



Full wwPDB EM Validation Report ⓘ

Mar 27, 2026 – 03:57 PM UTC

PDB ID : 7P6K / pdb_00007p6k
EMDB ID : EMD-13230
Title : Structure of homomeric LRRC8A Volume-Regulated Anion Channel in complex with synthetic nanobody Sb5
Authors : Deneka, D.; Rutz, S.; Sawicka, M.
Deposited on : 2021-07-16
Resolution : 3.80 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:

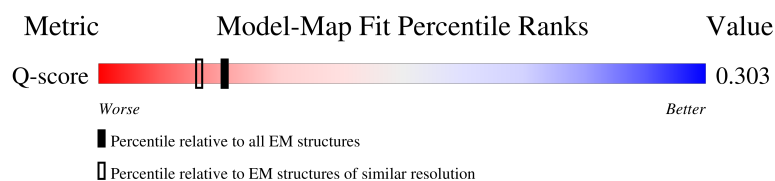
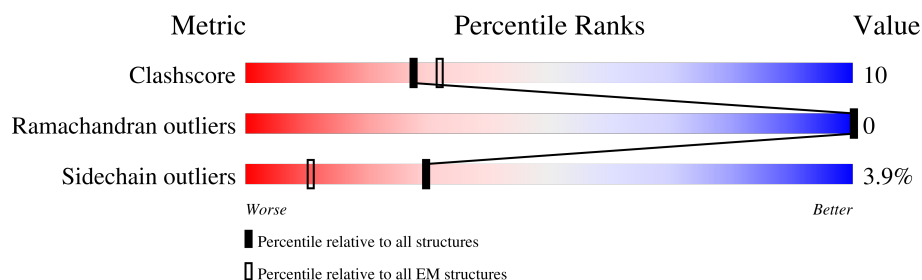
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.80 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	10198 (3.30 - 4.30)

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	810	16% 64% 24% • 11%
1	B	810	20% 64% 23% • 11%
1	C	810	17% 63% 24% • 11%
1	D	810	19% 64% 24% 11%

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Mol	Chain	Length	Quality of chain
1	E	810	17% 62% 26% • 11%
1	F	810	20% 65% 22% • 11%
2	G	144	66% 55% 22% • 21%
2	H	144	67% 56% 22% • 21%
2	I	144	65% 58% 19% • 21%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 38106 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 Volume-regulated anion channel subunit LRRC8A.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	718	Total	C	N	O	S	0	0
			5922	3853	997	1047	25		
1	B	718	Total	C	N	O	S	0	0
			5922	3853	997	1047	25		
1	C	718	Total	C	N	O	S	0	0
			5922	3853	997	1047	25		
1	D	718	Total	C	N	O	S	0	0
			5922	3853	997	1047	25		
1	E	718	Total	C	N	O	S	0	0
			5922	3853	997	1047	25		
1	F	718	Total	C	N	O	S	0	0
			5922	3853	997	1047	25		

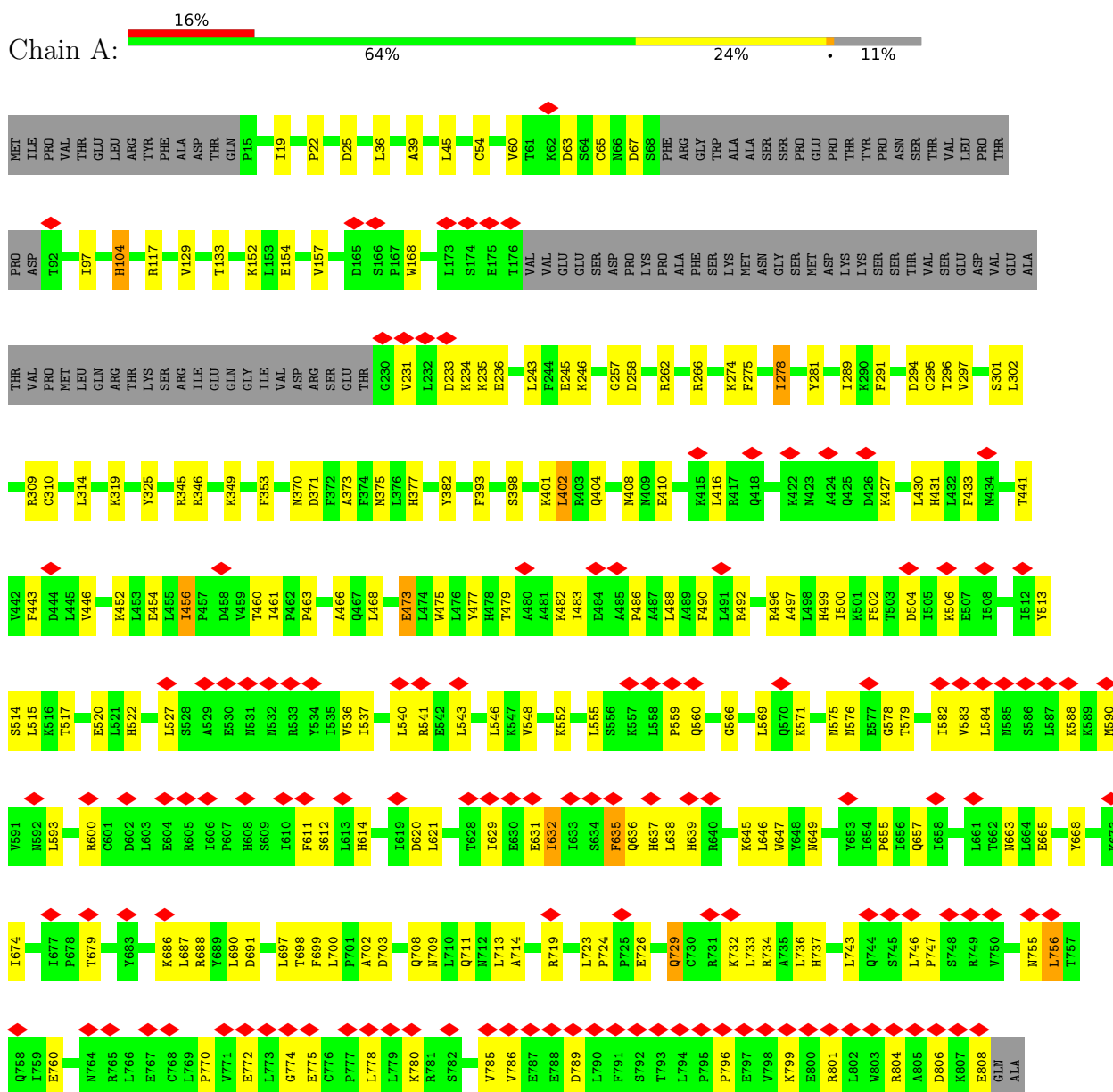
- Molecule 2 is a protein called synthetic nanobody Sb5.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	I	114	Total	C	N	O	S	0	0
			858	537	148	169	4		
2	G	114	Total	C	N	O	S	0	0
			858	537	148	169	4		
2	H	114	Total	C	N	O	S	0	0
			858	537	148	169	4		

3 Residue-property plots

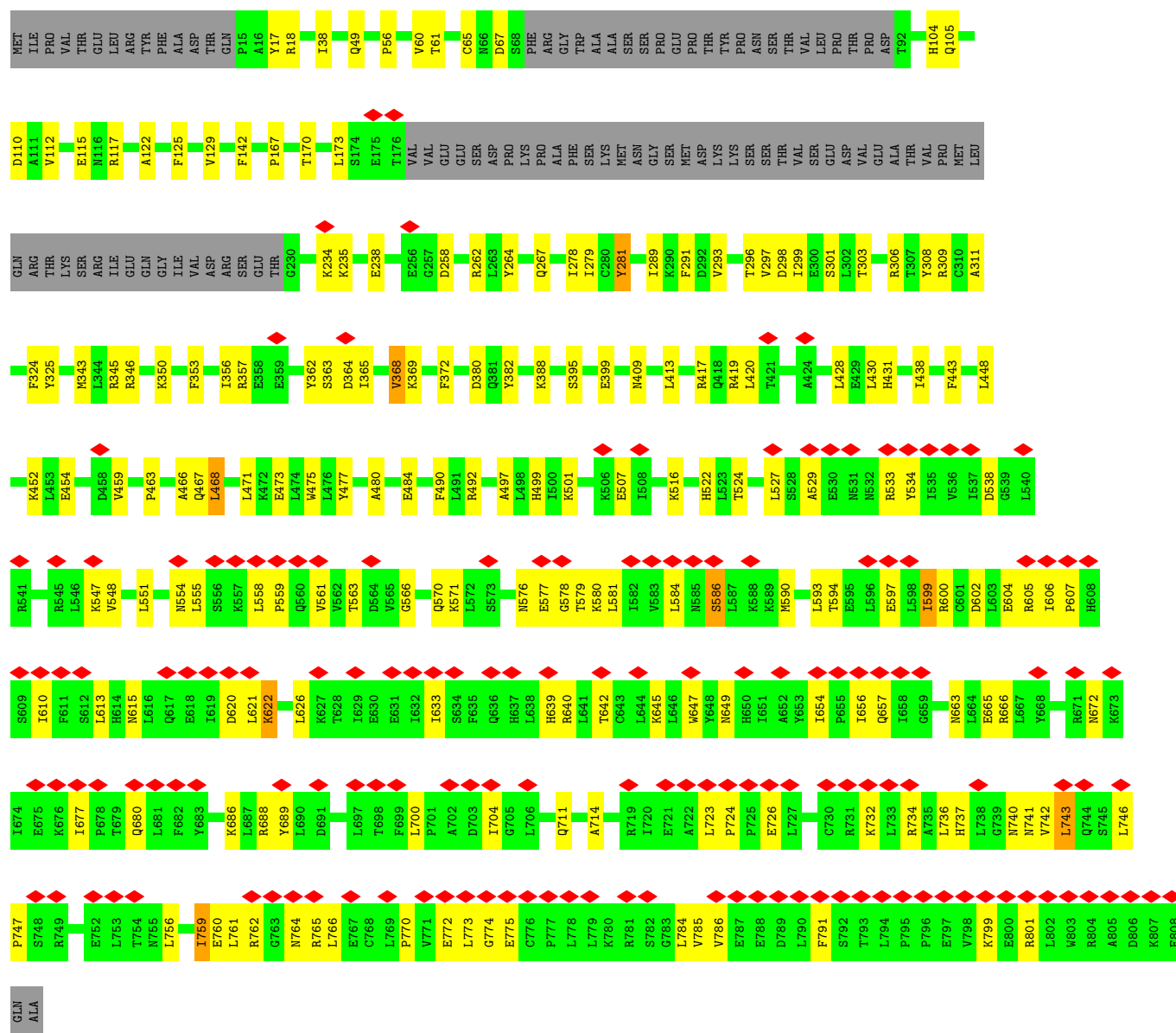
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Volume-regulated anion channel subunit LRRC8A



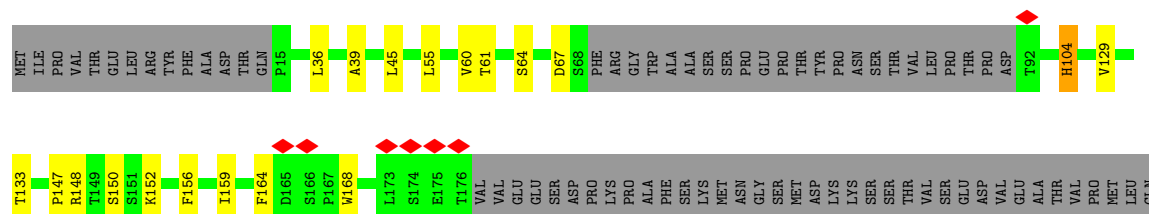
- Molecule 1: Volume-regulated anion channel subunit LRRC8A

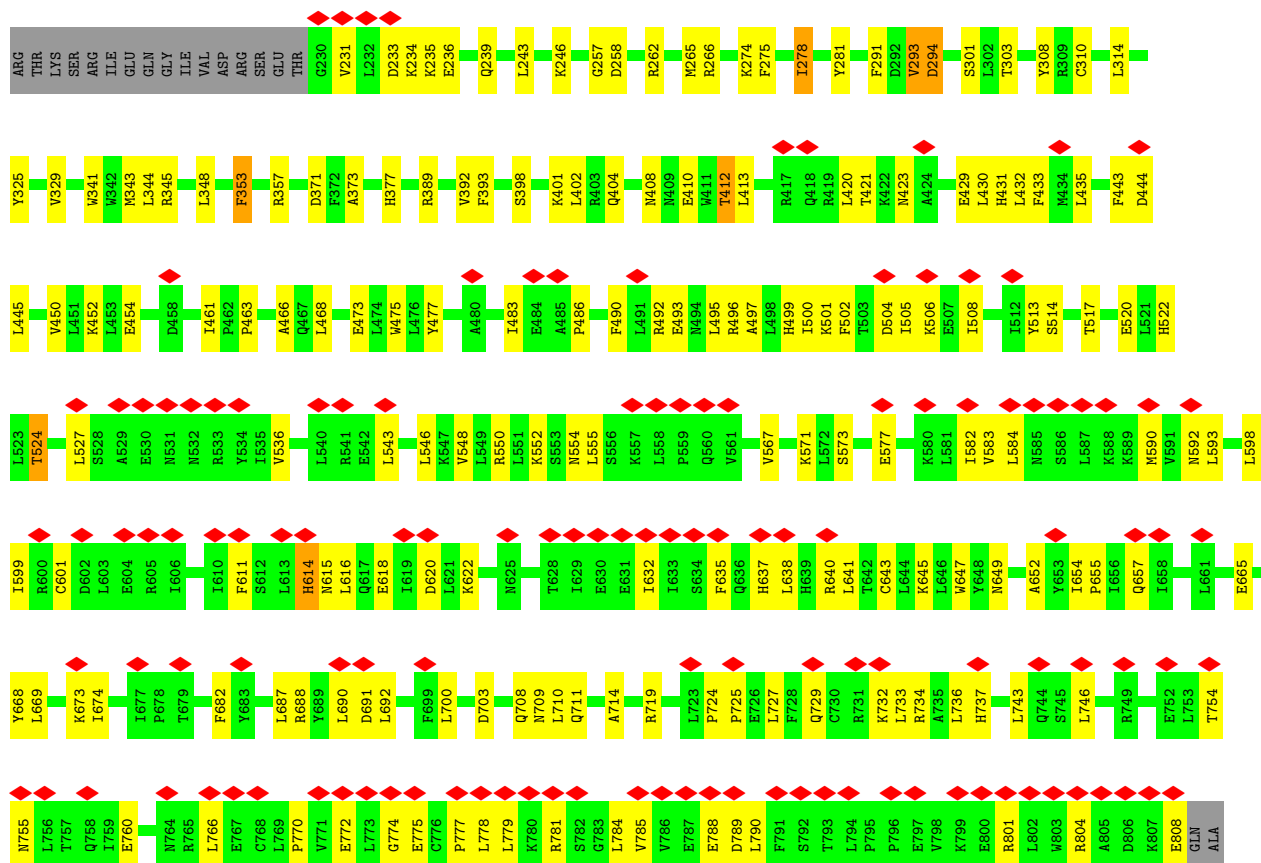
Chain B: 20% 64% 23% 11%



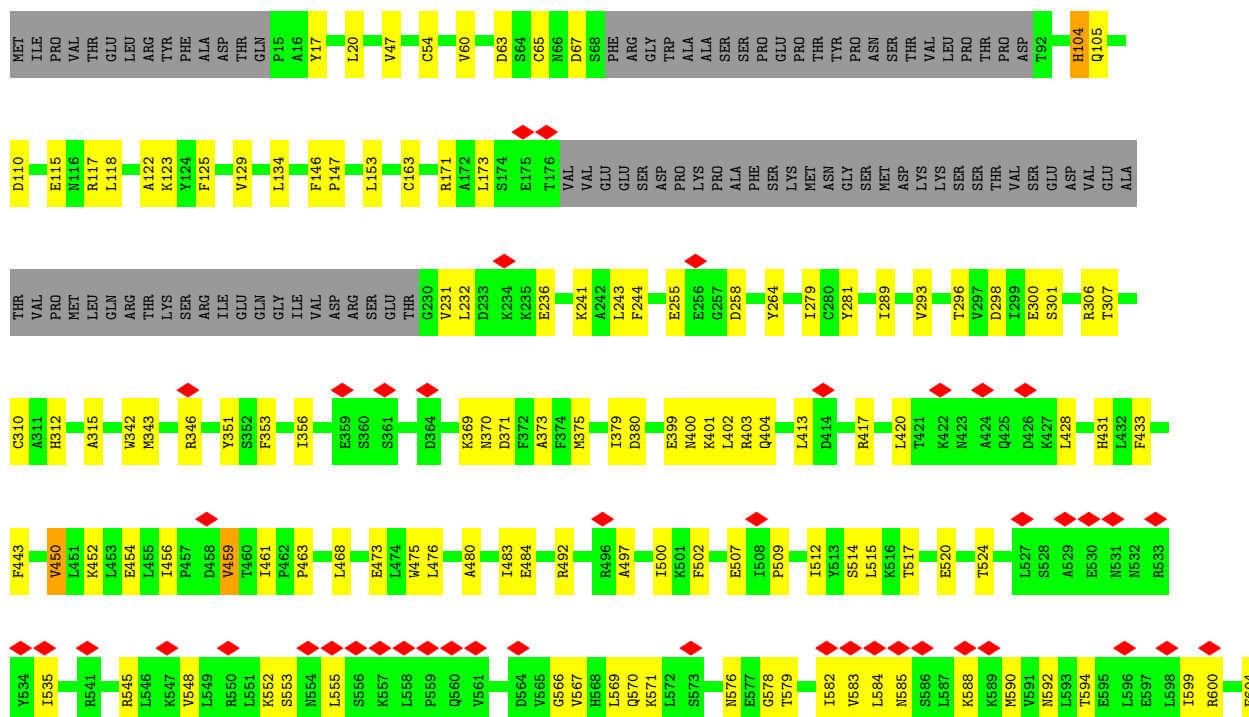
• Molecule 1: Volume-regulated anion channel subunit LRRC8A

