score mapped to coordinate model



X Y Z

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

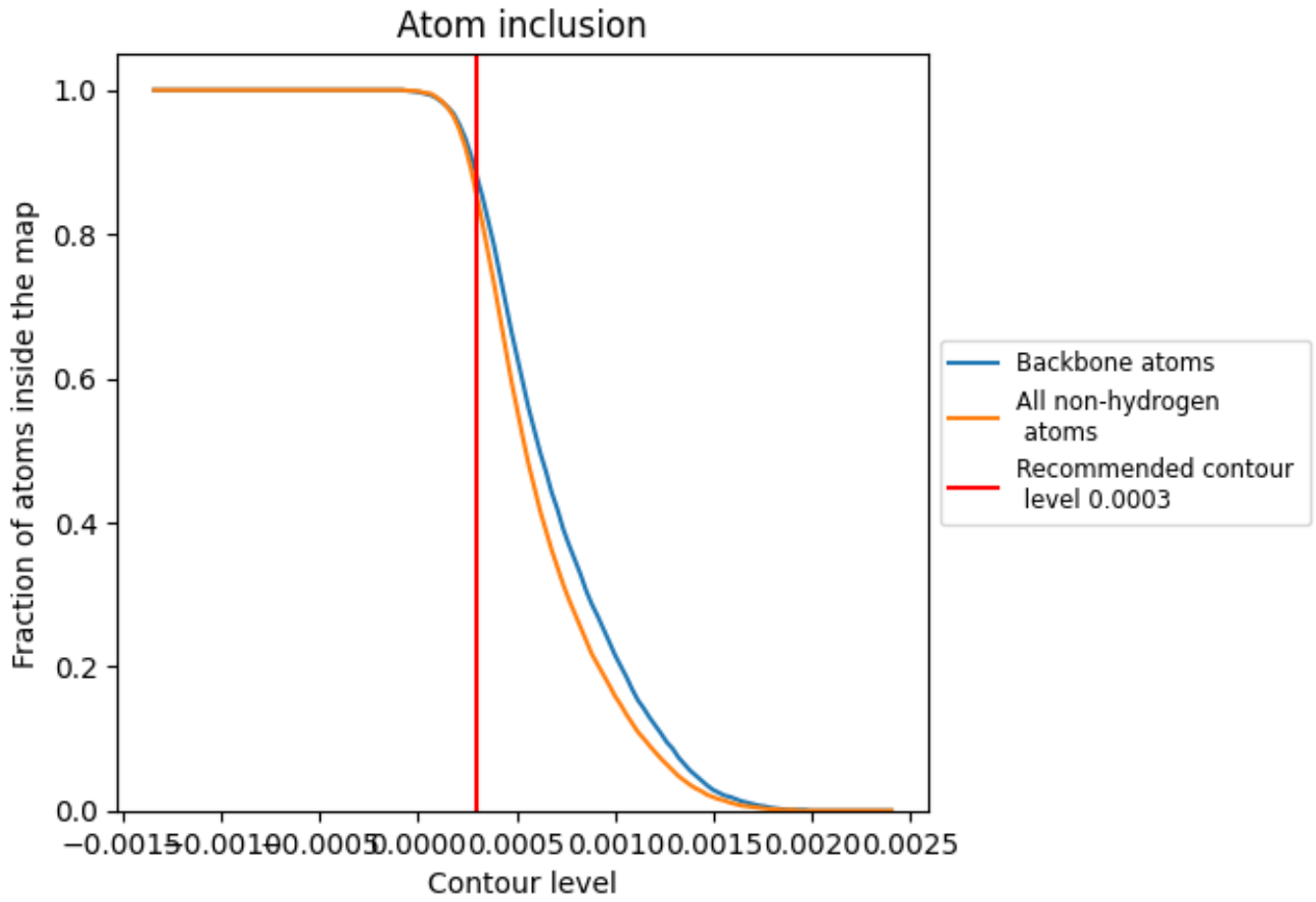
9.3 Atom inclusion mapped to coordinate model



X Y Z

The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.0003).

9.4 Atom inclusion (i)



At the recommended contour level, 88% of all backbone atoms, 85% 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 (0.0003) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8510	0.4840
A	0.8520	0.4850
B	0.8490	0.4820
C	0.8530	0.4850