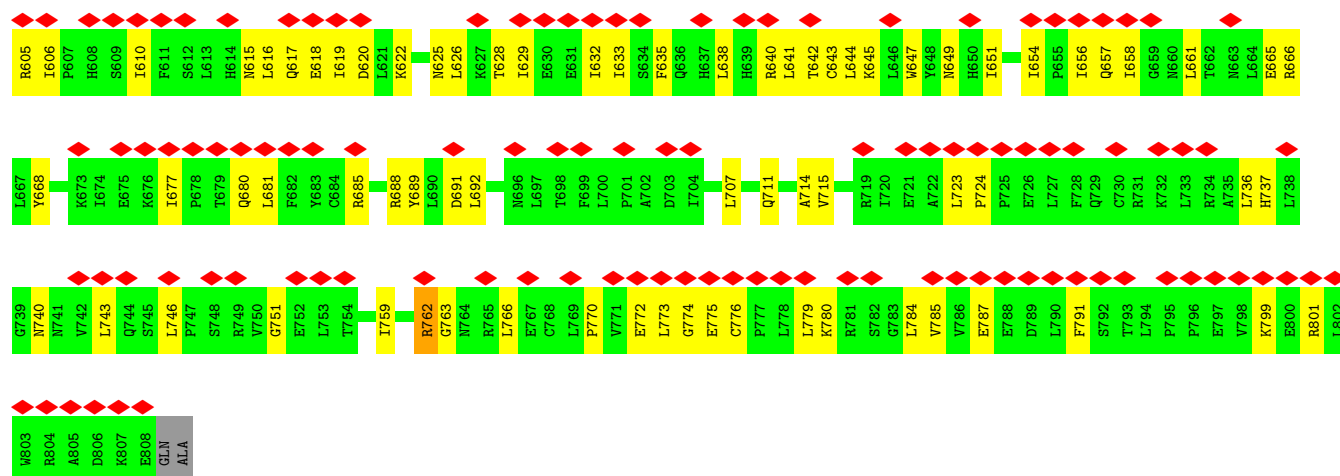
Chain C: 17% 63% 24% 11%



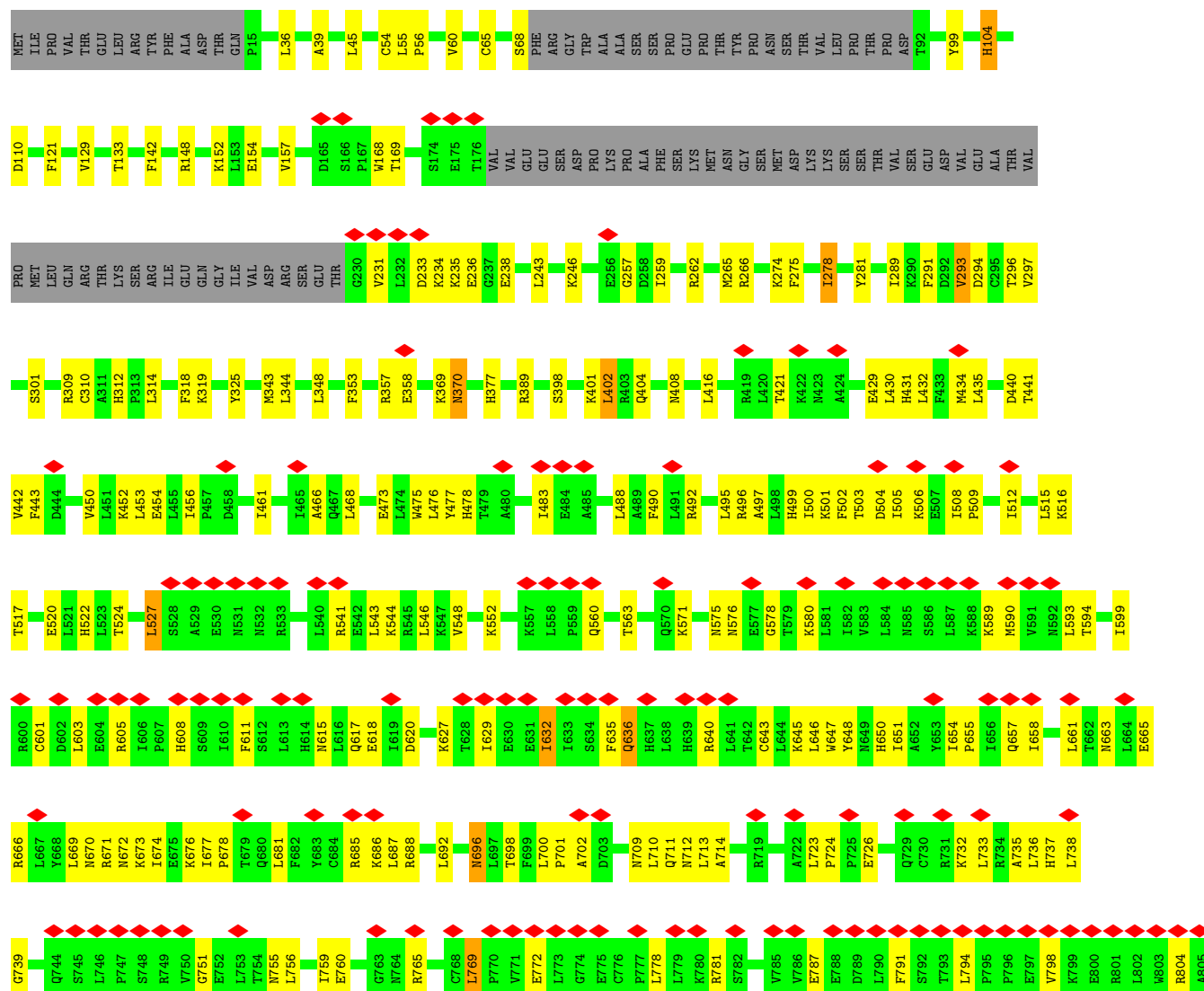


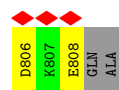
● Molecule 1: Volume-regulated anion channel subunit LRRC8A



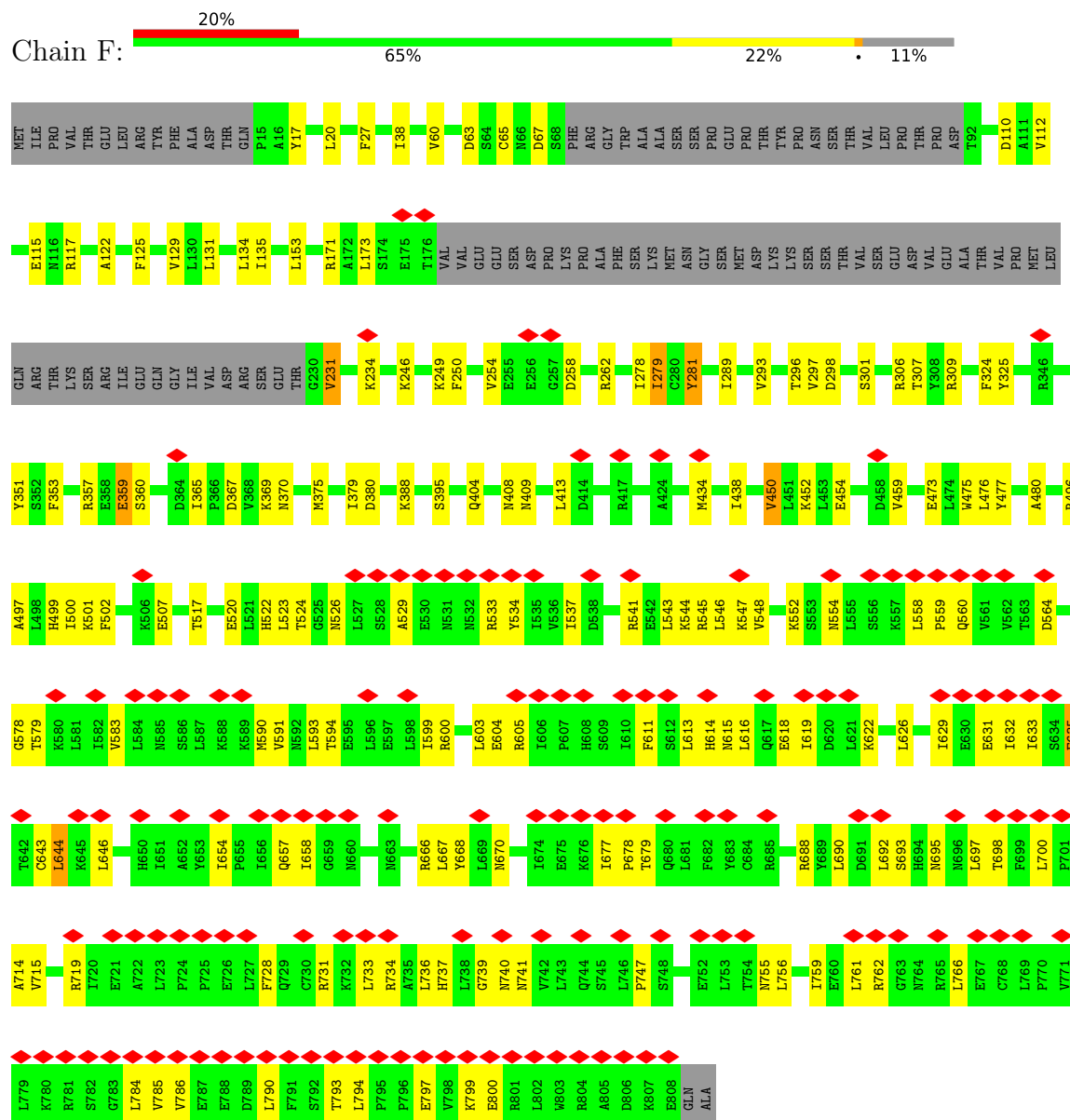


• Molecule 1: Volume-regulated anion channel subunit LRRC8A

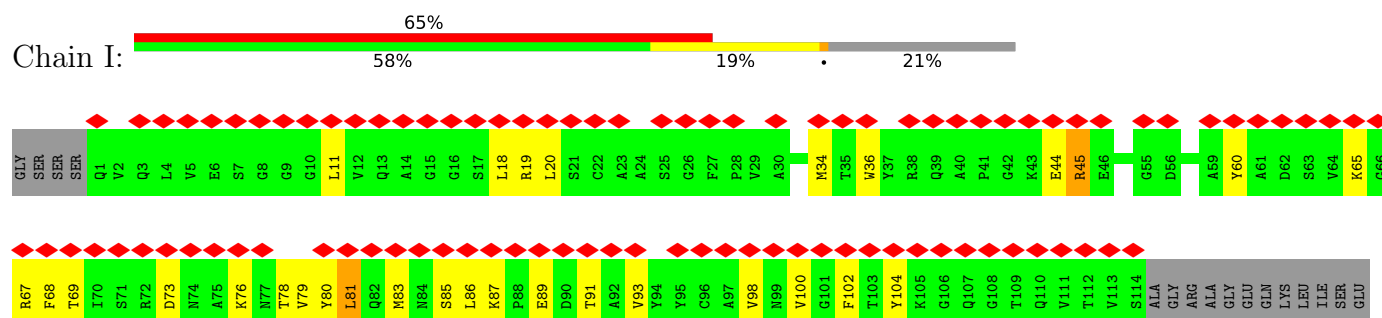




• Molecule 1: Volume-regulated anion channel subunit LRRC8A

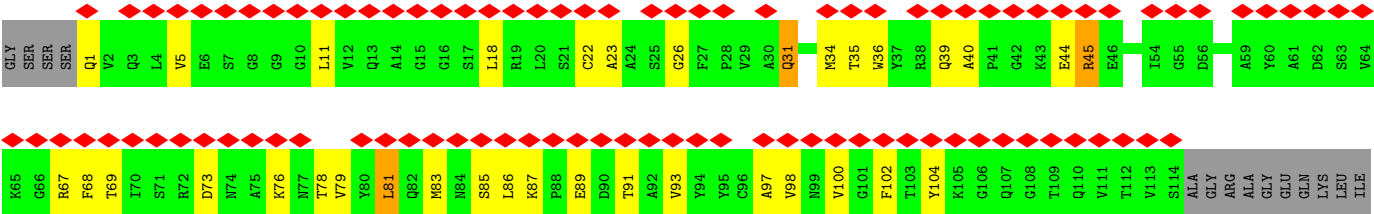


• Molecule 2: synthetic nanobody Sb5



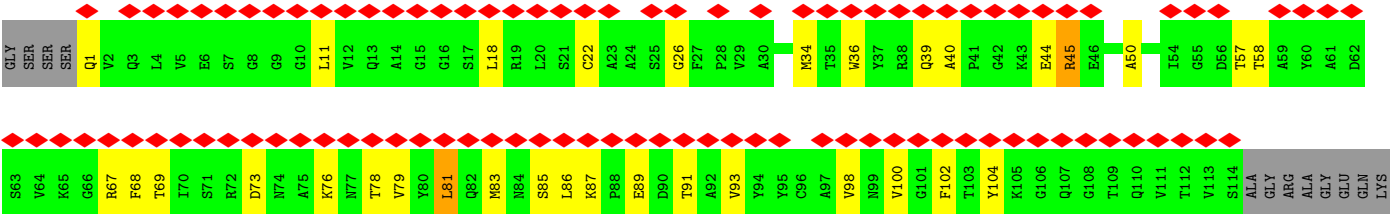
GLU
ASP
LEU
SER
SER
ALA
VAL
ASP
HIS
HIS
HIS
HIS
HIS

● Molecule 2: synthetic nanobody Sb5



SER
GLU
GLU
ASP
LEU
SER
SER
ALA
VAL
ASP
HIS
HIS
HIS
HIS
HIS

● Molecule 2: synthetic nanobody Sb5



LEU
ILE
SER
GLU
GLU
LEU
LEU
ASP
ASN
SER
SER
ALA
VAL
ASP
HIS
HIS
HIS
HIS
HIS

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	48917	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	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.116	Depositor
Minimum map value	-0.049	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.022	Depositor
Map size (Å)	437.47202, 437.47202, 437.47202	wwPDB
Map dimensions	336, 336, 336	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.302, 1.302, 1.302	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	0.10	0/6055	0.27	0/8207
1	B	0.10	0/6055	0.29	0/8207
1	C	0.10	0/6055	0.27	0/8207
1	D	0.10	0/6055	0.28	0/8207
1	E	0.10	0/6055	0.28	0/8207
1	F	0.10	0/6055	0.28	0/8207
2	G	0.10	0/873	0.25	0/1184
2	H	0.09	0/873	0.26	0/1184
2	I	0.09	0/873	0.25	0/1184
All	All	0.10	0/38949	0.27	0/52794

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	5922	0	6081	120	0
1	B	5922	0	6081	115	0
1	C	5922	0	6081	125	0
1	D	5922	0	6081	116	0
1	E	5922	0	6081	144	0
1	F	5922	0	6081	107	0
2	G	858	0	841	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	858	0	841	16	0
2	I	858	0	841	14	0
All	All	38106	0	39009	747	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 10.

All (747) 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:647:TRP:HA	1:E:672:ASN:HD21	1.54	0.72
1:F:298:ASP:HA	1:F:306:ARG:HH12	1.55	0.72
1:E:733:LEU:HD13	1:E:736:LEU:HD21	1.72	0.71
1:E:520:GLU:HG2	1:E:548:VAL:HB	1.72	0.71
1:E:700:LEU:HB3	1:E:724:PRO:HG3	1.73	0.70
1:F:692:LEU:HB2	1:F:715:VAL:HG12	1.74	0.70
1:C:620:ASP:HA	1:C:645:LYS:HB2	1.74	0.69
1:B:298:ASP:HA	1:B:306:ARG:HH12	1.58	0.68
1:C:45:LEU:HG	1:C:314:LEU:HD21	1.76	0.68
1:A:714:ALA:HA	1:A:737:HIS:HB2	1.76	0.68
1:B:452:LYS:HA	1:B:475:TRP:HB2	1.76	0.68
1:D:171:ARG:HH22	1:D:231:VAL:H	1.41	0.68
1:D:632:ILE:HD13	1:D:658:ILE:HA	1.76	0.68
1:B:723:LEU:HB2	1:B:743:LEU:HD11	1.76	0.67
1:E:524:THR:HA	1:E:552:LYS:HB2	1.76	0.67
1:C:520:GLU:HG2	1:C:548:VAL:HB	1.77	0.67
1:C:714:ALA:HA	1:C:737:HIS:HB2	1.77	0.67
1:A:520:GLU:HG2	1:A:548:VAL:HB	1.77	0.67
1:A:632:ILE:HA	1:A:635:PHE:HB2	1.76	0.67
1:D:616:LEU:HD21	1:D:619:ILE:HB	1.77	0.66
1:A:443:PHE:HA	1:A:468:LEU:HD21	1.78	0.66
2:G:67:ARG:NH1	2:G:85:SER:O	2.29	0.66
1:C:779:LEU:HD11	1:C:784:LEU:HB2	1.78	0.66
1:A:262:ARG:HH21	1:A:266:ARG:HH12	1.42	0.66
1:D:400:ASN:OD1	1:D:403:ARG:NH2	2.29	0.66
1:C:473:GLU:HG3	1:C:497:ALA:HB3	1.77	0.66
1:A:708:GLN:NE2	1:A:729:GLN:O	2.27	0.66
1:A:665:GLU:HA	1:A:687:LEU:HA	1.78	0.66
1:B:448:LEU:HB3	1:B:468:LEU:HD11	1.78	0.65
1:F:632:ILE:HA	1:F:635:PHE:HB2	1.78	0.65
1:F:666:ARG:NH1	1:F:668:TYR:OH	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:590:MET:HB3	1:E:593:LEU:HB2	1.79	0.65
1:D:452:LYS:HA	1:D:475:TRP:HB2	1.78	0.65
1:A:639:HIS:O	1:A:663:ASN:ND2	2.30	0.65
1:B:419:ARG:NH1	1:B:431:HIS:O	2.30	0.65
1:D:638:LEU:HD13	1:D:641:LEU:HD13	1.80	0.64
1:C:496:ARG:NH2	1:C:517:THR:OG1	2.30	0.64
1:D:638:LEU:HB3	1:D:641:LEU:HB2	1.79	0.64
1:E:781:ARG:HB3	1:E:806:ASP:HB2	1.80	0.64
1:F:524:THR:HG23	1:F:552:LYS:HD3	1.80	0.64
1:F:529:ALA:H	1:F:533:ARG:HB3	1.62	0.64
1:D:517:THR:HA	1:D:545:ARG:HD3	1.80	0.64
1:F:526:ASN:ND2	1:F:554:ASN:OD1	2.30	0.63
1:F:593:LEU:HB3	1:F:613:LEU:HD11	1.80	0.63
1:E:615:ASN:HA	1:E:640:ARG:HD2	1.80	0.63
1:B:345:ARG:HG3	1:B:346:ARG:HG2	1.80	0.63
1:B:774:GLY:O	1:B:801:ARG:NH2	2.32	0.63
1:E:294:ASP:OD1	1:E:309:ARG:NH1	2.31	0.63
1:C:522:HIS:HA	1:C:550:ARG:HB3	1.79	0.63
1:C:647:TRP:O	1:C:649:ASN:ND2	2.32	0.63
1:D:774:GLY:O	1:D:801:ARG:NH2	2.32	0.63
1:E:301:SER:HB3	1:F:67:ASP:HA	1.80	0.63
1:F:473:GLU:HG2	1:F:497:ALA:HB3	1.80	0.63
1:F:794:LEU:O	1:F:799:LYS:NZ	2.32	0.63
1:A:496:ARG:NH2	1:A:517:THR:OG1	2.32	0.62
1:D:576:ASN:HB3	1:D:579:THR:HB	1.81	0.62
1:A:404:GLN:O	1:A:408:ASN:ND2	2.32	0.62
1:C:262:ARG:HH21	1:C:266:ARG:HH12	1.47	0.62
1:E:696:ASN:N	1:E:696:ASN:OD1	2.30	0.62
1:B:606:ILE:HG23	1:B:610:ILE:HD12	1.81	0.62
2:I:34:MET:HG3	2:I:79:VAL:HG21	1.82	0.62
1:E:401:LYS:HG3	1:E:402:LEU:HD12	1.82	0.62
2:H:67:ARG:HH12	2:H:87:LYS:HG2	1.65	0.62
1:B:576:ASN:HB3	1:B:579:THR:HB	1.80	0.62
1:B:756:LEU:HD11	1:B:759:ILE:HG23	1.81	0.61
1:C:152:LYS:HZ1	1:C:257:GLY:HA3	1.66	0.61
1:F:711:GLN:HA	1:F:733:LEU:HA	1.82	0.61
1:E:670:ASN:O	1:E:672:ASN:ND2	2.33	0.61
1:B:234:LYS:HG3	1:B:409:ASN:HB3	1.81	0.61
1:C:404:GLN:O	1:C:408:ASN:ND2	2.33	0.61
1:E:496:ARG:NH2	1:E:517:THR:OG1	2.33	0.61
1:B:473:GLU:HG2	1:B:497:ALA:HB3	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:262:ARG:HH21	1:E:266:ARG:HH12	1.47	0.61
1:A:452:LYS:HA	1:A:475:TRP:HB2	1.83	0.61
1:E:152:LYS:HZ1	1:E:257:GLY:HA3	1.65	0.61
1:C:611:PHE:HA	1:C:638:LEU:HD11	1.83	0.61
1:E:578:GLY:HA3	1:F:544:LYS:HG3	1.83	0.61
1:C:401:LYS:HG3	1:C:402:LEU:HD12	1.81	0.60
1:E:404:GLN:O	1:E:408:ASN:ND2	2.34	0.60
1:A:774:GLY:O	1:A:801:ARG:NH2	2.30	0.60
1:B:593:LEU:HB3	1:B:613:LEU:HD11	1.82	0.60
1:B:711:GLN:NE2	1:B:732:LYS:O	2.34	0.60
2:G:39:GLN:NE2	2:G:40:ALA:O	2.34	0.60
1:E:632:ILE:HA	1:E:635:PHE:HB2	1.83	0.60
2:I:67:ARG:HH12	2:I:87:LYS:HG2	1.64	0.60
1:D:524:THR:HG23	1:D:552:LYS:HD3	1.82	0.60
1:F:616:LEU:HD21	1:F:619:ILE:HB	1.82	0.60
1:F:632:ILE:HG12	1:F:658:ILE:HA	1.82	0.60
1:B:499:HIS:HA	1:B:522:HIS:HB2	1.82	0.60
1:A:294:ASP:OD1	1:A:309:ARG:NH1	2.35	0.60
1:A:647:TRP:O	1:A:649:ASN:ND2	2.35	0.60
1:F:714:ALA:HA	1:F:737:HIS:HB2	1.84	0.60
1:F:677:ILE:HD11	1:F:697:LEU:HD13	1.84	0.60
2:G:31:GLN:O	2:G:31:GLN:NE2	2.35	0.60
1:B:663:ASN:HA	1:B:686:LYS:HD3	1.83	0.59
1:F:452:LYS:HA	1:F:475:TRP:HB2	1.83	0.59
1:F:688:ARG:HH22	1:F:734:ARG:HH12	1.50	0.59
2:H:67:ARG:NH1	2:H:85:SER:O	2.35	0.59
1:E:663:ASN:HA	1:E:686:LYS:HD2	1.83	0.59
1:A:504:ASP:OD2	1:A:506:LYS:NZ	2.35	0.59
1:A:702:ALA:HA	1:A:726:GLU:HG3	1.83	0.59
1:F:258:ASP:OD2	1:F:262:ARG:NH1	2.34	0.59
1:C:504:ASP:OD2	1:C:506:LYS:NZ	2.32	0.59
1:E:452:LYS:HA	1:E:475:TRP:HB2	1.85	0.59
1:A:632:ILE:HD11	1:A:655:PRO:HB2	1.83	0.58
1:B:700:LEU:HB2	1:B:724:PRO:HG3	1.85	0.58
1:E:676:LYS:NZ	1:E:698:THR:OG1	2.30	0.58
1:F:611:PHE:HZ	1:F:631:GLU:HB3	1.69	0.58
1:A:576:ASN:HB3	1:A:579:THR:HB	1.84	0.58
1:A:719:ARG:HH12	1:B:640:ARG:HA	1.68	0.58
1:B:761:LEU:HB2	1:B:786:VAL:HG23	1.84	0.58
1:C:590:MET:HB3	1:C:593:LEU:HB2	1.84	0.58
1:E:461:ILE:HB	1:E:483:ILE:HG13	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:ALA:HB2	1:A:129:VAL:HG12	1.85	0.58
1:B:639:HIS:O	1:B:663:ASN:ND2	2.33	0.58
1:C:39:ALA:HB2	1:C:129:VAL:HG12	1.85	0.58
1:F:576:ASN:HB3	1:F:579:THR:HB	1.85	0.58
1:F:759:ILE:HD11	1:F:784:LEU:HG	1.84	0.58
1:F:667:LEU:HD23	1:F:690:LEU:HD13	1.85	0.58
1:B:420:LEU:HD13	1:B:428:LEU:HD21	1.86	0.57
1:C:733:LEU:HD13	1:C:736:LEU:HD21	1.86	0.57
1:E:563:THR:HB	1:E:589:LYS:HD2	1.86	0.57
1:A:427:LYS:NZ	1:A:473:GLU:OE2	2.31	0.57
1:A:552:LYS:HG3	1:A:575:ASN:HB3	1.85	0.57
1:B:770:PRO:HG2	1:B:773:LEU:HB2	1.87	0.57
1:A:431:HIS:NE2	1:A:454:GLU:OE1	2.36	0.57
1:D:759:ILE:HD11	1:D:784:LEU:HG	1.86	0.57
1:A:401:LYS:HG3	1:A:402:LEU:HD12	1.86	0.57
1:E:398:SER:HA	1:E:401:LYS:HG2	1.86	0.57
2:H:22:CYS:HB3	2:H:79:VAL:HB	1.86	0.57
1:A:45:LEU:HG	1:A:314:LEU:HD21	1.87	0.57
1:D:241:LYS:HA	1:D:402:LEU:HD21	1.87	0.57
1:F:638:LEU:HD23	1:F:641:LEU:HD13	1.87	0.57
1:D:751:GLY:HA3	1:D:772:GLU:HB3	1.87	0.56
1:E:148:ARG:NH1	1:F:380:ASP:O	2.38	0.56
1:B:454:GLU:HA	1:B:477:TYR:HB2	1.87	0.56
1:E:499:HIS:HA	1:E:522:HIS:HB2	1.87	0.56
1:E:769:LEU:HD11	1:E:794:LEU:HA	1.88	0.56
1:F:591:VAL:O	1:F:615:ASN:ND2	2.38	0.56
1:C:301:SER:HB3	1:D:67:ASP:HA	1.86	0.56
1:C:461:ILE:HB	1:C:483:ILE:HG13	1.88	0.56
1:C:711:GLN:HB3	1:C:734:ARG:HG2	1.88	0.56
1:D:629:ILE:HD13	1:D:651:ILE:HD12	1.86	0.56
1:D:714:ALA:HA	1:D:737:HIS:HB2	1.88	0.56
1:E:233:ASP:HB3	1:E:236:GLU:HG3	1.87	0.56
1:E:492:ARG:HG2	1:E:515:LEU:HA	1.88	0.56
2:H:39:GLN:NE2	2:H:40:ALA:O	2.38	0.56
1:F:654:ILE:HD12	1:F:677:ILE:HG23	1.87	0.56
1:F:756:LEU:HD11	1:F:759:ILE:HG23	1.87	0.56
1:A:711:GLN:HB3	1:A:734:ARG:HG2	1.88	0.56
1:E:169:THR:OG1	1:E:389:ARG:NH2	2.39	0.56
1:F:638:LEU:HB3	1:F:641:LEU:HB2	1.88	0.56
1:A:104:HIS:NE2	1:B:110:ASP:OD2	2.36	0.56
1:A:691:ASP:HA	1:A:714:ALA:HB3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:480:ALA:HB2	1:B:507:GLU:HA	1.88	0.56
1:A:243:LEU:HB3	1:A:393:PHE:HE2	1.70	0.55
1:A:629:ILE:HG21	1:A:646:LEU:HD13	1.88	0.55
1:C:373:ALA:O	1:C:377:HIS:ND1	2.40	0.55
2:H:34:MET:HG3	2:H:79:VAL:HG21	1.88	0.55
1:A:488:LEU:HG	1:A:492:ARG:HE	1.71	0.55
1:B:737:HIS:HA	1:B:760:GLU:HB3	1.88	0.55
1:C:760:GLU:HA	1:C:785:VAL:HB	1.89	0.55
1:D:654:ILE:HD12	1:D:677:ILE:HG23	1.88	0.55
1:E:39:ALA:HB2	1:E:129:VAL:HG12	1.88	0.55
1:E:541:ARG:NH2	1:E:560:GLN:OE1	2.40	0.55
1:E:45:LEU:HG	1:E:314:LEU:HD21	1.88	0.55
1:D:617:GLN:OE1	1:D:640:ARG:NH2	2.34	0.55
1:A:301:SER:HB3	1:B:67:ASP:HA	1.88	0.55
1:F:700:LEU:HB3	1:F:704:ILE:HD13	1.87	0.55
1:D:566:GLY:HA3	1:D:590:MET:HE2	1.89	0.55
1:A:168:TRP:HH2	1:A:398:SER:HB3	1.73	0.54
1:B:554:ASN:ND2	1:B:577:GLU:OE2	2.40	0.54
1:B:438:ILE:HG13	1:B:459:VAL:HG21	1.87	0.54
1:C:243:LEU:HA	1:C:246:LYS:HG2	1.90	0.54
1:C:700:LEU:HB2	1:C:724:PRO:HG3	1.88	0.54
1:D:117:ARG:NH1	1:D:296:THR:O	2.38	0.54
1:E:688:ARG:HA	1:E:710:LEU:HA	1.88	0.54
1:F:480:ALA:HB2	1:F:507:GLU:HA	1.89	0.54
2:I:67:ARG:NH1	2:I:85:SER:O	2.40	0.54
2:G:34:MET:HG3	2:G:79:VAL:HG21	1.88	0.54
1:C:443:PHE:HA	1:C:468:LEU:HD21	1.90	0.54
1:D:298:ASP:HA	1:D:306:ARG:HH12	1.72	0.54
1:C:234:LYS:NZ	1:C:410:GLU:OE1	2.38	0.54
1:A:590:MET:HB3	1:A:593:LEU:HB2	1.88	0.54
1:B:420:LEU:HB3	1:B:428:LEU:HD11	1.90	0.54
1:A:243:LEU:HD23	1:A:246:LYS:HD2	1.88	0.54
1:B:621:LEU:H	1:B:622:LYS:HZ1	1.56	0.54
1:B:368:VAL:HG11	1:B:372:PHE:HD2	1.73	0.54
1:D:618:GLU:HG2	1:D:666:ARG:HH22	1.73	0.54
1:F:541:ARG:HH12	1:F:564:ASP:HB2	1.72	0.54
1:A:663:ASN:HA	1:A:686:LYS:HD2	1.90	0.54
1:C:398:SER:HA	1:C:401:LYS:HG2	1.90	0.54
1:C:691:ASP:HA	1:C:714:ALA:HB3	1.90	0.54
1:F:438:ILE:HG13	1:F:459:VAL:HG21	1.90	0.54
1:C:755:ASN:HA	1:C:778:LEU:HD12	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:492:ARG:HG2	1:A:515:LEU:HA	1.91	0.53
1:E:431:HIS:HA	1:E:452:LYS:HB2	1.90	0.53
1:C:635:PHE:HD1	1:C:638:LEU:HD13	1.72	0.53
1:B:105:GLN:HA	1:C:55:LEU:HD11	1.89	0.53
1:F:790:LEU:O	1:F:793:THR:OG1	2.25	0.53
1:B:740:ASN:N	1:B:762:ARG:O	2.41	0.53
1:F:654:ILE:HB	1:F:678:PRO:HD2	1.90	0.53
1:F:618:GLU:HG3	1:F:643:CYS:HB3	1.90	0.53
2:I:73:ASP:OD1	2:I:73:ASP:N	2.42	0.53
1:B:117:ARG:NH1	1:B:296:THR:O	2.41	0.53
1:D:740:ASN:N	1:D:762:ARG:O	2.33	0.53
1:E:473:GLU:HG3	1:E:497:ALA:HB3	1.91	0.53
1:F:766:LEU:HD23	1:F:790:LEU:HD13	1.91	0.52
1:C:719:ARG:HH22	1:D:640:ARG:HB2	1.73	0.52
1:E:580:LYS:NZ	1:E:603:LEU:O	2.42	0.52
1:E:712:ASN:HA	1:E:735:ALA:HB3	1.90	0.52
1:F:693:SER:O	1:F:695:ASN:ND2	2.42	0.52
1:A:233:ASP:HB3	1:A:236:GLU:HG3	1.92	0.52
1:B:566:GLY:HA3	1:B:590:MET:HE2	1.91	0.52
1:B:791:PHE:O	1:B:799:LYS:NZ	2.29	0.52
1:D:492:ARG:NH1	1:D:514:SER:O	2.43	0.52
1:D:632:ILE:HA	1:D:635:PHE:HD1	1.74	0.52
1:E:714:ALA:HA	1:E:737:HIS:HB2	1.90	0.52
1:A:234:LYS:NZ	1:A:410:GLU:OE1	2.38	0.52
1:A:541:ARG:NH2	1:A:560:GLN:OE1	2.42	0.52
1:D:763:GLY:N	1:D:787:GLU:OE2	2.34	0.52
1:D:770:PRO:HG2	1:D:773:LEU:HB2	1.91	0.52
1:E:665:GLU:HG3	1:E:666:ARG:HG3	1.90	0.52
1:C:463:PRO:HB3	1:C:486:PRO:HB2	1.92	0.52
1:E:416:LEU:HD21	1:E:442:VAL:HA	1.92	0.52
1:A:709:ASN:HA	1:A:732:LYS:HD2	1.91	0.52
1:C:432:LEU:HB3	1:C:435:LEU:HD12	1.92	0.52
1:D:635:PHE:HB3	1:D:661:LEU:HD21	1.92	0.52
1:A:668:TYR:HA	1:A:691:ASP:HB3	1.91	0.52
1:B:656:ILE:HB	1:B:680:GLN:HG3	1.92	0.52
1:A:700:LEU:HB2	1:A:724:PRO:HG3	1.92	0.52
1:B:580:LYS:HA	1:B:602:ASP:HB2	1.92	0.52
1:C:638:LEU:HD22	1:C:641:LEU:HD22	1.92	0.52
1:F:570:GLN:OE1	1:F:594:THR:OG1	2.27	0.52
1:C:233:ASP:HB3	1:C:236:GLU:HG3	1.92	0.51
2:G:1:GLN:HE21	2:G:26:GLY:HA3	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:713:LEU:HB3	1:E:736:LEU:HD23	1.93	0.51
1:F:360:SER:HB2	1:F:388:LYS:HZ2	1.75	0.51
1:E:632:ILE:HD11	1:E:655:PRO:HB2	1.90	0.51
2:I:44:GLU:HG3	2:I:45:ARG:H	1.75	0.51
1:A:463:PRO:HB3	1:A:486:PRO:HB2	1.92	0.51
1:D:463:PRO:HD3	1:D:484:GLU:HG3	1.93	0.51
1:A:67:ASP:HA	1:F:301:SER:HB2	1.92	0.51
1:C:433:PHE:HA	1:C:454:GLU:HG2	1.93	0.51
1:C:781:ARG:NH2	1:C:788:GLU:OE1	2.44	0.51
1:E:620:ASP:HA	1:E:645:LYS:HB2	1.91	0.51
1:E:685:ARG:O	1:E:709:ASN:ND2	2.44	0.51
1:A:416:LEU:HD22	1:A:441:THR:HG23	1.92	0.51
1:B:122:ALA:HB2	1:B:289:ILE:HG22	1.93	0.51
1:C:492:ARG:HG3	1:C:493:GLU:HG2	1.92	0.51
1:A:760:GLU:HA	1:A:785:VAL:HB	1.92	0.51
1:B:38:ILE:HD11	1:B:324:PHE:HD2	1.75	0.51
1:E:543:LEU:HB3	1:E:546:LEU:HB2	1.92	0.51
1:E:787:GLU:OE1	1:E:787:GLU:N	2.41	0.51
1:D:615:ASN:HA	1:D:640:ARG:HD3	1.92	0.51
1:E:421:THR:OG1	1:E:429:GLU:OE2	2.29	0.51
1:C:168:TRP:HH2	1:C:398:SER:HB3	1.76	0.51
1:C:543:LEU:HD12	1:C:546:LEU:HD22	1.92	0.51
1:C:632:ILE:HD11	1:C:655:PRO:HB2	1.93	0.51
1:B:529:ALA:H	1:B:533:ARG:HB3	1.74	0.50
1:E:416:LEU:HD22	1:E:441:THR:HG23	1.91	0.50
1:B:633:ILE:HB	1:B:657:GLN:HG2	1.93	0.50
1:D:450:VAL:HG23	1:D:473:GLU:HB2	1.93	0.50
1:F:365:ILE:HG12	1:F:395:SER:HA	1.94	0.50
1:F:632:ILE:HG23	1:F:657:GLN:HB3	1.94	0.50
1:B:258:ASP:O	1:B:262:ARG:NH1	2.45	0.50
1:C:550:ARG:HD2	1:C:573:SER:HB3	1.93	0.50
1:E:755:ASN:HA	1:E:778:LEU:HD12	1.92	0.50
1:C:492:ARG:HH21	1:C:514:SER:HB2	1.77	0.50
1:C:615:ASN:HA	1:C:640:ARG:HD3	1.94	0.50
1:E:432:LEU:HB3	1:E:435:LEU:HD12	1.93	0.50
2:H:44:GLU:HG3	2:H:45:ARG:H	1.76	0.50
1:C:412:THR:OG1	1:C:413:LEU:N	2.44	0.50
1:E:629:ILE:HG21	1:E:646:LEU:HD13	1.94	0.50
1:A:54:CYS:HA	1:A:310:CYS:HA	1.94	0.50
1:B:501:LYS:HG2	1:B:524:THR:HB	1.93	0.50
1:D:566:GLY:HA2	1:D:569:LEU:HD23	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:476:LEU:HB2	1:D:500:ILE:HD13	1.93	0.50
1:D:618:GLU:HG3	1:D:643:CYS:HB3	1.94	0.50
1:F:520:GLU:HA	1:F:548:VAL:HB	1.94	0.50
1:F:731:ARG:O	1:F:755:ASN:ND2	2.44	0.50
1:A:345:ARG:HG3	1:A:346:ARG:HG2	1.94	0.49
1:C:104:HIS:NE2	1:D:110:ASP:OD2	2.45	0.49
1:E:618:GLU:HG2	1:E:643:CYS:HB3	1.93	0.49
2:H:83:MET:HB3	2:H:86:LEU:HD21	1.94	0.49
2:I:83:MET:HB3	2:I:86:LEU:HD21	1.94	0.49
2:H:73:ASP:N	2:H:73:ASP:OD1	2.44	0.49
1:A:243:LEU:HA	1:A:246:LYS:HG2	1.94	0.49
1:A:258:ASP:N	1:A:371:ASP:OD2	2.46	0.49
1:D:420:LEU:HD13	1:D:428:LEU:HD11	1.95	0.49
1:F:599:ILE:HD12	1:F:622:LYS:HB3	1.94	0.49
1:B:570:GLN:OE1	1:B:594:THR:OG1	2.29	0.49
1:E:443:PHE:HA	1:E:468:LEU:HD21	1.93	0.49
1:B:452:LYS:HG2	1:B:475:TRP:CG	2.47	0.49
1:B:746:LEU:HD23	1:B:766:LEU:HD22	1.93	0.49
1:C:746:LEU:HD23	1:C:770:PRO:HD2	1.94	0.49
1:D:301:SER:OG	1:E:68:SER:N	2.45	0.49
1:E:353:PHE:O	1:E:357:ARG:NH1	2.46	0.49
1:A:698:THR:HG23	1:A:699:PHE:HD1	1.78	0.49
1:C:265:MET:HA	1:C:343:MET:HE1	1.94	0.49
1:D:555:LEU:O	1:D:576:ASN:ND2	2.45	0.49
1:E:488:LEU:HG	1:E:492:ARG:HE	1.77	0.49
1:C:754:THR:HG23	1:C:777:PRO:HG2	1.94	0.49
1:D:656:ILE:HB	1:D:680:GLN:HG3	1.93	0.49
1:F:117:ARG:NH1	1:F:296:THR:O	2.45	0.49
1:D:255:GLU:OE2	1:D:369:LYS:HB3	2.13	0.49
1:D:762:ARG:HH12	1:D:785:VAL:HG12	1.78	0.49
1:F:688:ARG:HA	1:F:710:LEU:HA	1.94	0.49
1:A:555:LEU:HD21	1:A:559:PRO:HD3	1.95	0.48
1:C:61:THR:O	1:C:64:SER:OG	2.28	0.48
1:C:413:LEU:HD13	1:C:444:ASP:HB2	1.94	0.48
1:C:548:VAL:HG22	1:C:571:LYS:HB3	1.95	0.48
1:C:599:ILE:HG23	1:C:622:LYS:HD2	1.93	0.48
1:E:265:MET:HA	1:E:343:MET:HE1	1.95	0.48
1:C:420:LEU:HD11	1:C:445:LEU:HD22	1.95	0.48
1:D:480:ALA:HB2	1:D:507:GLU:HA	1.96	0.48
1:D:688:ARG:HE	1:D:711:GLN:HG3	1.77	0.48
1:E:36:LEU:HB2	1:E:133:THR:HG21	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:431:HIS:HB3	1:E:452:LYS:HD2	1.95	0.48
1:F:246:LYS:HD2	1:F:249:LYS:HD3	1.95	0.48
1:F:761:LEU:HB2	1:F:786:VAL:HG23	1.95	0.48
1:B:622:LYS:HE3	1:B:645:LYS:HB3	1.93	0.48
1:B:665:GLU:O	1:B:688:ARG:N	2.39	0.48
1:C:614:HIS:CE1	1:C:637:HIS:HB3	2.48	0.48
1:E:233:ASP:OD2	1:E:235:LYS:NZ	2.44	0.48
1:E:618:GLU:OE2	1:E:666:ARG:NH2	2.47	0.48
1:B:413:LEU:HG	1:B:417:ARG:HH12	1.78	0.48
1:B:599:ILE:HD13	1:B:622:LYS:HB2	1.95	0.48
1:B:726:GLU:OE1	1:B:726:GLU:N	2.43	0.48
1:C:294:ASP:OD1	1:C:294:ASP:N	2.45	0.48
1:E:104:HIS:NE2	1:F:110:ASP:OD2	2.41	0.48
1:E:711:GLN:HA	1:E:733:LEU:HA	1.95	0.48
1:C:258:ASP:N	1:C:371:ASP:OD2	2.46	0.48
1:D:443:PHE:HA	1:D:468:LEU:HD21	1.96	0.48
2:I:18:LEU:HB2	2:I:83:MET:HE3	1.96	0.48
1:A:129:VAL:HG13	1:A:325:TYR:CE1	2.48	0.48
1:C:36:LEU:HB2	1:C:133:THR:HG21	1.95	0.48
1:F:543:LEU:HB3	1:F:546:LEU:HB2	1.94	0.48
1:A:278:ILE:HA	1:A:281:TYR:CD2	2.49	0.48
1:B:700:LEU:HB3	1:B:704:ILE:HD13	1.95	0.48
1:E:504:ASP:OD2	1:E:506:LYS:NZ	2.32	0.48
1:C:774:GLY:O	1:C:801:ARG:NH2	2.38	0.48
1:E:168:TRP:HH2	1:E:398:SER:HB3	1.78	0.48
1:A:416:LEU:HD11	1:A:430:LEU:HD21	1.96	0.48
1:C:688:ARG:HA	1:C:710:LEU:HA	1.95	0.48
1:D:431:HIS:HD2	2:H:57:THR:HG21	1.79	0.48
1:D:570:GLN:OE1	1:D:594:THR:OG1	2.31	0.48
1:E:129:VAL:HG13	1:E:325:TYR:CE1	2.49	0.48
1:F:679:THR:OG1	1:F:703:ASP:OD2	2.32	0.48
1:B:742:VAL:O	1:B:765:ARG:NE	2.36	0.48
1:D:620:ASP:HB2	1:D:645:LYS:HE3	1.96	0.48
1:F:122:ALA:HB2	1:F:289:ILE:HG22	1.96	0.48
1:A:635:PHE:HD1	1:A:638:LEU:HD13	1.79	0.47
1:B:647:TRP:HA	1:B:672:ASN:HD21	1.79	0.47
1:D:399:GLU:OE1	1:D:399:GLU:N	2.42	0.47
1:A:398:SER:HA	1:A:401:LYS:HG2	1.96	0.47
1:C:129:VAL:HG13	1:C:325:TYR:CE1	2.49	0.47
1:C:577:GLU:OE1	1:D:492:ARG:NH2	2.39	0.47
1:D:104:HIS:NE2	1:E:110:ASP:OD2	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:452:LYS:HE3	1:E:475:TRP:CZ3	2.49	0.47
1:F:112:VAL:HG12	1:F:297:VAL:HG11	1.96	0.47
2:G:44:GLU:HG3	2:G:45:ARG:H	1.79	0.47
1:C:348:LEU:HD22	1:C:377:HIS:CG	2.49	0.47
1:D:582:ILE:HG22	1:D:584:LEU:H	1.79	0.47
1:E:453:LEU:HB2	1:E:476:LEU:HD13	1.97	0.47
1:A:755:ASN:HA	1:A:778:LEU:HD22	1.97	0.47
1:C:772:GLU:O	1:C:775:GLU:HG2	2.14	0.47
1:D:20:LEU:HD13	1:D:153:LEU:HD22	1.97	0.47
1:D:125:PHE:O	1:D:129:VAL:HG12	2.15	0.47
1:A:780:LYS:HD2	1:A:806:ASP:HA	1.97	0.47
1:B:463:PRO:HD3	1:B:484:GLU:HG3	1.96	0.47
1:C:164:PHE:O	1:C:389:ARG:NH1	2.40	0.47
1:C:495:LEU:H	1:C:517:THR:HG21	1.78	0.47
1:C:652:ALA:HA	1:C:673:LYS:HB2	1.96	0.47
1:C:789:ASP:OD1	1:C:789:ASP:N	2.47	0.47
1:E:243:LEU:HA	1:E:246:LYS:HG2	1.96	0.47
1:E:505:ILE:O	1:E:508:ILE:HG22	2.15	0.47
1:E:658:ILE:HG21	1:E:681:LEU:HD12	1.96	0.47
1:F:692:LEU:N	1:F:714:ALA:O	2.36	0.47
1:B:538:ASP:HA	1:B:561:VAL:HG21	1.97	0.47
1:B:642:THR:O	1:B:665:GLU:N	2.39	0.47
1:D:654:ILE:HG21	1:D:681:LEU:HD22	1.95	0.47
1:B:555:LEU:O	1:B:576:ASN:ND2	2.47	0.47
1:C:278:ILE:HA	1:C:281:TYR:CD2	2.50	0.47
1:C:766:LEU:HB3	1:C:790:LEU:HD21	1.95	0.47
1:B:18:ARG:NH1	1:B:382:TYR:OH	2.48	0.47
1:E:370:ASN:C	1:E:370:ASN:HD22	2.21	0.47
1:E:688:ARG:HD2	1:E:711:GLN:HE22	1.80	0.47
1:F:454:GLU:HA	1:F:477:TYR:HB2	1.97	0.47
1:F:578:GLY:H	1:F:600:ARG:HB3	1.79	0.47
2:G:35:THR:N	2:G:97:ALA:O	2.45	0.47
1:B:597:GLU:HG2	1:B:620:ASP:HB2	1.95	0.47
1:B:666:ARG:HG2	1:B:689:TYR:HD1	1.80	0.47
1:D:605:ARG:NH1	1:D:628:THR:OG1	2.48	0.47
1:E:791:PHE:HA	1:E:794:LEU:HD13	1.97	0.47
1:F:546:LEU:HG	1:F:569:LEU:HD11	1.96	0.47
1:C:450:VAL:HG22	1:C:473:GLU:HB2	1.97	0.47
1:C:524:THR:HG22	1:C:552:LYS:HD3	1.96	0.47
1:A:63:ASP:HB3	1:A:296:THR:HG21	1.97	0.46
1:C:709:ASN:HA	1:C:732:LYS:HE3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:416:LEU:HD11	1:E:430:LEU:HD21	1.97	0.46
1:F:476:LEU:HB2	1:F:500:ILE:HD13	1.96	0.46
1:B:647:TRP:O	1:B:649:ASN:ND2	2.47	0.46
1:B:654:ILE:HD12	1:B:677:ILE:HG23	1.97	0.46
1:F:499:HIS:HA	1:F:522:HIS:HB2	1.96	0.46
1:D:351:TYR:HB2	1:D:373:ALA:HB1	1.96	0.46
1:F:698:THR:HG22	1:F:719:ARG:HB2	1.98	0.46
1:A:473:GLU:HB3	1:A:497:ALA:HB3	1.97	0.46
1:A:665:GLU:HB2	1:A:688:ARG:HG2	1.97	0.46
1:B:578:GLY:H	1:B:600:ARG:HB3	1.80	0.46
1:F:670:ASN:HA	1:F:695:ASN:HD21	1.80	0.46
1:A:537:ILE:HG22	1:A:540:LEU:HB2	1.96	0.46
1:B:301:SER:HB2	1:C:67:ASP:HA	1.97	0.46
1:B:443:PHE:HA	1:B:468:LEU:HD23	1.97	0.46
1:C:724:PRO:HG2	1:C:727:LEU:HG	1.96	0.46
1:E:278:ILE:HA	1:E:281:TYR:CD2	2.49	0.46
1:C:341:TRP:CD1	1:C:345:ARG:HE	2.33	0.46
1:C:477:TYR:HA	1:C:501:LYS:HG2	1.97	0.46
1:E:454:GLU:HB3	1:E:477:TYR:HB2	1.97	0.46
2:G:73:ASP:OD1	2:G:73:ASP:N	2.45	0.46
1:D:666:ARG:HB3	1:D:689:TYR:HB2	1.96	0.46
1:A:674:ILE:HG22	1:A:697:LEU:HD21	1.98	0.46
1:B:468:LEU:HD12	1:B:471:LEU:HB2	1.98	0.46
1:D:369:LYS:HG2	1:D:370:ASN:H	1.81	0.46
1:D:375:MET:O	1:D:379:ILE:HG12	2.15	0.46
1:F:20:LEU:HD13	1:F:153:LEU:HD22	1.98	0.46
1:F:605:ARG:HA	1:F:605:ARG:HH11	1.81	0.46
1:A:154:GLU:HA	1:A:157:VAL:HG12	1.98	0.46
1:A:289:ILE:O	1:A:319:LYS:NZ	2.46	0.46
1:E:723:LEU:HD12	1:E:724:PRO:HD2	1.97	0.46
2:G:76:LYS:HG3	2:G:78:THR:HG22	1.97	0.46
1:B:772:GLU:O	1:B:775:GLU:HG2	2.17	0.45
1:E:751:GLY:HA3	1:E:772:GLU:HB3	1.99	0.45
1:F:633:ILE:HG12	1:F:657:GLN:HG2	1.97	0.45
2:I:19:ARG:HB3	2:I:80:TYR:HD2	1.81	0.45
1:C:148:ARG:NH1	1:D:380:ASP:O	2.49	0.45
1:C:554:ASN:O	1:C:554:ASN:ND2	2.49	0.45
1:D:351:TYR:OH	1:D:380:ASP:OD2	2.23	0.45
1:E:477:TYR:HA	1:E:501:LYS:HB3	1.98	0.45
1:B:115:GLU:HG3	1:C:291:PHE:CD1	2.51	0.45
1:D:456:ILE:HG22	1:D:459:VAL:HB	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:475:TRP:HA	1:E:499:HIS:HB2	1.97	0.45
2:I:20:LEU:HB2	2:I:81:LEU:HB3	1.99	0.45
1:A:620:ASP:HA	1:A:645:LYS:HB2	1.97	0.45
1:B:303:THR:OG1	1:B:308:TYR:OH	2.23	0.45
1:C:505:ILE:O	1:C:508:ILE:HG22	2.16	0.45
1:D:118:LEU:HD23	1:D:123:LYS:HG3	1.98	0.45
1:D:685:ARG:HA	1:D:707:LEU:HD12	1.98	0.45
1:F:762:ARG:HH12	1:F:785:VAL:HG12	1.81	0.45
1:B:714:ALA:HA	1:B:737:HIS:HB2	1.99	0.45
1:D:401:LYS:O	1:D:404:GLN:HG3	2.17	0.45
1:A:736:LEU:HD12	1:A:756:LEU:HD11	1.99	0.45
1:C:466:ALA:HA	1:C:490:PHE:CD1	2.51	0.45
1:D:606:ILE:HG23	1:D:610:ILE:HD12	1.98	0.45
1:F:357:ARG:NH1	1:F:367:ASP:OD1	2.40	0.45
2:H:18:LEU:HB2	2:H:83:MET:HE3	1.97	0.45
1:B:399:GLU:OE1	1:B:399:GLU:N	2.41	0.45
1:B:594:THR:HG22	1:B:615:ASN:HB3	1.99	0.45
1:D:791:PHE:O	1:D:799:LYS:NZ	2.39	0.45
1:E:450:VAL:HG22	1:E:473:GLU:HB2	1.99	0.45
2:G:36:TRP:CH2	2:G:81:LEU:HB2	2.52	0.45
1:C:499:HIS:HA	1:C:522:HIS:HB2	1.99	0.45
1:D:473:GLU:HG2	1:D:497:ALA:HB3	1.99	0.45
1:F:38:ILE:HD11	1:F:324:PHE:HD2	1.80	0.45
1:A:499:HIS:HA	1:A:522:HIS:HB2	1.99	0.45
1:C:243:LEU:HB3	1:C:393:PHE:HE2	1.82	0.45
1:C:804:ARG:HH12	1:C:808:GLU:HB3	1.81	0.45
1:E:739:GLY:N	1:E:760:GLU:OE2	2.49	0.45
1:A:233:ASP:OD2	1:A:235:LYS:NZ	2.44	0.45
1:A:460:THR:HG22	1:A:482:LYS:HE2	1.98	0.45
1:C:452:LYS:HA	1:C:475:TRP:HB2	1.99	0.45
1:D:163:CYS:HA	1:D:243:LEU:HD21	1.99	0.45
1:D:232:LEU:HB2	1:D:236:GLU:HG2	1.99	0.45
1:E:643:CYS:SG	1:E:666:ARG:NH1	2.90	0.45
1:E:648:TYR:CE1	1:E:671:ARG:HG3	2.52	0.45
1:A:723:LEU:HB3	1:A:747:PRO:HG2	1.98	0.44
1:C:687:LEU:HD21	1:C:690:LEU:HD13	1.99	0.44
1:E:576:ASN:HB2	1:E:601:CYS:HA	1.99	0.44
1:E:648:TYR:HE1	1:E:671:ARG:HG3	1.82	0.44
1:A:152:LYS:HZ3	1:A:257:GLY:HA3	1.81	0.44
1:B:688:ARG:HE	1:B:711:GLN:HG3	1.83	0.44
1:D:691:ASP:OD1	1:D:691:ASP:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:625:ASN:N	1:D:649:ASN:OD1	2.49	0.44
1:D:642:THR:O	1:D:665:GLU:N	2.34	0.44
1:D:643:CYS:SG	1:D:666:ARG:HB2	2.58	0.44
1:F:63:ASP:HA	1:F:307:THR:HG21	1.99	0.44
1:F:234:LYS:HG2	1:F:409:ASN:HB3	2.00	0.44
1:F:258:ASP:O	1:F:262:ARG:NH1	2.44	0.44
1:A:466:ALA:HA	1:A:490:PHE:CD1	2.53	0.44
1:B:365:ILE:HG12	1:B:395:SER:HA	2.00	0.44
1:E:435:LEU:HB2	1:E:456:ILE:HD12	2.00	0.44
1:E:677:ILE:HB	1:E:701:PRO:HG2	2.00	0.44
1:D:599:ILE:HD12	1:D:622:LYS:HB2	2.00	0.44
1:A:621:LEU:HB2	1:A:646:LEU:HD23	2.00	0.44
1:A:713:LEU:HB3	1:A:736:LEU:HD23	1.99	0.44
1:B:364:ASP:OD1	1:B:364:ASP:N	2.50	0.44
1:C:343:MET:HB3	1:C:343:MET:HE2	1.83	0.44
1:C:454:GLU:HB3	1:C:477:TYR:HB2	1.98	0.44
2:H:1:GLN:HE21	2:H:26:GLY:HA3	1.82	0.44
1:A:302:LEU:HD11	1:B:309:ARG:HG2	2.00	0.44
1:C:513:TYR:OH	1:C:536:VAL:O	2.23	0.44
1:C:645:LYS:HB3	1:C:647:TRP:CZ3	2.53	0.44
1:D:645:LYS:HG2	1:D:668:TYR:CD2	2.53	0.44
1:F:501:LYS:HG2	1:F:524:THR:HB	2.00	0.44
1:A:548:VAL:HG13	1:A:571:LYS:HB3	2.00	0.44
1:B:264:TYR:CD2	1:B:343:MET:HE1	2.53	0.44
1:D:570:GLN:O	1:D:594:THR:OG1	2.35	0.44
1:E:509:PRO:HB2	1:E:512:ILE:HG23	2.00	0.44
1:F:728:PHE:HZ	1:F:747:PRO:HD2	1.83	0.44
1:B:741:ASN:N	1:B:764:ASN:OD1	2.44	0.44
1:C:665:GLU:O	1:C:688:ARG:N	2.34	0.44
1:A:611:PHE:HA	1:A:638:LEU:HD11	2.00	0.43
1:B:558:LEU:HD23	1:B:586:SER:HB3	1.99	0.43
1:C:293:VAL:HG13	1:C:310:CYS:HB2	1.99	0.43
1:D:509:PRO:HB2	1:D:512:ILE:HG23	1.99	0.43
1:E:275:PHE:HA	1:E:278:ILE:HG22	2.00	0.43
1:E:289:ILE:O	1:E:319:LYS:NZ	2.46	0.43
1:E:501:LYS:HA	1:E:524:THR:HB	1.99	0.43
1:A:454:GLU:HA	1:A:477:TYR:HB2	2.00	0.43
1:B:56:PRO:HA	1:B:308:TYR:HA	1.99	0.43
1:B:297:VAL:HG23	1:B:299:ILE:HG12	2.00	0.43
1:E:627:LYS:HG2	1:E:650:HIS:HB2	2.00	0.43
1:F:502:PHE:HB2	1:F:507:GLU:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:PHE:O	1:B:129:VAL:HG12	2.18	0.43
1:C:303:THR:OG1	1:C:308:TYR:OH	2.24	0.43
1:C:421:THR:OG1	1:C:429:GLU:OE1	2.36	0.43
1:D:632:ILE:HA	1:D:635:PHE:CD1	2.52	0.43
1:E:594:THR:O	1:E:617:GLN:N	2.44	0.43
1:E:645:LYS:HB3	1:E:647:TRP:CZ3	2.53	0.43
1:F:643:CYS:SG	1:F:644:LEU:N	2.92	0.43
1:A:275:PHE:HA	1:A:278:ILE:HG22	2.00	0.43
1:A:456:ILE:O	1:A:479:THR:HG23	2.18	0.43
1:A:492:ARG:HD3	1:A:514:SER:HB2	2.00	0.43
1:B:356:ILE:HG21	1:B:388:LYS:HA	2.00	0.43
1:C:682:PHE:CD2	1:C:703:ASP:HB3	2.53	0.43
1:D:342:TRP:CD1	1:D:346:ARG:HH21	2.37	0.43
1:F:171:ARG:NH1	1:F:231:VAL:O	2.39	0.43
1:F:351:TYR:OH	1:F:380:ASP:OD2	2.24	0.43
1:B:235:LYS:HA	1:B:238:GLU:HG2	2.00	0.43
1:C:353:PHE:O	1:C:357:ARG:NH1	2.51	0.43
1:E:654:ILE:HG13	1:E:674:ILE:HD12	2.01	0.43
1:F:125:PHE:O	1:F:129:VAL:HG12	2.19	0.43
1:F:629:ILE:HG21	1:F:646:LEU:HD13	2.01	0.43
1:F:739:GLY:O	1:F:741:ASN:ND2	2.52	0.43
1:C:392:VAL:HG23	1:C:393:PHE:HD1	1.84	0.43
2:G:18:LEU:HB2	2:G:83:MET:HE3	2.01	0.43
1:A:566:GLY:HA2	1:A:569:LEU:HG	1.99	0.43
1:B:534:TYR:CE1	1:B:559:PRO:HG3	2.54	0.43
1:D:692:LEU:HB2	1:D:715:VAL:HG12	2.01	0.43
1:E:608:HIS:HA	1:E:611:PHE:CE2	2.54	0.43
1:E:665:GLU:HA	1:E:687:LEU:HA	2.00	0.43
1:E:756:LEU:HD11	1:E:759:ILE:HG23	2.01	0.43
1:A:152:LYS:NZ	1:A:257:GLY:HA3	2.34	0.43
1:C:423:ASN:ND2	1:C:450:VAL:HG21	2.33	0.43
1:C:724:PRO:HA	1:C:725:PRO:HD3	1.94	0.43
1:C:804:ARG:NH1	1:C:808:GLU:HB3	2.33	0.43
1:D:746:LEU:HD13	1:D:766:LEU:HD13	2.00	0.43
1:D:772:GLU:O	1:D:775:GLU:HG2	2.18	0.43
1:E:636:GLN:HG2	1:E:661:LEU:HD22	2.01	0.43
1:F:450:VAL:HG23	1:F:473:GLU:HB2	2.01	0.43
1:F:558:LEU:HD22	1:F:583:VAL:HG12	2.01	0.43
1:A:746:LEU:HD23	1:A:770:PRO:HD2	2.01	0.43
1:D:122:ALA:HB2	1:D:289:ILE:HG22	2.01	0.43
1:D:633:ILE:HG12	1:D:657:GLN:HG2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:517:THR:HA	1:F:545:ARG:HD3	2.00	0.43
1:A:687:LEU:HD21	1:A:690:LEU:HD13	1.99	0.43
1:A:796:PRO:HA	1:A:799:LYS:HD3	2.00	0.43
1:C:598:LEU:HB3	1:C:601:CYS:SG	2.59	0.43
1:A:461:ILE:HB	1:A:483:ILE:HG13	2.00	0.42
1:A:566:GLY:HA3	1:A:590:MET:HG3	2.01	0.42
1:C:654:ILE:HG13	1:C:674:ILE:HD12	2.01	0.42
1:D:264:TYR:CD2	1:D:343:MET:HE1	2.54	0.42
1:D:548:VAL:HG13	1:D:571:LYS:HD3	2.01	0.42
1:E:121:PHE:HD2	1:E:289:ILE:HD11	1.84	0.42
1:E:548:VAL:HG22	1:E:571:LYS:HB3	2.00	0.42
1:A:168:TRP:CE2	1:A:402:LEU:HD22	2.54	0.42
1:A:291:PHE:CD1	1:F:115:GLU:HG3	2.54	0.42
1:A:804:ARG:HH12	1:A:808:GLU:HB3	1.84	0.42
1:B:420:LEU:HD23	1:B:430:LEU:HD13	2.02	0.42
1:B:759:ILE:HG13	1:B:784:LEU:HA	2.02	0.42
1:E:408:ASN:OD1	1:E:440:ASP:N	2.51	0.42
1:A:555:LEU:O	1:A:576:ASN:ND2	2.53	0.42
1:A:688:ARG:NH1	1:A:709:ASN:HD22	2.17	0.42
1:A:688:ARG:HD2	1:A:711:GLN:HE22	1.84	0.42
1:A:772:GLU:O	1:A:775:GLU:HG2	2.18	0.42
1:B:350:LYS:HE3	1:B:350:LYS:HB3	1.90	0.42
1:B:711:GLN:O	1:B:734:ARG:N	2.52	0.42
1:B:762:ARG:HH12	1:B:785:VAL:HG12	1.84	0.42
1:C:500:ILE:HG12	1:C:508:ILE:HG13	2.01	0.42
1:D:47:VAL:HG13	1:E:45:LEU:HD13	2.01	0.42
1:D:255:GLU:CD	1:D:369:LYS:HB3	2.44	0.42
1:E:45:LEU:HD21	1:E:314:LEU:HD11	2.01	0.42
1:E:500:ILE:HG21	1:E:508:ILE:HG13	2.01	0.42
1:E:516:LYS:HE3	1:E:544:LYS:H	1.84	0.42
1:E:575:ASN:HA	1:E:599:ILE:O	2.19	0.42
1:E:794:LEU:HD23	1:E:798:VAL:HG12	2.02	0.42
2:G:67:ARG:HH12	2:G:87:LYS:HG2	1.84	0.42
1:A:373:ALA:O	1:A:377:HIS:ND1	2.52	0.42
1:A:582:ILE:HG23	1:A:584:LEU:HG	2.01	0.42
1:B:357:ARG:NE	1:B:363:SER:O	2.39	0.42
1:B:736:LEU:HB2	1:B:759:ILE:HG22	2.02	0.42
1:C:420:LEU:HD23	1:C:430:LEU:HD13	2.02	0.42
1:E:154:GLU:HA	1:E:157:VAL:HG12	2.01	0.42
1:E:673:LYS:O	1:E:696:ASN:ND2	2.52	0.42
1:F:797:GLU:HA	1:F:800:GLU:HG2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:711:GLN:HA	1:A:733:LEU:HA	2.01	0.42
1:E:56:PRO:HG2	1:E:99:TYR:CD1	2.54	0.42
1:E:651:ILE:HG22	1:E:674:ILE:HD11	2.02	0.42
1:F:279:ILE:HD13	1:F:279:ILE:HA	1.85	0.42
2:H:36:TRP:CH2	2:H:81:LEU:HB2	2.55	0.42
1:A:513:TYR:OH	1:A:536:VAL:O	2.27	0.42
1:E:152:LYS:NZ	1:E:257:GLY:HA3	2.32	0.42
1:F:129:VAL:HG23	1:F:325:TYR:CE1	2.55	0.42
1:F:360:SER:HB2	1:F:388:LYS:NZ	2.34	0.42
1:F:547:LYS:O	1:F:570:GLN:HG3	2.19	0.42
2:I:76:LYS:HG3	2:I:78:THR:HG22	2.02	0.42
2:H:102:PHE:HB2	2:H:104:TYR:CZ	2.55	0.42
1:D:63:ASP:HA	1:D:307:THR:HG21	2.02	0.42
1:F:772:GLU:O	1:F:775:GLU:HG2	2.19	0.42
1:A:117:ARG:HG3	1:A:295:CYS:HA	2.02	0.42
1:A:578:GLY:H	1:A:600:ARG:HB3	1.83	0.42
1:B:167:PRO:O	1:B:170:THR:OG1	2.35	0.42
1:D:616:LEU:HB3	1:D:638:LEU:HD21	2.00	0.42
1:E:709:ASN:HA	1:E:732:LYS:HE2	2.00	0.42
2:G:22:CYS:HB2	2:G:36:TRP:CH2	2.55	0.42
1:A:22:PRO:HD2	1:A:25:ASP:HB2	2.01	0.42
1:A:36:LEU:HB2	1:A:133:THR:HG21	2.01	0.42
1:A:45:LEU:HD21	1:A:314:LEU:HD11	2.00	0.42
1:A:679:THR:OG1	1:A:703:ASP:OD2	2.36	0.42
1:B:563:THR:HA	1:B:590:MET:HE3	2.01	0.42
1:C:353:PHE:HD1	1:C:353:PHE:HA	1.80	0.42
1:E:168:TRP:CE2	1:E:402:LEU:HD22	2.55	0.42
2:G:83:MET:HB3	2:G:86:LEU:HD11	2.02	0.42
2:H:50:ALA:O	2:H:58:THR:OG1	2.28	0.42
1:B:570:GLN:HA	1:B:593:LEU:HA	2.01	0.42
1:C:233:ASP:OD2	1:C:235:LYS:NZ	2.53	0.42
1:C:502:PHE:CZ	1:C:527:LEU:HB3	2.55	0.42
1:D:736:LEU:O	1:D:759:ILE:HA	2.20	0.42
1:E:291:PHE:O	1:E:312:HIS:N	2.38	0.42
1:E:466:ALA:HA	1:E:490:PHE:CD1	2.55	0.42
1:F:534:TYR:CE1	1:F:559:PRO:HG3	2.54	0.42
1:A:543:LEU:HD12	1:A:546:LEU:HD22	2.01	0.41
1:B:142:PHE:HB3	1:B:267:GLN:OE1	2.19	0.41
1:B:548:VAL:HG13	1:B:571:LYS:HD3	2.02	0.41
1:D:567:VAL:O	1:D:592:ASN:ND2	2.44	0.41
1:D:583:VAL:HG23	1:D:585:ASN:H	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:723:LEU:HD12	1:D:724:PRO:HD2	2.02	0.41
1:E:702:ALA:O	1:E:726:GLU:HG3	2.20	0.41
1:E:804:ARG:NH1	1:E:808:GLU:HB3	2.35	0.41
1:F:375:MET:O	1:F:379:ILE:HG12	2.20	0.41
1:A:446:VAL:HG12	1:A:468:LEU:HD23	2.00	0.41
1:B:129:VAL:HG23	1:B:325:TYR:CE1	2.55	0.41
1:C:275:PHE:HA	1:C:278:ILE:HG22	2.02	0.41
1:C:567:VAL:O	1:C:592:ASN:ND2	2.53	0.41
1:D:146:PHE:HD1	1:D:147:PRO:HD2	1.85	0.41
1:F:278:ILE:HA	1:F:281:TYR:HD1	1.86	0.41
1:F:359:GLU:OE2	1:F:388:LYS:NZ	2.52	0.41
1:A:502:PHE:HZ	1:A:527:LEU:HD23	1.85	0.41
1:D:115:GLU:HG3	1:E:291:PHE:CD1	2.55	0.41
1:F:369:LYS:HG2	1:F:370:ASN:H	1.85	0.41
2:I:60:TYR:HE1	2:I:65:LYS:HA	1.85	0.41
1:A:349:LYS:HD2	1:A:370:ASN:HB3	2.02	0.41
1:A:632:ILE:HG13	1:A:657:GLN:HB2	2.01	0.41
1:B:547:LYS:O	1:B:570:GLN:HG3	2.20	0.41
1:D:54:CYS:HA	1:D:310:CYS:HA	2.01	0.41
1:D:578:GLY:H	1:D:600:ARG:HB3	1.85	0.41
1:E:293:VAL:HG13	1:E:310:CYS:HB2	2.00	0.41
1:A:614:HIS:NE2	1:A:637:HIS:HB3	2.36	0.41
1:D:258:ASP:N	1:D:371:ASP:OD2	2.50	0.41
1:B:492:ARG:HH12	1:B:516:LYS:HE2	1.86	0.41
1:B:604:GLU:O	1:B:605:ARG:NH1	2.54	0.41
1:B:607:PRO:HG2	1:B:610:ILE:HG13	2.03	0.41
1:F:688:ARG:NH2	1:F:734:ARG:HH12	2.16	0.41
1:D:105:GLN:HA	1:E:55:LEU:HD11	2.02	0.41
1:D:604:GLU:O	1:D:605:ARG:NE	2.30	0.41
1:E:495:LEU:H	1:E:517:THR:HG21	1.85	0.41
1:E:502:PHE:HZ	1:E:527:LEU:HD23	1.85	0.41
1:E:603:LEU:C	1:E:605:ARG:H	2.29	0.41
2:I:36:TRP:CH2	2:I:81:LEU:HB2	2.56	0.41
2:I:83:MET:HB3	2:I:86:LEU:HD11	2.03	0.41
1:A:789:ASP:OD1	1:A:789:ASP:N	2.53	0.41
1:C:156:PHE:O	1:C:159:ILE:HG13	2.21	0.41
1:C:582:ILE:HG23	1:C:584:LEU:HG	2.02	0.41
1:D:502:PHE:HB2	1:D:507:GLU:HB2	2.02	0.41
1:D:626:LEU:HB2	1:D:649:ASN:ND2	2.36	0.41
1:D:780:LYS:HE3	1:D:780:LYS:HB2	1.95	0.41
1:E:54:CYS:HA	1:E:310:CYS:HA	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:142:PHE:HE1	1:F:27:PHE:HZ	1.68	0.41
1:E:348:LEU:HD22	1:E:377:HIS:CG	2.56	0.41
1:F:404:GLN:HE22	1:F:408:ASN:HD21	1.68	0.41
1:F:500:ILE:HB	1:F:523:LEU:HD23	2.03	0.41
2:G:5:VAL:HG12	2:G:23:ALA:HB3	2.03	0.41
2:G:83:MET:HB3	2:G:86:LEU:HD21	2.01	0.41
1:A:588:LYS:HB2	1:A:612:SER:HB3	2.02	0.41
1:B:343:MET:HB2	1:B:343:MET:HE3	1.84	0.41
1:B:466:ALA:HA	1:B:490:PHE:CG	2.56	0.41
1:B:723:LEU:HD23	1:B:747:PRO:HD2	2.03	0.41
1:C:152:LYS:NZ	1:C:257:GLY:HA3	2.34	0.41
1:C:262:ARG:NH2	1:C:266:ARG:HH22	2.19	0.41
1:C:618:GLU:HG2	1:C:643:CYS:HB3	2.02	0.41
1:C:668:TYR:HA	1:C:691:ASP:HB3	2.03	0.41
1:C:719:ARG:NH2	1:D:640:ARG:HB2	2.35	0.41
1:D:312:HIS:HB3	1:D:315:ALA:HB2	2.02	0.41
1:D:433:PHE:HD1	1:D:454:GLU:HB3	1.86	0.41
1:D:647:TRP:H	1:D:647:TRP:CD1	2.39	0.41
1:D:776:CYS:HB2	1:D:779:LEU:HB3	2.03	0.41
1:E:233:ASP:OD1	1:E:234:LYS:N	2.54	0.41
1:E:548:VAL:HG13	1:E:571:LYS:HB3	2.03	0.41
1:E:654:ILE:HB	1:E:678:PRO:HD3	2.03	0.41
1:E:669:LEU:HB2	1:E:692:LEU:HD23	2.03	0.41
2:H:76:LYS:HG3	2:H:78:THR:HG22	2.01	0.41
1:B:112:VAL:HG12	1:B:297:VAL:HG11	2.02	0.41
1:C:236:GLU:O	1:C:239:GLN:HG3	2.21	0.41
1:F:740:ASN:N	1:F:762:ARG:O	2.52	0.41
1:B:278:ILE:HA	1:B:281:TYR:HD1	1.86	0.40
1:C:669:LEU:HB2	1:C:692:LEU:HD23	2.03	0.40
1:D:461:ILE:HB	1:D:483:ILE:HA	2.03	0.40
1:E:262:ARG:NH2	1:E:266:ARG:HH22	2.19	0.40
1:E:709:ASN:O	1:E:711:GLN:NE2	2.54	0.40
1:F:131:LEU:O	1:F:135:ILE:HG13	2.22	0.40
1:F:616:LEU:HD23	1:F:638:LEU:HD21	2.03	0.40
1:B:357:ARG:HG3	1:B:362:TYR:O	2.21	0.40
1:C:147:PRO:HA	1:C:150:SER:HB3	2.03	0.40
1:C:708:GLN:NE2	1:C:729:GLN:O	2.45	0.40
1:E:478:HIS:CD2	1:E:503:THR:HG23	2.55	0.40
1:F:554:ASN:ND2	1:F:576:ASN:O	2.54	0.40
2:H:22:CYS:HB2	2:H:36:TRP:CH2	2.56	0.40
1:A:433:PHE:CE2	1:B:467:GLN:HB2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:611:PHE:CZ	1:A:631:GLU:HB3	2.56	0.40
1:D:244:PHE:CD2	1:D:402:LEU:HD22	2.56	0.40
1:E:369:LYS:HE2	1:E:369:LYS:HB3	1.89	0.40
1:E:765:ARG:H	1:E:765:ARG:HG2	1.70	0.40
1:A:19:ILE:HG13	1:A:382:TYR:CD2	2.57	0.40
1:B:291:PHE:CZ	1:B:311:ALA:HB1	2.57	0.40
1:C:431:HIS:HA	1:C:452:LYS:HG3	2.03	0.40
1:C:433:PHE:HD2	1:C:454:GLU:HG3	1.86	0.40
1:E:259:ILE:HD12	1:E:259:ILE:HA	1.95	0.40
1:F:250:PHE:CZ	1:F:254:VAL:HG21	2.56	0.40
1:F:614:HIS:HB3	1:F:639:HIS:CE1	2.56	0.40
2:I:102:PHE:HB2	2:I:104:TYR:CZ	2.56	0.40
1:A:548:VAL:HG22	1:A:571:LYS:HB3	2.03	0.40
1:B:578:GLY:N	1:B:600:ARG:HB3	2.36	0.40
1:B:580:LYS:HD3	1:B:581:LEU:N	2.36	0.40
1:D:417:ARG:HA	1:D:420:LEU:HG	2.04	0.40
1:D:588:LYS:HE2	1:D:588:LYS:HB3	1.94	0.40
1:E:686:LYS:HA	1:E:709:ASN:HD22	1.86	0.40
2:G:102:PHE:HB2	2:G:104:TYR:CZ	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	712/810 (88%)	692 (97%)	20 (3%)	0	100	100
1	B	712/810 (88%)	683 (96%)	29 (4%)	0	100	100
1	C	712/810 (88%)	692 (97%)	20 (3%)	0	100	100
1	D	712/810 (88%)	689 (97%)	23 (3%)	0	100	100
1	E	712/810 (88%)	692 (97%)	20 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	712/810 (88%)	685 (96%)	27 (4%)	0	100	100
2	G	112/144 (78%)	102 (91%)	10 (9%)	0	100	100
2	H	112/144 (78%)	103 (92%)	9 (8%)	0	100	100
2	I	112/144 (78%)	103 (92%)	9 (8%)	0	100	100
All	All	4608/5292 (87%)	4441 (96%)	167 (4%)	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	666/749 (89%)	643 (96%)	23 (4%)	32	54
1	B	666/749 (89%)	642 (96%)	24 (4%)	31	54
1	C	666/749 (89%)	648 (97%)	18 (3%)	39	59
1	D	666/749 (89%)	644 (97%)	22 (3%)	33	56
1	E	666/749 (89%)	643 (96%)	23 (4%)	32	54
1	F	666/749 (89%)	641 (96%)	25 (4%)	29	53
2	G	90/114 (79%)	79 (88%)	11 (12%)	5	21
2	H	90/114 (79%)	80 (89%)	10 (11%)	6	24
2	I	90/114 (79%)	80 (89%)	10 (11%)	6	24
All	All	4266/4836 (88%)	4100 (96%)	166 (4%)	30	52

All (166) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	VAL
1	A	65	CYS
1	A	97	ILE
1	A	104	HIS
1	A	231	VAL

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Mol	Chain	Res	Type
1	A	245	GLU
1	A	274	LYS
1	A	278	ILE
1	A	297	VAL
1	A	353	PHE
1	A	375	MET
1	A	402	LEU
1	A	456	ILE
1	A	473	GLU
1	A	500	ILE
1	A	583	VAL
1	A	632	ILE
1	A	635	PHE
1	A	636	GLN
1	A	729	GLN
1	A	743	LEU
1	A	756	LEU
1	A	786	VAL
1	B	17	TYR
1	B	49	GLN
1	B	60	VAL
1	B	61	THR
1	B	65	CYS
1	B	104	HIS
1	B	173	LEU
1	B	279	ILE
1	B	281	TYR
1	B	293	VAL
1	B	353	PHE
1	B	368	VAL
1	B	369	LYS
1	B	380	ASP
1	B	468	LEU
1	B	527	LEU
1	B	551	LEU
1	B	584	LEU
1	B	586	SER
1	B	599	ILE
1	B	622	LYS
1	B	626	LEU
1	B	743	LEU
1	B	759	ILE

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Mol	Chain	Res	Type
1	C	60	VAL
1	C	104	HIS
1	C	231	VAL
1	C	274	LYS
1	C	278	ILE
1	C	293	VAL
1	C	294	ASP
1	C	329	VAL
1	C	344	LEU
1	C	353	PHE
1	C	412	THR
1	C	524	THR
1	C	555	LEU
1	C	583	VAL
1	C	614	HIS
1	C	616	LEU
1	C	657	GLN
1	C	743	LEU
1	D	17	TYR
1	D	60	VAL
1	D	65	CYS
1	D	104	HIS
1	D	134	LEU
1	D	173	LEU
1	D	279	ILE
1	D	281	TYR
1	D	293	VAL
1	D	300	GLU
1	D	353	PHE
1	D	356	ILE
1	D	413	LEU
1	D	450	VAL
1	D	459	VAL
1	D	515	LEU
1	D	520	GLU
1	D	535	ILE
1	D	553	SER
1	D	644	LEU
1	D	743	LEU
1	D	762	ARG
1	E	60	VAL
1	E	65	CYS

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Mol	Chain	Res	Type
1	E	104	HIS
1	E	231	VAL
1	E	238	GLU
1	E	274	LYS
1	E	278	ILE
1	E	293	VAL
1	E	296	THR
1	E	297	VAL
1	E	318	PHE
1	E	344	LEU
1	E	358	GLU
1	E	370	ASN
1	E	402	LEU
1	E	434	MET
1	E	527	LEU
1	E	632	ILE
1	E	636	GLN
1	E	657	GLN
1	E	696	ASN
1	E	738	LEU
1	E	769	LEU
1	F	17	TYR
1	F	60	VAL
1	F	65	CYS
1	F	134	LEU
1	F	173	LEU
1	F	231	VAL
1	F	279	ILE
1	F	281	TYR
1	F	293	VAL
1	F	309	ARG
1	F	353	PHE
1	F	359	GLU
1	F	413	LEU
1	F	434	MET
1	F	450	VAL
1	F	496	ARG
1	F	537	ILE
1	F	560	GLN
1	F	590	MET
1	F	603	LEU
1	F	604	GLU

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Mol	Chain	Res	Type
1	F	626	LEU
1	F	635	PHE
1	F	644	LEU
1	F	736	LEU
2	I	11	LEU
2	I	45	ARG
2	I	68	PHE
2	I	69	THR
2	I	81	LEU
2	I	89	GLU
2	I	91	THR
2	I	93	VAL
2	I	98	VAL
2	I	100	VAL
2	G	11	LEU
2	G	31	GLN
2	G	45	ARG
2	G	68	PHE
2	G	69	THR
2	G	81	LEU
2	G	89	GLU
2	G	91	THR
2	G	93	VAL
2	G	98	VAL
2	G	100	VAL
2	H	11	LEU
2	H	45	ARG
2	H	68	PHE
2	H	69	THR
2	H	81	LEU
2	H	89	GLU
2	H	91	THR
2	H	93	VAL
2	H	98	VAL
2	H	100	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	ASN
1	A	267	GLN
1	A	576	ASN

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Mol	Chain	Res	Type
1	B	267	GLN
1	B	370	ASN
1	B	406	ASN
1	B	592	ASN
1	B	615	ASN
1	B	624	ASN
1	B	636	GLN
1	B	709	ASN
1	C	494	ASN
1	C	660	ASN
1	C	696	ASN
1	D	253	HIS
1	D	267	GLN
1	D	431	HIS
1	D	467	GLN
1	D	494	ASN
1	D	663	ASN
1	D	695	ASN
1	D	744	GLN
1	D	755	ASN
1	E	370	ASN
1	E	672	ASN
1	E	709	ASN
1	E	755	ASN
1	F	267	GLN
1	F	404	GLN
1	F	708	GLN
2	I	82	GLN
2	G	1	GLN
2	G	3	GLN
2	G	99	ASN
2	G	110	GLN
2	H	3	GLN
2	H	82	GLN
2	H	99	ASN

5.3.3 RNA ⓘ

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.

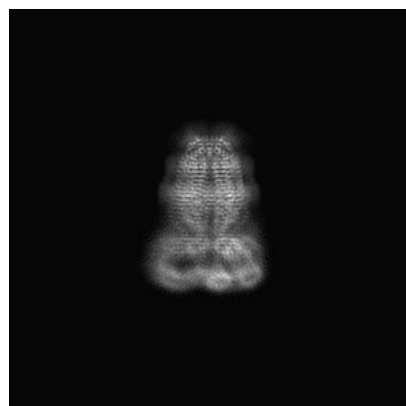
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-13230. These allow visual inspection of the internal detail of the map and identification of artifacts.

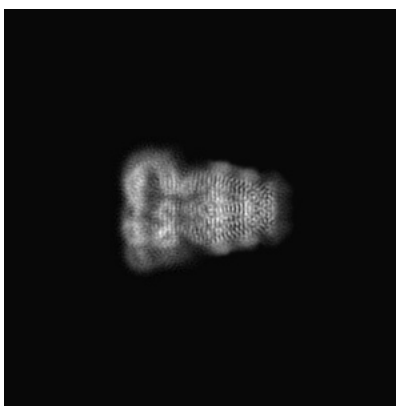
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

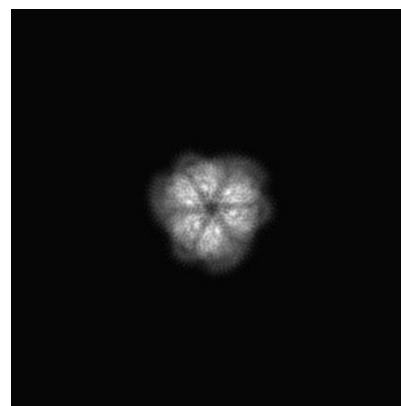
6.1.1 Primary map



X

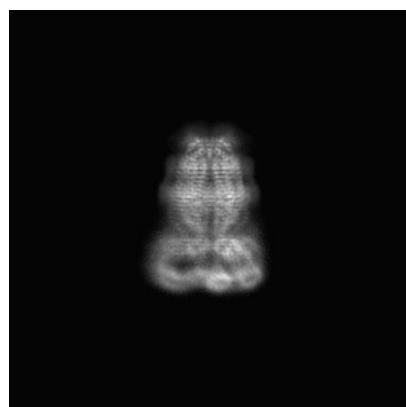


Y

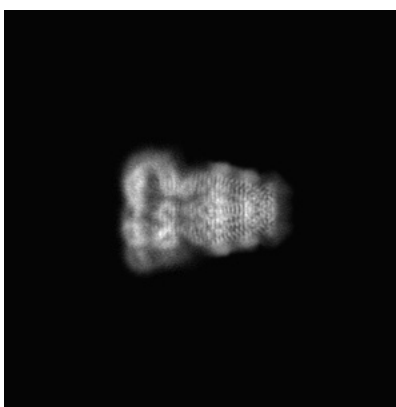


Z

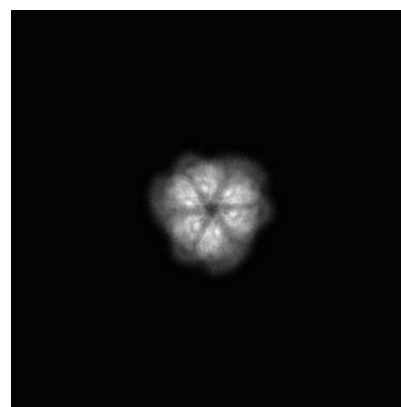
6.1.2 Raw map



X



Y

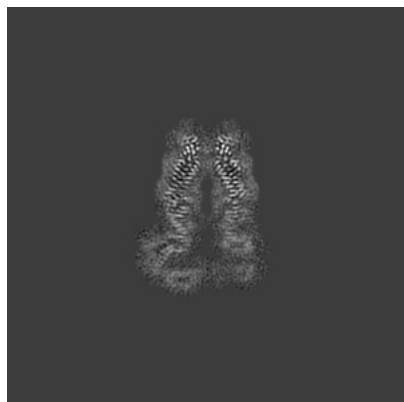


Z

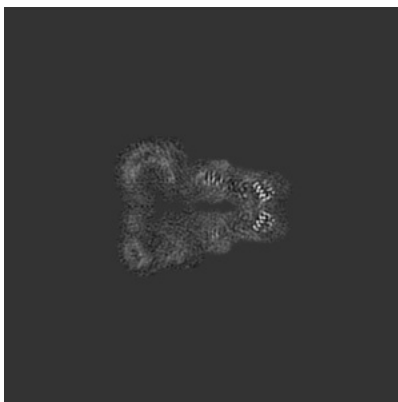
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

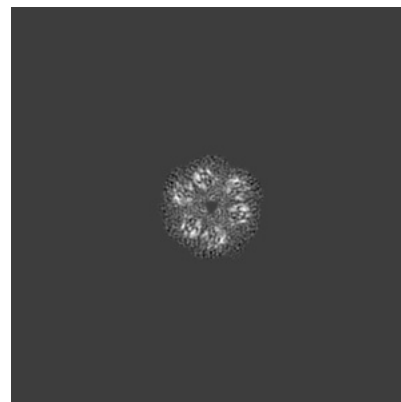
6.2.1 Primary map



X Index: 168

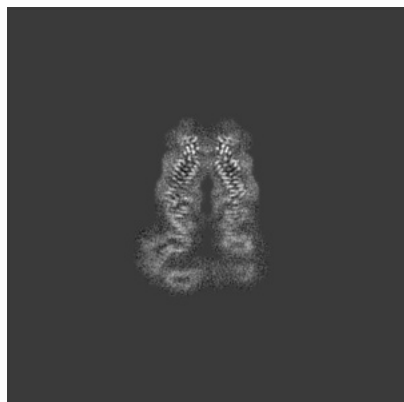


Y Index: 168

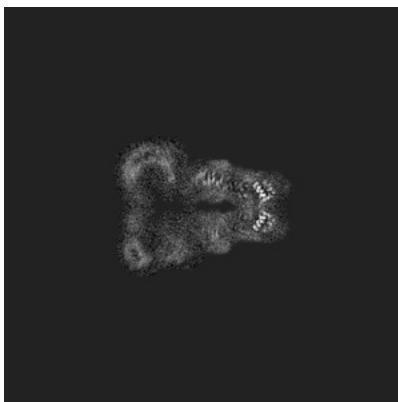


Z Index: 168

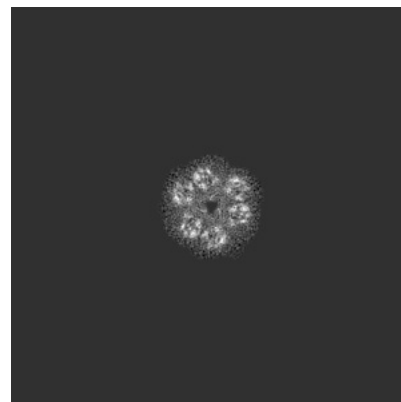
6.2.2 Raw map



X Index: 168



Y Index: 168



Z Index: 168

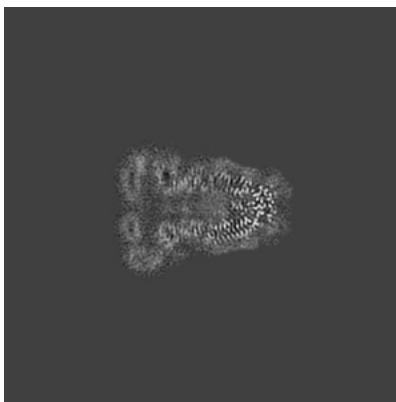
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

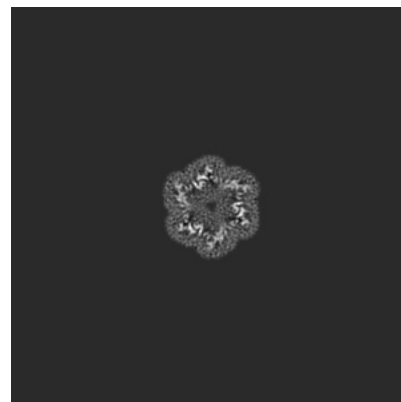
6.3.1 Primary map



X Index: 167

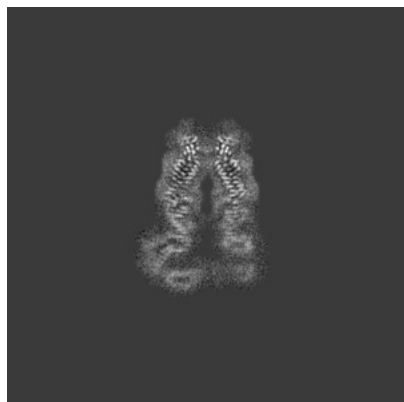


Y Index: 180

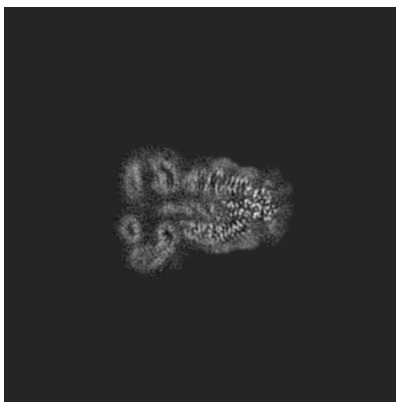


Z Index: 179

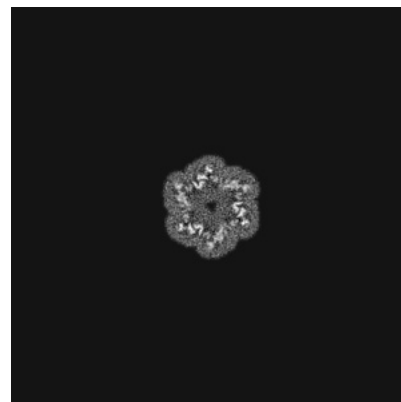
6.3.2 Raw map



X Index: 168



Y Index: 183

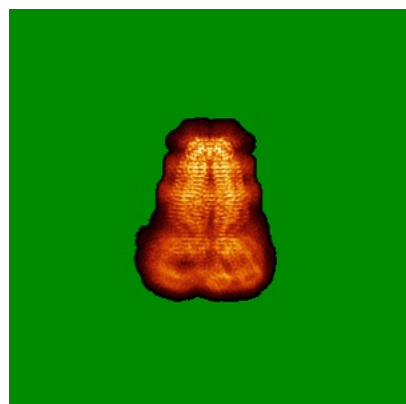


Z Index: 179

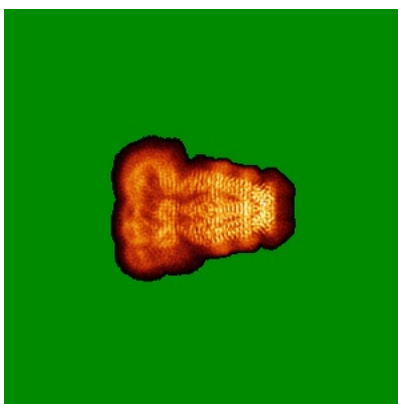
The images above show the largest variance slices of the map in three orthogonal directions.

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

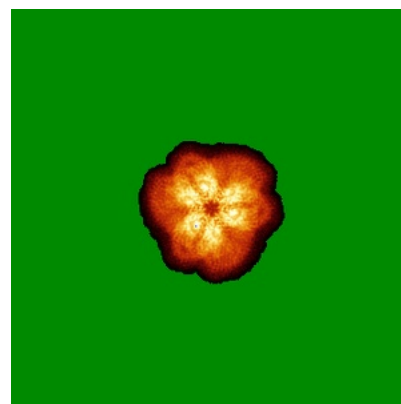
6.4.1 Primary map



X

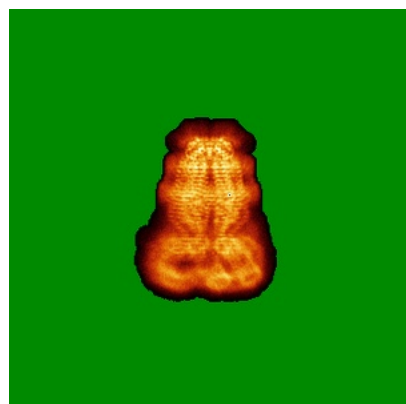


Y

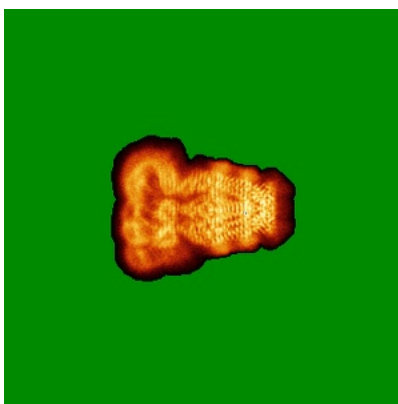


Z

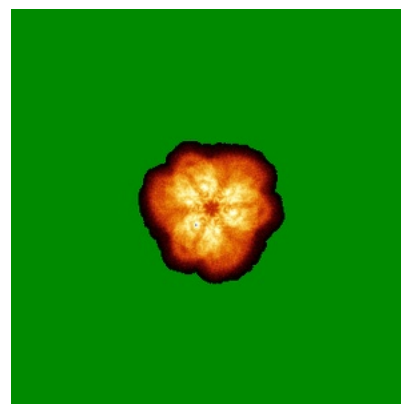
6.4.2 Raw map



X



Y

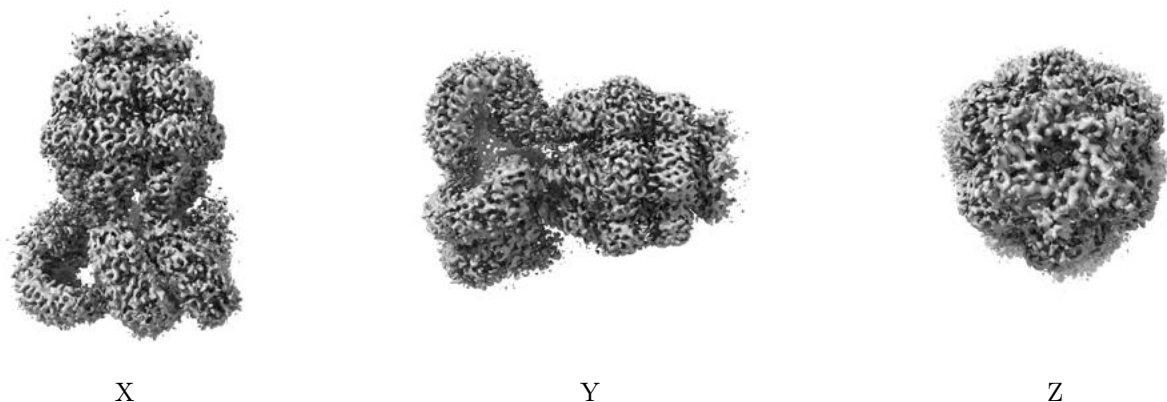


Z

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.

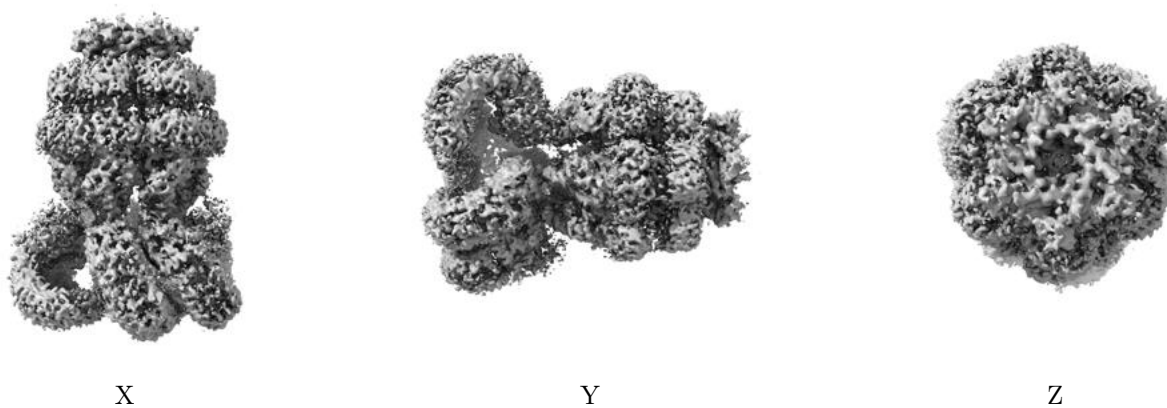
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.022. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

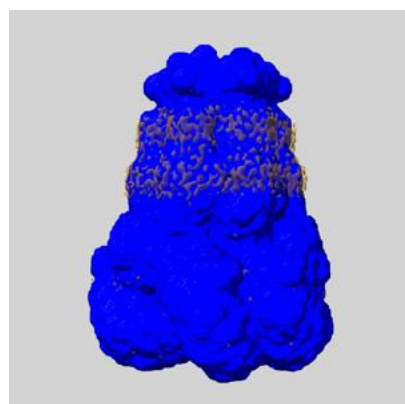
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

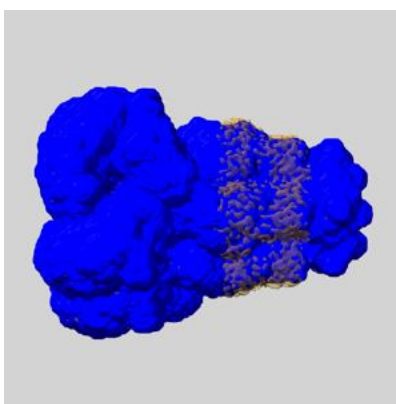
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

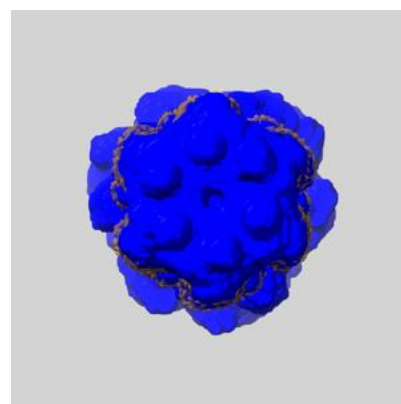
6.6.1 emd_13230_msk_1.map [i](#)



X



Y

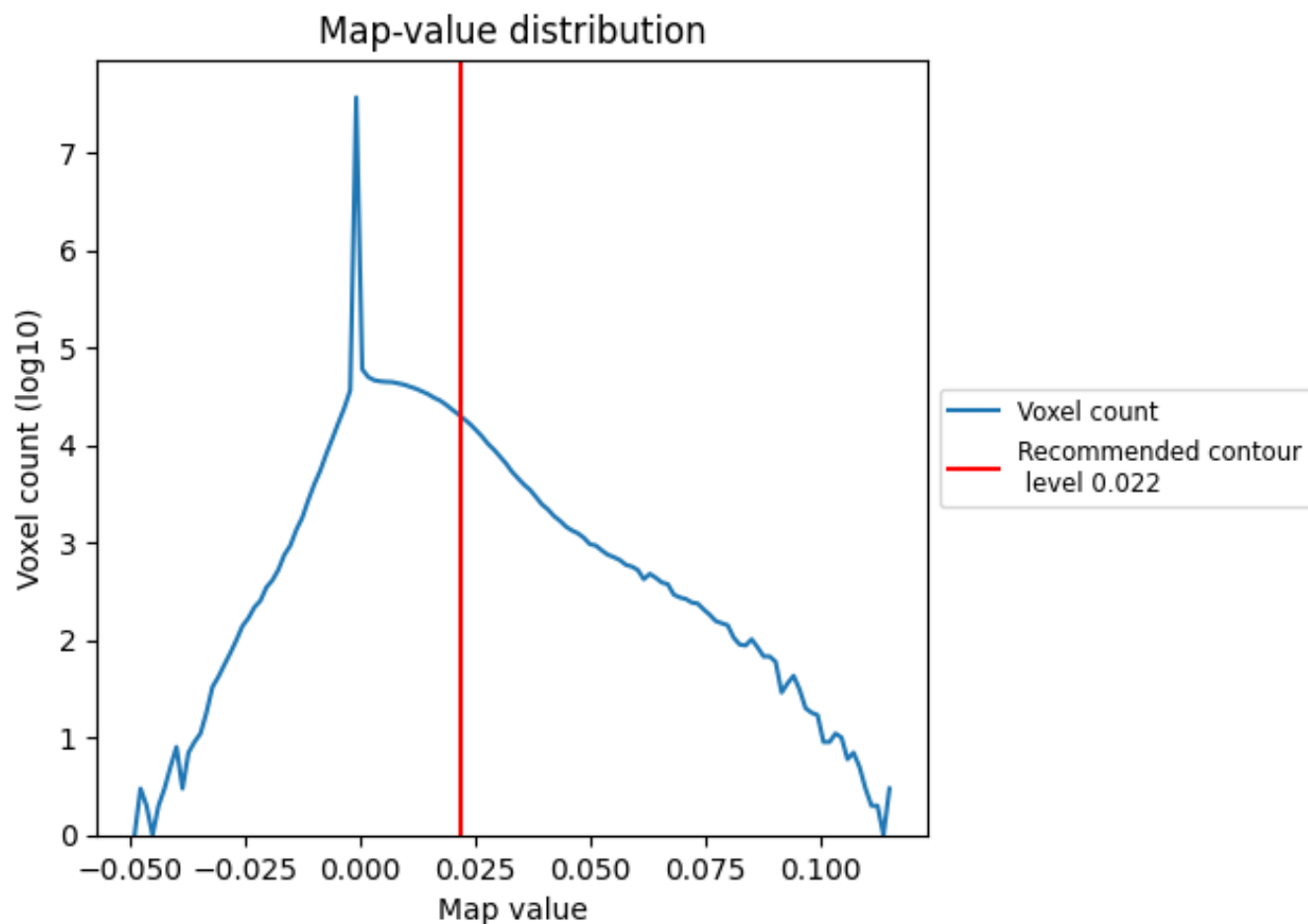


Z

7 Map analysis [i](#)

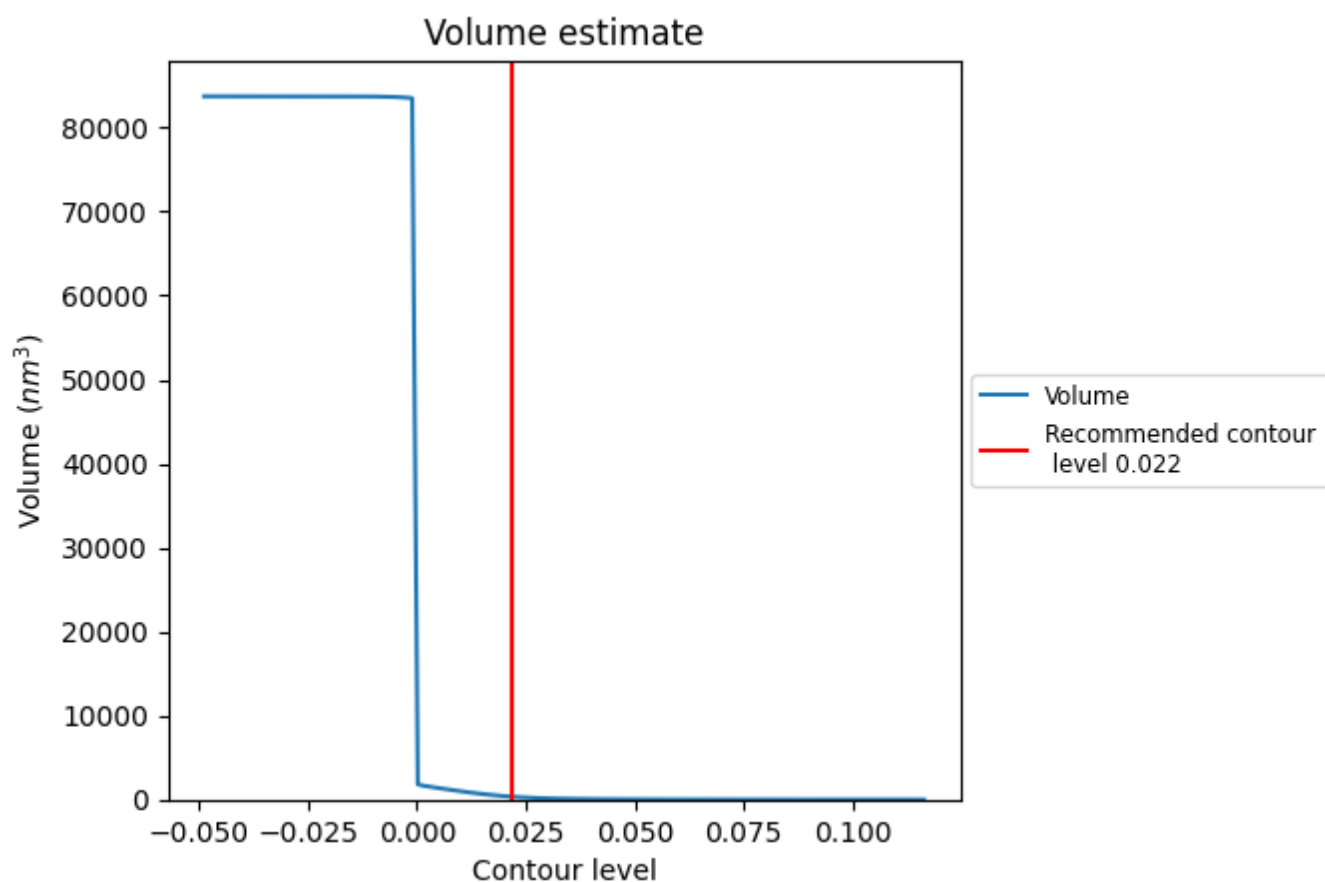
This section contains the results of statistical analysis of the map.

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.

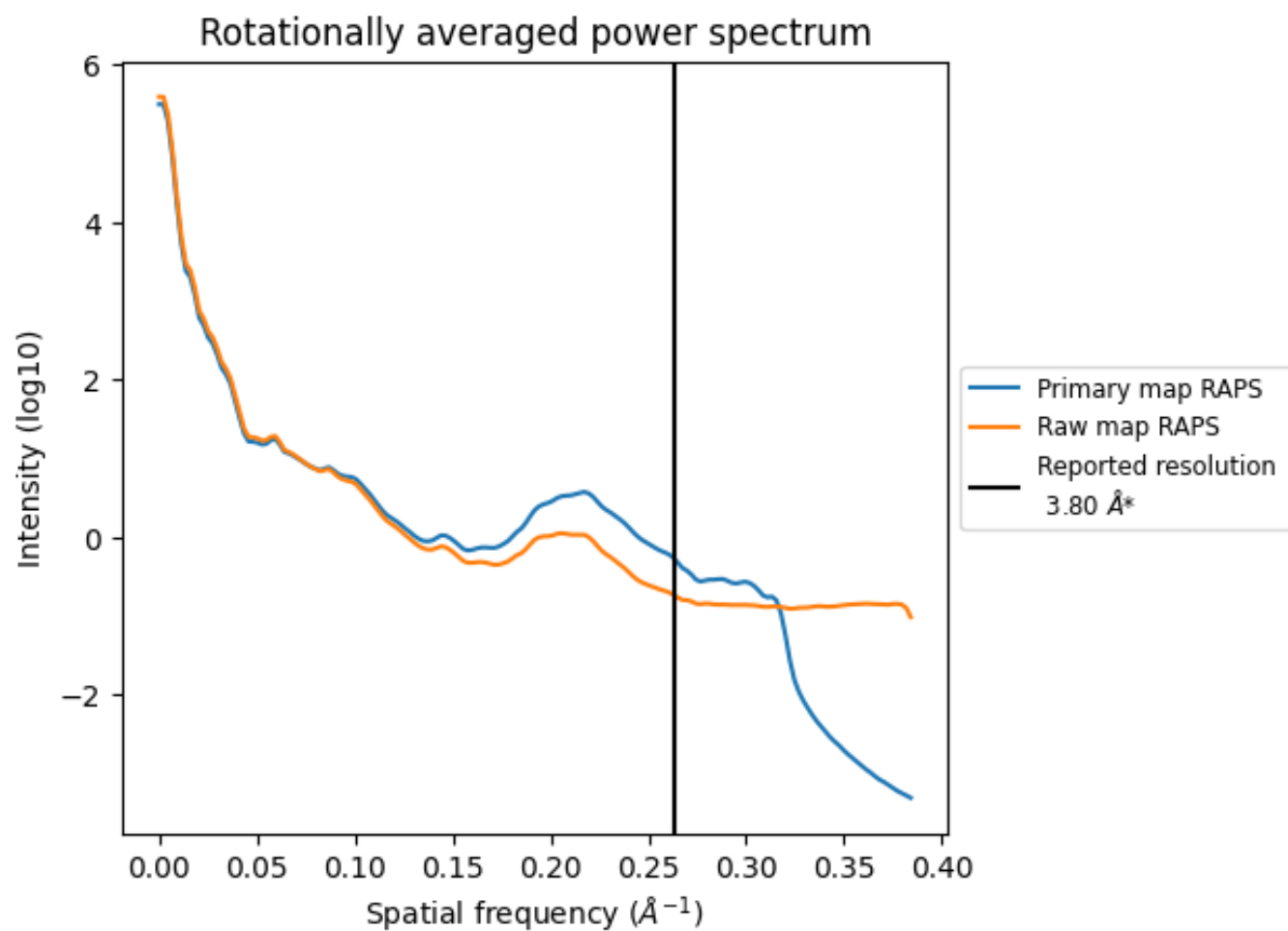
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 334 nm^3 ; this corresponds to an approximate mass of 302 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 ⓘ

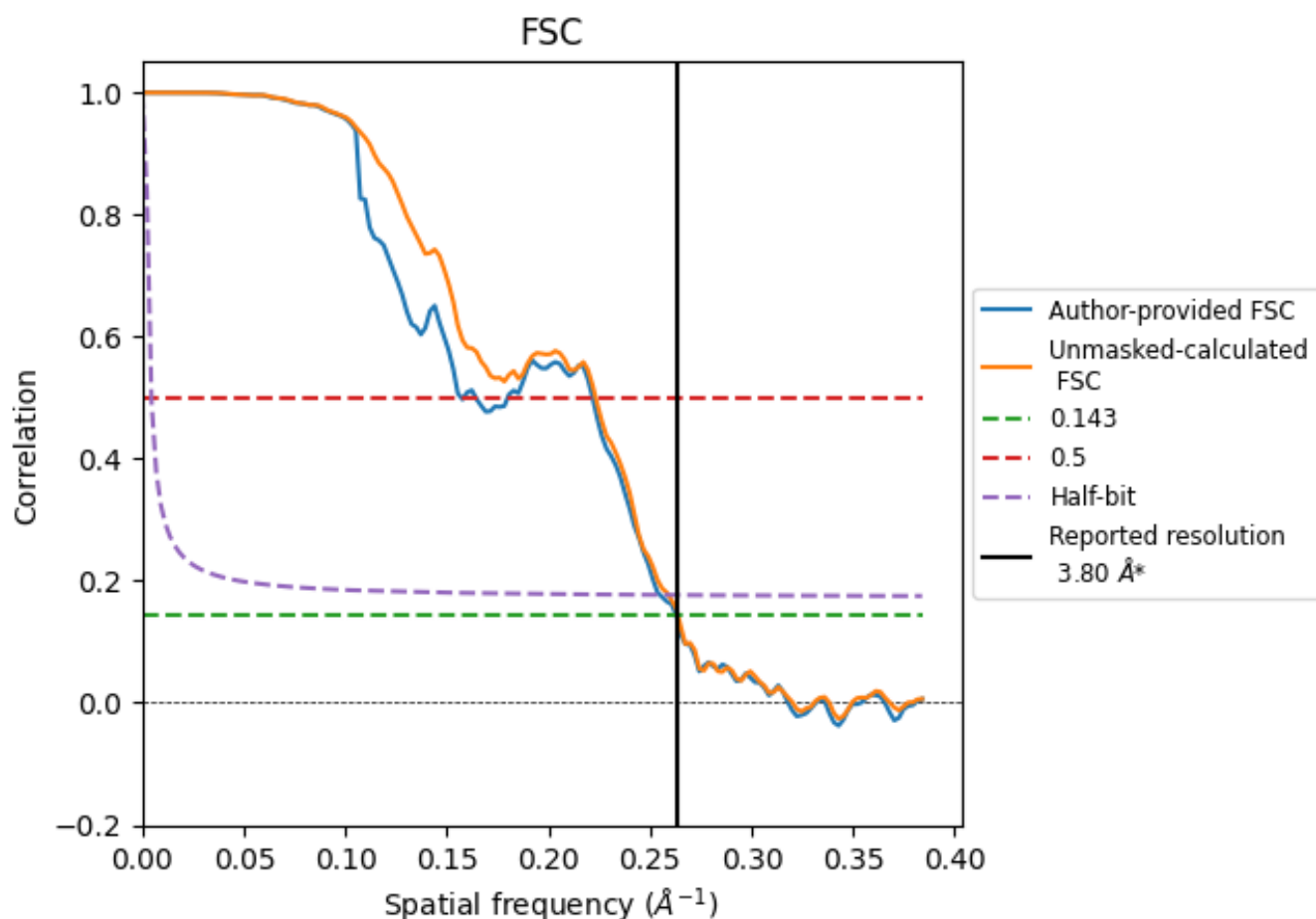


*Reported resolution corresponds to spatial frequency of $0.263 \text{ \AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.263 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

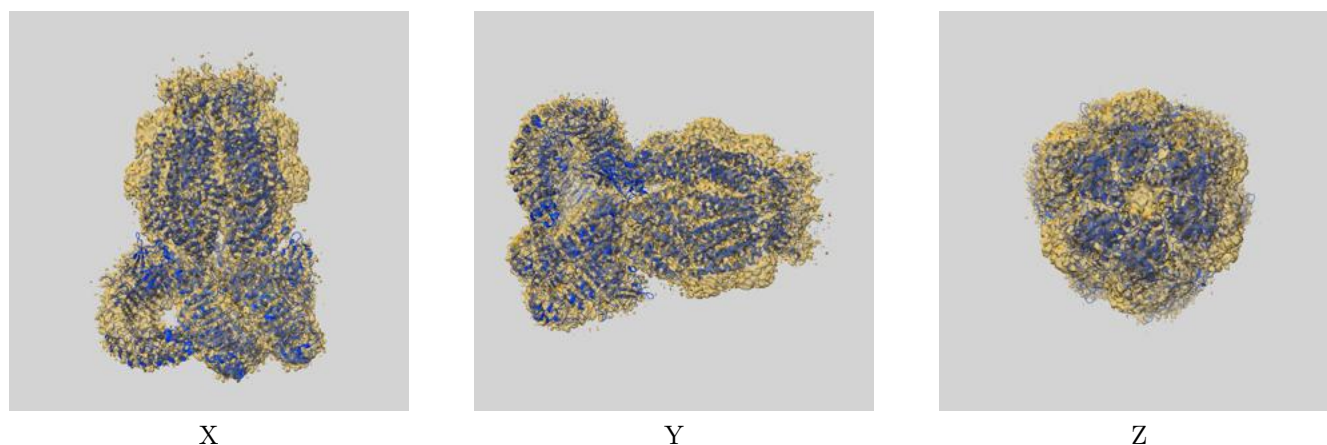
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.80	-	-
Author-provided FSC curve	3.80	6.37	3.92
Unmasked-calculated*	3.79	4.48	3.86

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

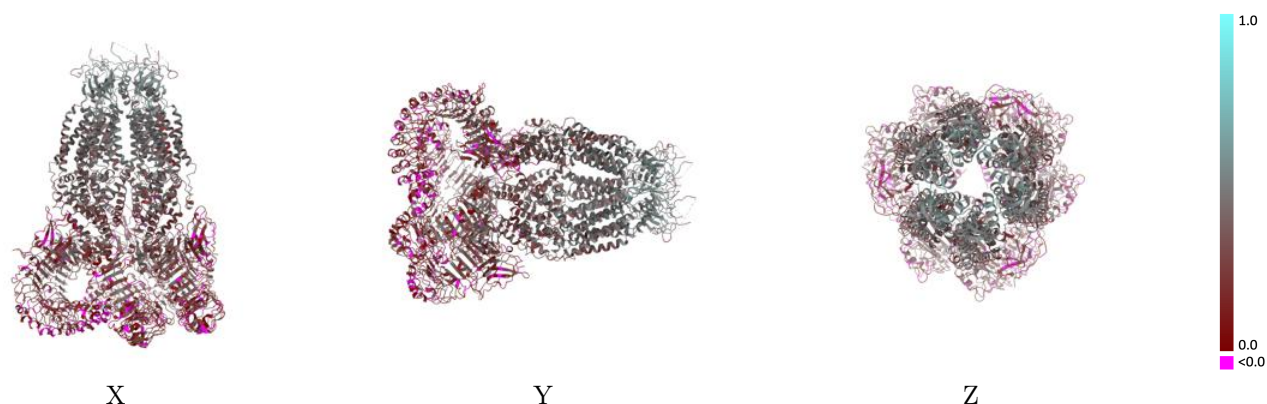
This section contains information regarding the fit between EMDB map EMD-13230 and PDB model 7P6K. Per-residue inclusion information can be found in [section 3](#) on [page 5](#).

9.1 Map-model overlay [i](#)



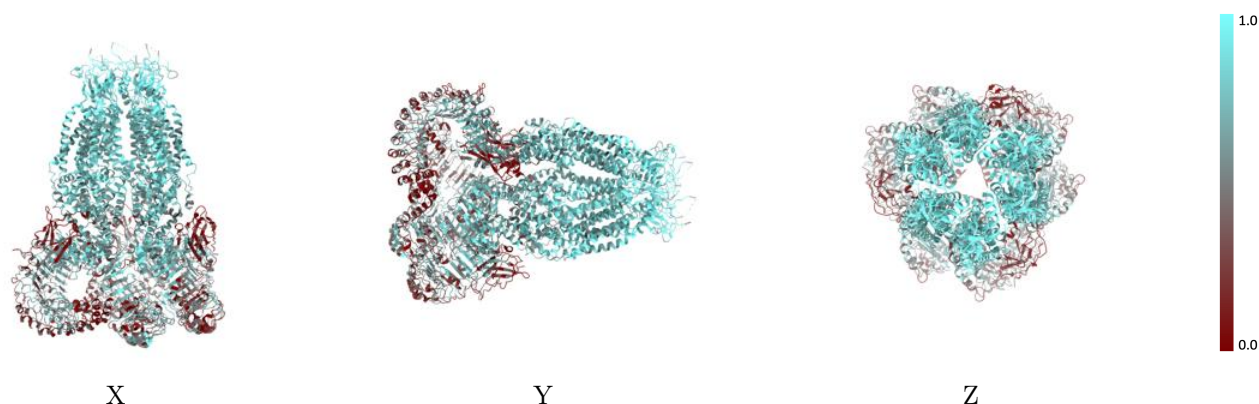
The images above show the 3D surface view of the map at the recommended contour level 0.022 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



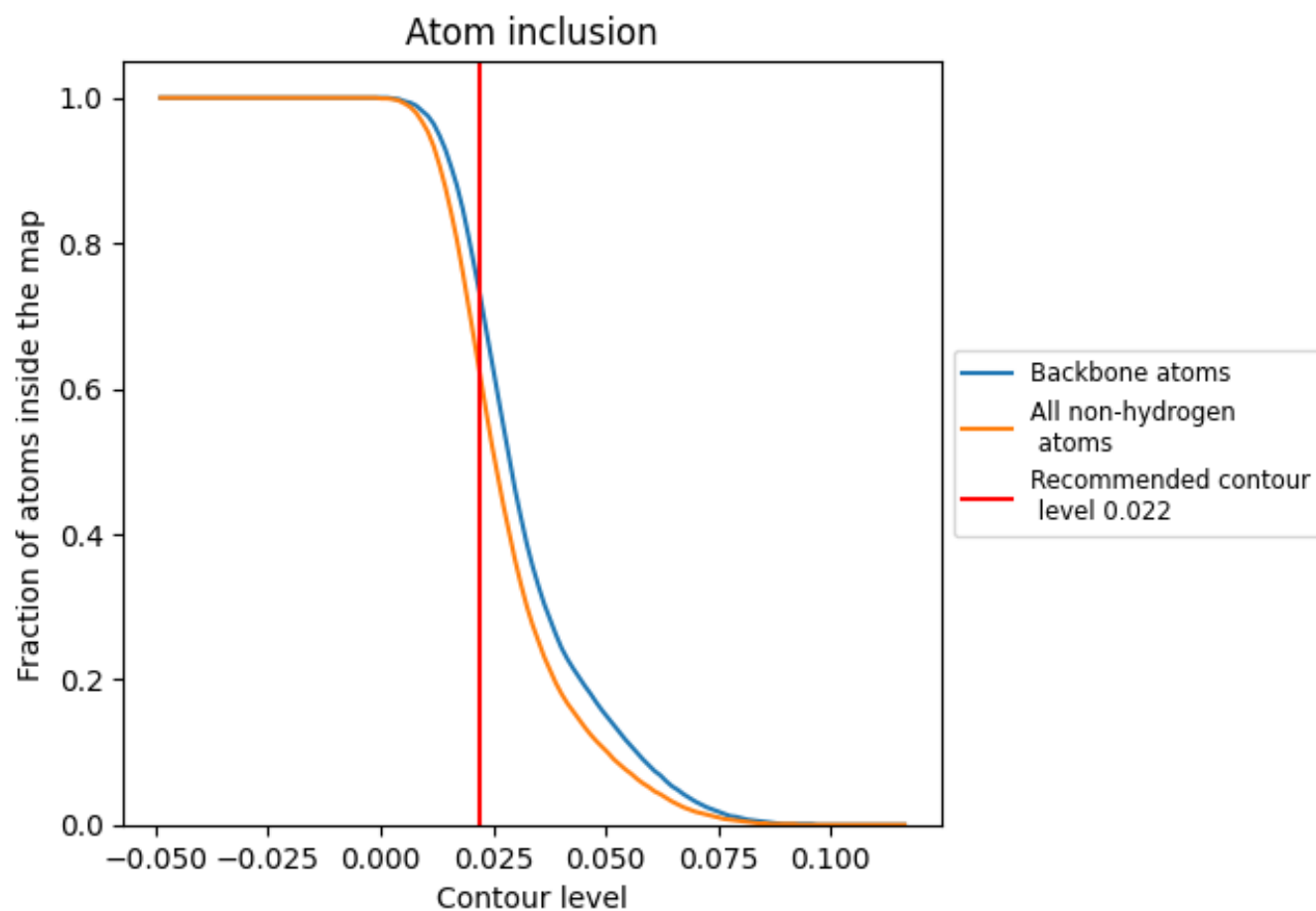
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.022).

9.4 Atom inclusion [i](#)



At the recommended contour level, 73% 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 (0.022) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6220	0.3030
A	0.6660	0.3160
B	0.6460	0.3100
C	0.6620	0.3140
D	0.6390	0.3100
E	0.6660	0.3140
F	0.6440	0.3090
G	0.1870	0.1720
H	0.1830	0.1720
I	0.1910	0.1750

1.0

0.0

<0.0